



wwPDB EM Validation Summary Report ⓘ

Sep 16, 2026 – 11:01 PM EDT

PDB ID : 7UPH / pdb_00007uph
EMDB ID : EMD-26666
Title : Structure of a ribosome with tethered subunits
Authors : Kim, D.S.; Watkins, A.; Bidstrup, E.; Lee, J.; Topkar, V.V.; Kofman, C.;
Schwarz, K.J.; Liu, Y.; Pintilie, G.; Roney, E.; Das, R.; Jewett, M.C.
Deposited on : 2022-04-15
Resolution : 4.18 Å(reported)
Based on initial model : 4YBB

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.dev133
Mogul : 2022.3.0, CSD as543be (2022)
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) : 2.50

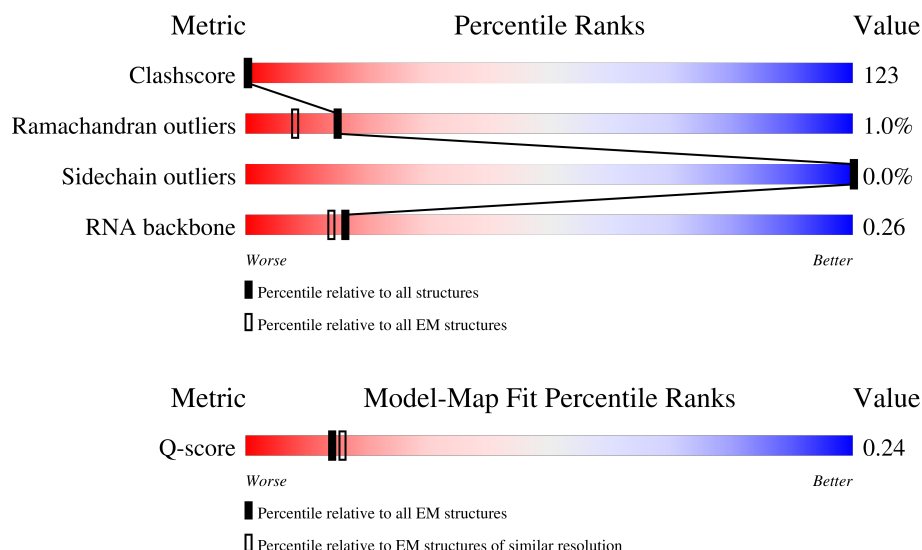
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 4.18 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.



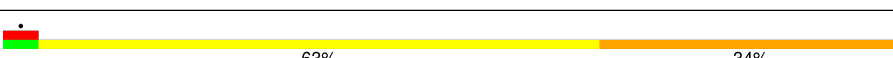
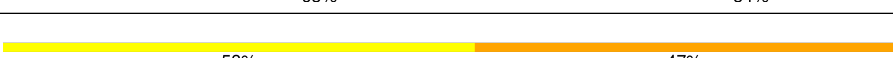
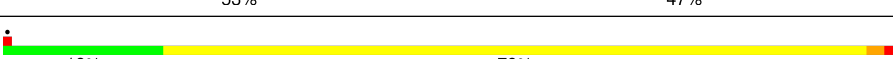
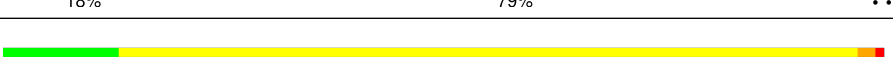
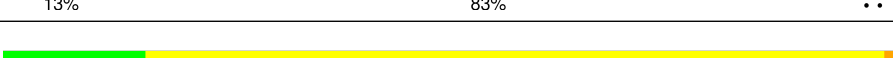
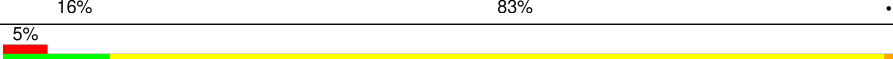
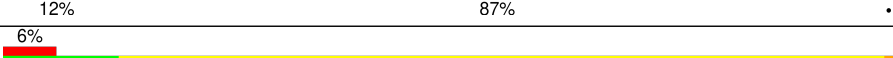
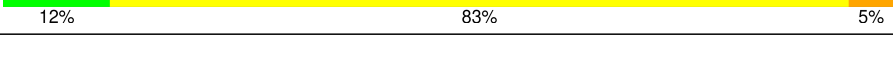
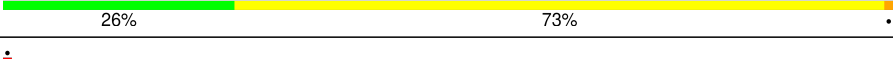
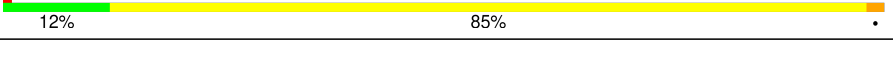
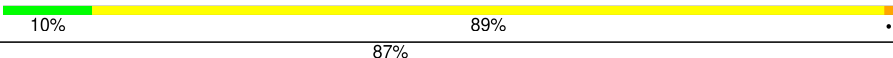
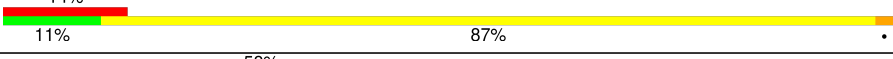
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	5389 (3.68 - 4.68)

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	100	
2	B	88	
3	C	82	

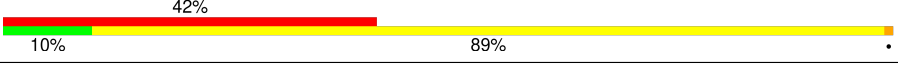
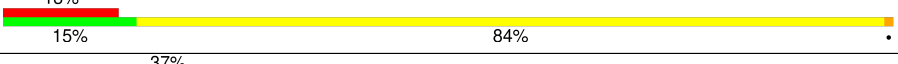
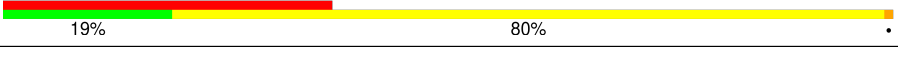
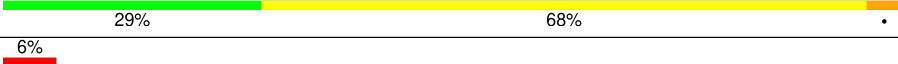
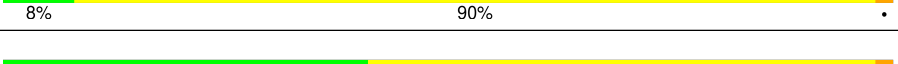
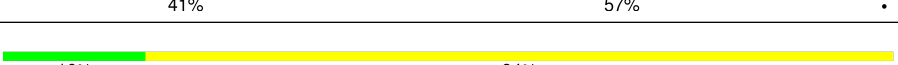
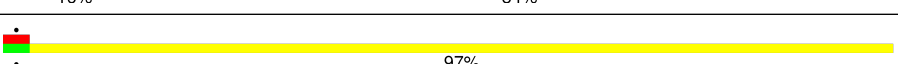
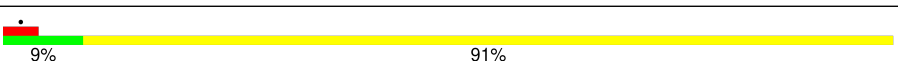
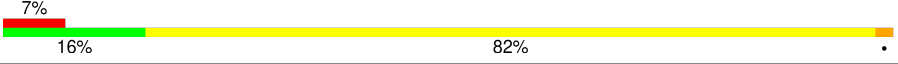
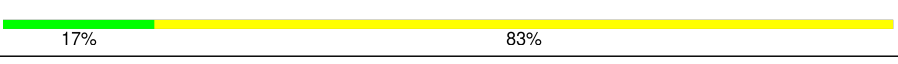
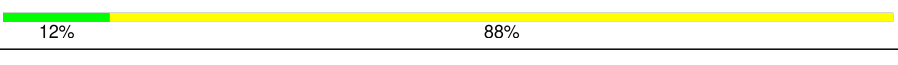
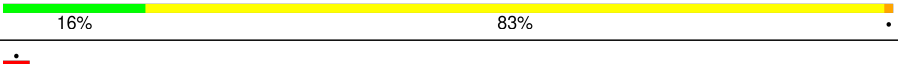
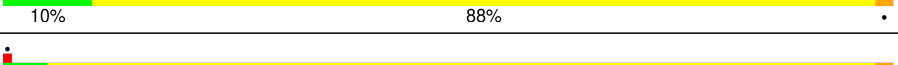
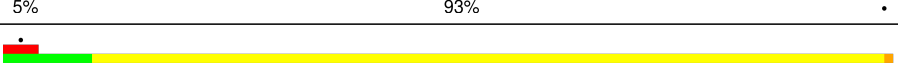
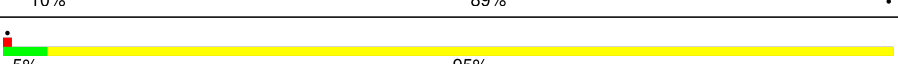
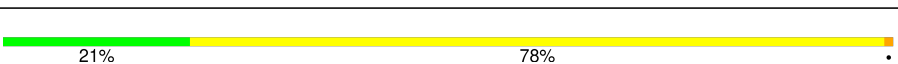
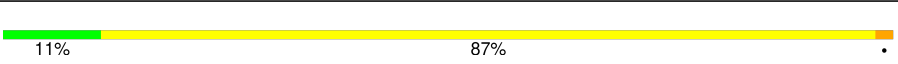
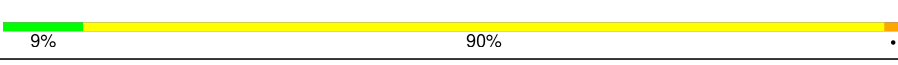
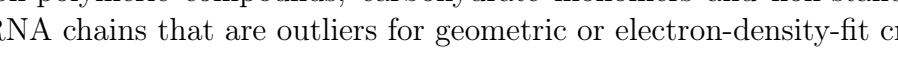
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Mol	Chain	Length	Quality of chain
4	D	80	
5	E	55	
6	F	79	
7	G	85	
8	H	56	
9	I	4438	
10	J	118	
11	K	271	
12	L	209	
13	M	201	
14	N	177	
15	O	176	
16	R	142	
17	S	122	
18	T	144	
19	U	136	
20	V	120	
21	W	224	
22	X	206	
23	Y	205	
24	Z	150	
25	a	100	
26	b	151	
27	c	129	
28	d	127	

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Mol	Chain	Length	Quality of chain
29	e	98	
30	f	117	
31	g	123	
32	h	114	
33	i	56	
34	j	50	
35	k	46	
36	l	64	
37	m	38	
38	n	116	
39	o	114	
40	p	117	
41	q	103	
42	r	110	
43	s	93	
44	t	102	
45	u	94	
46	v	75	
47	w	77	
48	x	62	
49	y	58	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
31	D2T	g	88	-	-	X	-
9	1MG	I	2245	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
9	PSU	I	2449	-	-	X	-
9	3TD	I	3409	-	-	X	-
9	PSU	I	3411	-	-	X	-
9	PSU	I	3951	-	-	X	-
9	OMC	I	3992	-	-	X	-
9	5MC	I	966	-	-	X	-

2 Entry composition

There are 49 unique types of molecules in this entry. The entry contains 141080 atoms, of which 625 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a protein called 30S ribosomal protein S14.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	100	Total	C	N	O	S	0	0
			805	499	164	139	3		

- Molecule 2 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	88	Total	C	N	O	S	0	0
			714	439	144	130	1		

- Molecule 3 is a protein called 30S ribosomal protein S16.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	82	Total	C	N	O	S	0	0
			649	406	128	114	1		

- Molecule 4 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	80	Total	C	N	O	S	0	0
			649	411	121	114	3		

- Molecule 5 is a protein called 30S ribosomal protein S18.

Mol	Chain	Residues	Atoms				AltConf	Trace
5	E	55	Total	C	N	O	0	0
			456	288	86	82		

- Molecule 6 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	79	Total	C	N	O	S	0	0
			638	408	120	108	2		

- Molecule 7 is a protein called 30S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	85	Total	C	N	O	S	0	0
			665	411	137	114	3		

- Molecule 8 is a protein called 30S ribosomal protein S21.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	56	Total	C	N	O	S	0	0
			465	290	96	78	1		

- Molecule 9 is a RNA chain called Tethered rRNA.

Mol	Chain	Residues	Atoms						AltConf	Trace
9	I	4438	Total	C	H	N	O	P	0	0
			95909	42519	625	17511	30816	4438		

- Molecule 10 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	118	Total	C	N	O	P	0	0
			2529	1126	464	821	118		

- Molecule 11 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	271	Total	C	N	O	S	0	0
			2083	1288	423	365	7		

- Molecule 12 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	209	Total	C	N	O	S	0	0
			1565	979	288	294	4		

- Molecule 13 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	201	Total	C	N	O	S	0	0
			1552	974	283	290	5		

- Molecule 14 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	N	177	Total	C	N	O	S	0	0
			1411	899	249	257	6		

- Molecule 15 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	O	176	Total	C	N	O	S	0	0
			1323	832	243	246	2		

- Molecule 16 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	R	142	Total	C	N	O	S	0	0
			1129	714	212	199	4		

- Molecule 17 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	S	122	Total	C	N	O	S	0	0
			938	587	180	165	6		

- Molecule 18 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	T	144	Total	C	N	O	S	0	0
			1053	654	207	190	2		

- Molecule 19 is a protein called 50S ribosomal protein L16.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	U	136	Total	C	N	O	S	0	0
			1075	686	205	178	6		

- Molecule 20 is a protein called 50S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	V	120	Total	C	N	O	S	0	0
			960	593	196	166	5		

- Molecule 21 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	W	224	Total	C	N	O	S	0	0
			1753	1109	315	321	8		

- Molecule 22 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	X	206	Total	C	N	O	S	0	0
			1625	1028	305	289	3		

- Molecule 23 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	Y	205	Total	C	N	O	S	0	0
			1643	1026	315	298	4		

- Molecule 24 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	Z	150	Total	C	N	O	S	0	0
			1105	687	211	201	6		

- Molecule 25 is a protein called 30S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	a	100	Total	C	N	O	S	0	0
			817	515	148	148	6		

- Molecule 26 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	b	151	Total	C	N	O	S	0	0
			1182	735	227	216	4		

- Molecule 27 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	c	129	Total	C	N	O	S	0	0
			979	616	173	184	6		

- Molecule 28 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	d	127	Total	C	N	O	S	0	0
			1022	634	206	179	3		

- Molecule 29 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	e	98	Total	C	N	O	S	0	0
			787	493	150	143	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
30	f	117	Total	C	N	O	S	0	0
			877	540	174	160	3		

- Molecule 31 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	g	123	Total	C	N	O	S	0	0
			957	591	196	165	5		

- Molecule 32 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	h	114	Total	C	N	O	S	0	0
			884	546	178	157	3		

- Molecule 33 is a protein called 50S ribosomal protein L32.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	i	56	Total	C	N	O	S	0	0
			444	269	94	80	1		

- Molecule 34 is a protein called 50S ribosomal protein L33.

Mol	Chain	Residues	Atoms				AltConf	Trace
34	j	50	Total	C	N	O	0	0
			409	263	75	71		

- Molecule 35 is a protein called 50S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	k	46	Total	C	N	O	S	0	0
			377	228	90	57	2		

- Molecule 36 is a protein called 50S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	l	64	Total	C	N	O	S	0	0
			504	323	105	74	2		

- Molecule 37 is a protein called 50S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	m	38	Total	C	N	O	S	0	0
			302	185	65	48	4		

- Molecule 38 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	n	116	Total	C	N	O		0	0
			892	552	178	162			

- Molecule 39 is a protein called 50S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	o	114	Total	C	N	O	S	0	0
			917	574	179	163	1		

- Molecule 40 is a protein called 50S ribosomal protein L20.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	p	117	Total	C	N	O		0	0
			947	604	192	151			

- Molecule 41 is a protein called 50S ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	q	103	Total	C	N	O	S	0	0
			816	516	153	145	2		

- Molecule 42 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	r	110	Total	C	N	O	S	0	0
			857	532	166	156	3		

- Molecule 43 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	s	93	Total	C	N	O	S	0	0
			739	466	139	132	2		

- Molecule 44 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	t	102	Total	C	N	O		0	0
			780	492	146	142			

- Molecule 45 is a protein called 50S ribosomal protein L25.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	u	94	Total	C	N	O	S	0	0
			753	479	137	134	3		

- Molecule 46 is a protein called 50S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	v	75	Total	C	N	O	S	0	0
			569	353	113	102	1		

- Molecule 47 is a protein called 50S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	w	77	Total	C	N	O	S	0	0
			625	388	129	106	2		

- Molecule 48 is a protein called 50S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	x	62	Total	C	N	O	S	0	0
			501	308	98	94	1		

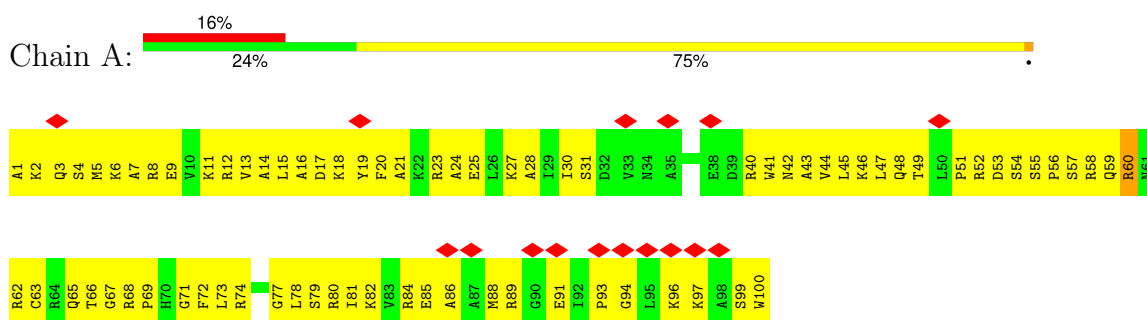
- Molecule 49 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	y	58	Total	C	N	O	S	0	0
			449	281	87	79	2		

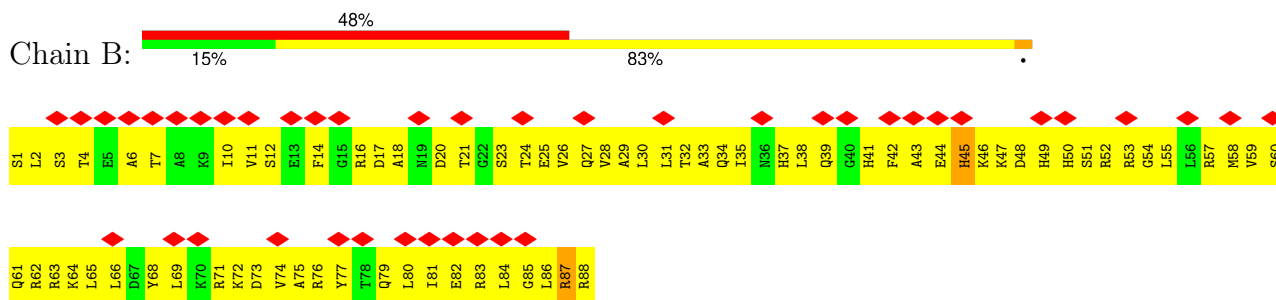
3 Residue-property plots

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.

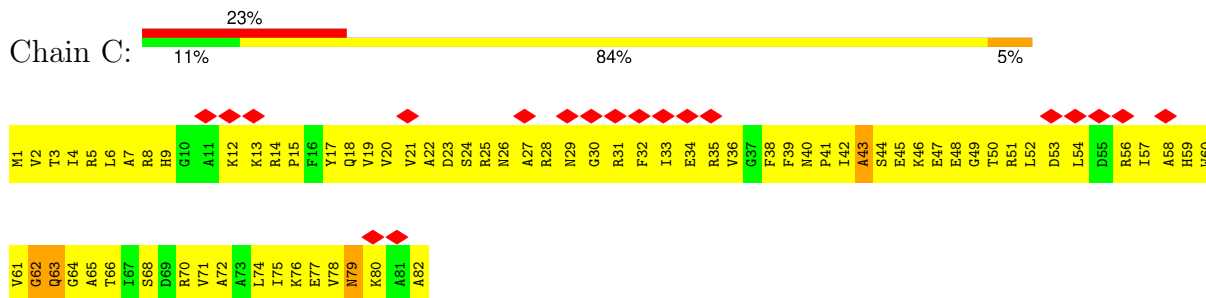
• Molecule 1: 30S ribosomal protein S14



• Molecule 2: 30S ribosomal protein S15

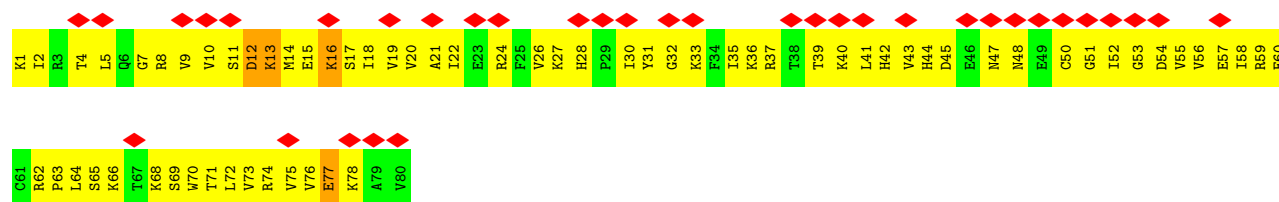


• Molecule 3: 30S ribosomal protein S16

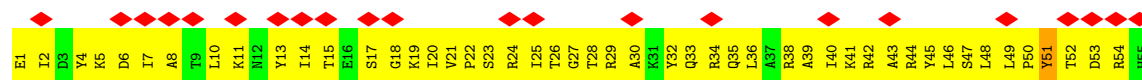


• Molecule 4: 30S ribosomal protein S17

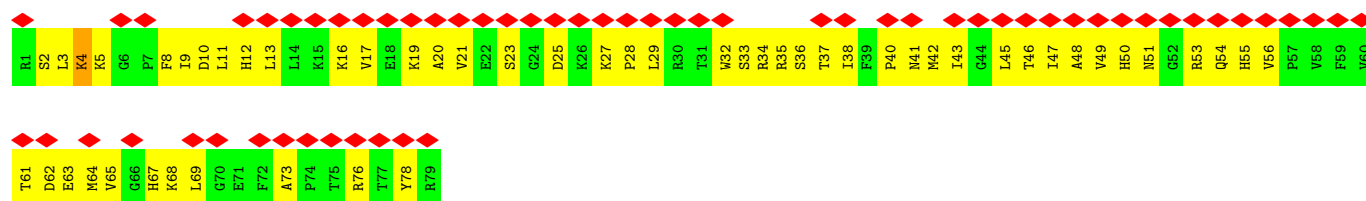
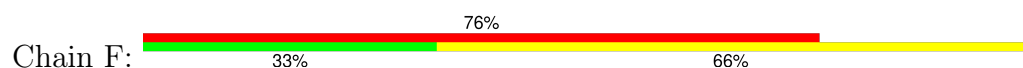




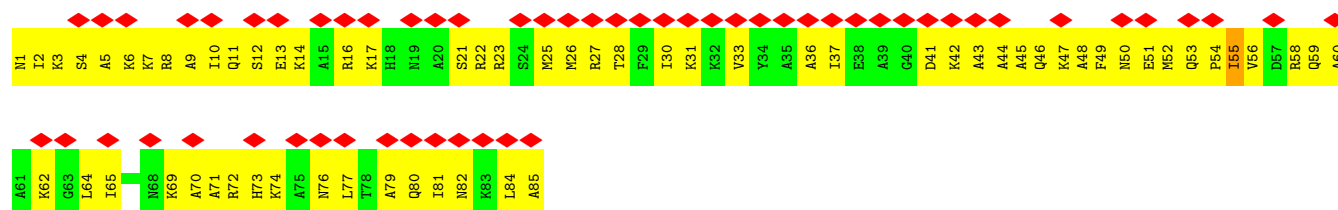
• Molecule 5: 30S ribosomal protein S18



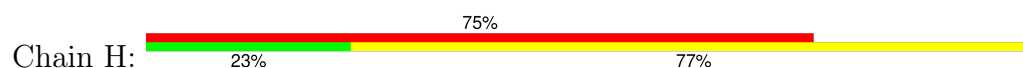
• Molecule 6: 30S ribosomal protein S19



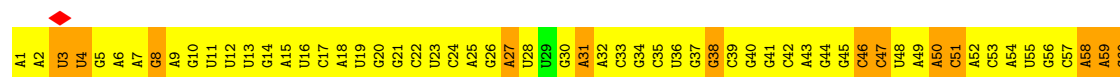
• Molecule 7: 30S ribosomal protein S20



• Molecule 8: 30S ribosomal protein S21



• Molecule 9: Tethered rRNA





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G3702	C3639	A3582	U3522	G3461	G3400	G3340	A3280	C3219	A3159	A3097	G3037	G2978	G2917	G2857	A2795
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G4401	G4402	G4403	A4404	A4405	A4406	A4407	G4408	A4409	U4410	A4411	A4412	C4413	C4414	C4415	U4416	U4417	C4418	C4419	C4420	C4421	A4422	A4423	C4424	C4425	U4426	C4427	C4428	G4429	C4430	C4431	C4432	C4433	C4434	A4435	U4436	C4437	A4438																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																											
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• Molecule 10: 5S rRNA

Chain J:

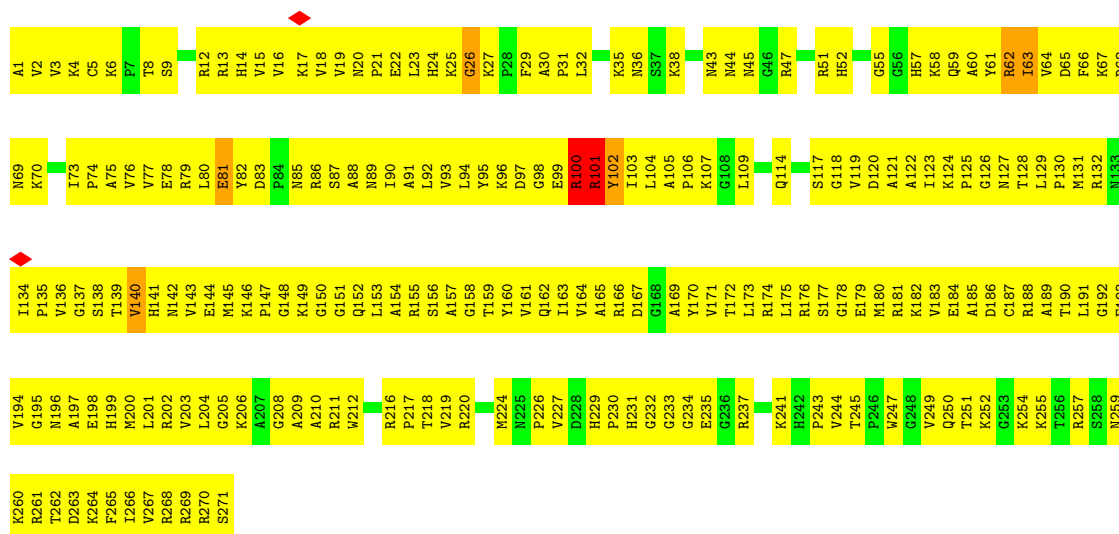


G1	G2	G3	G4	G5	G6	G7	G8	G9	G10	G11	G12	G13	G14	G15	G16	G17	G18	G19	G20	G21	G22	G23	G24	G25	G26	G27	G28	G29	G30	G31	G32	G33	G34	G35	G36	G37	G38	G39	G40	G41	G42	G43	G44	G45	G46	G47	G48	G49	G50	G51	G52	G53	G54	G55	G56	G57	G58	G59	G60
C61	C62	C63	C64	C65	C66	C67	C68	C69	C70	C71	C72	C73	C74	C75	C76	C77	C78	C79	C80	C81	C82	C83	C84	C85	C86	C87	C88	C89	C90	C91	C92	C93	C94	C95	C96	C97	C98	C99	C100	C101	C102	C103	C104	C105	C106	C107	C108	C109	C110	C111	C112	C113	C114	C115	C116	C117	C118		

• Molecule 11: 50S ribosomal protein L2

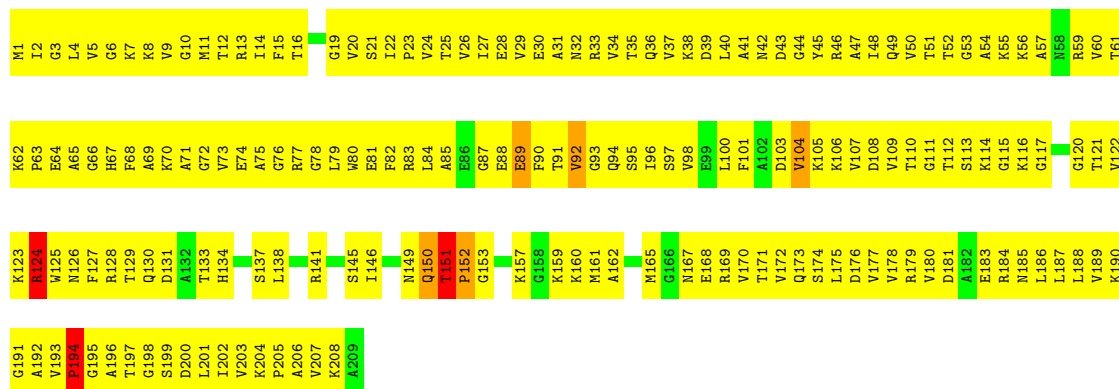
Chain K:





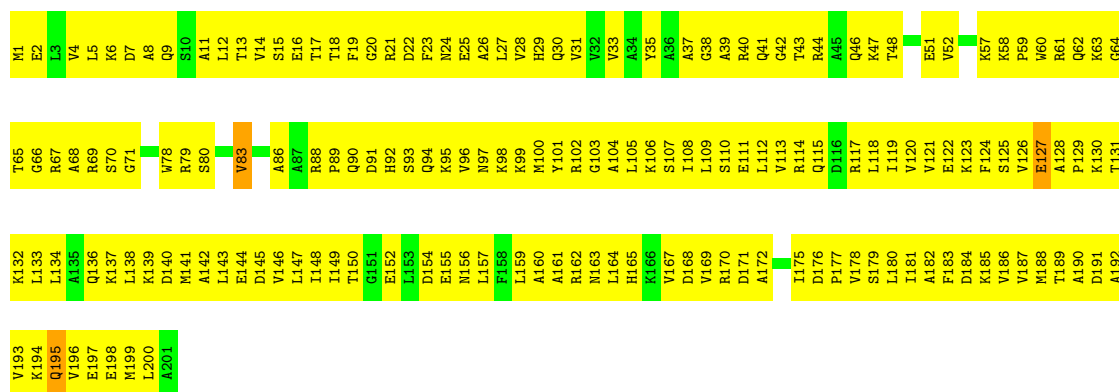
• Molecule 12: 50S ribosomal protein L3

Chain L: 13% 83% ..

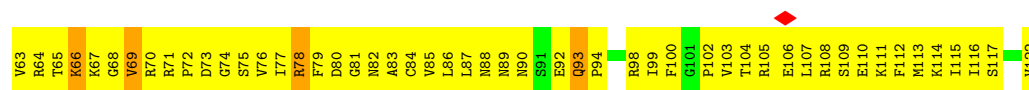


• Molecule 13: 50S ribosomal protein L4

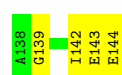
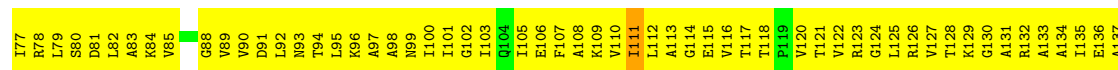
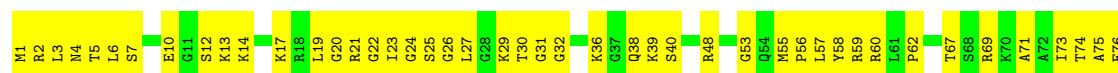
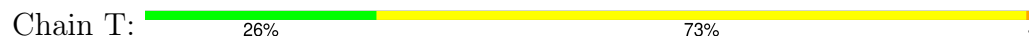
Chain M: 16% 83% .



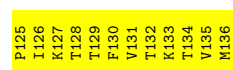
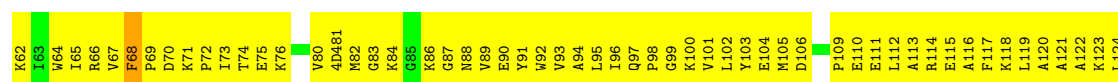
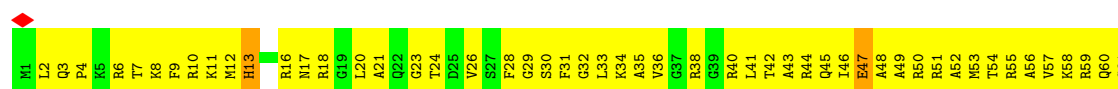
• Molecule 14: 50S ribosomal protein L5



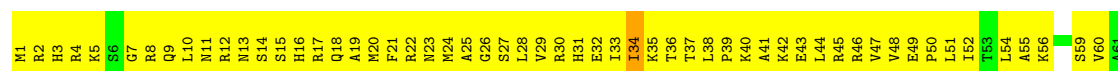
• Molecule 18: 50S ribosomal protein L15



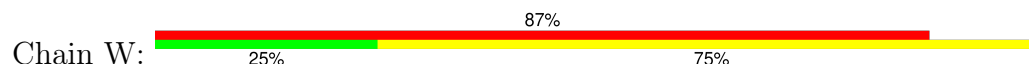
• Molecule 19: 50S ribosomal protein L16

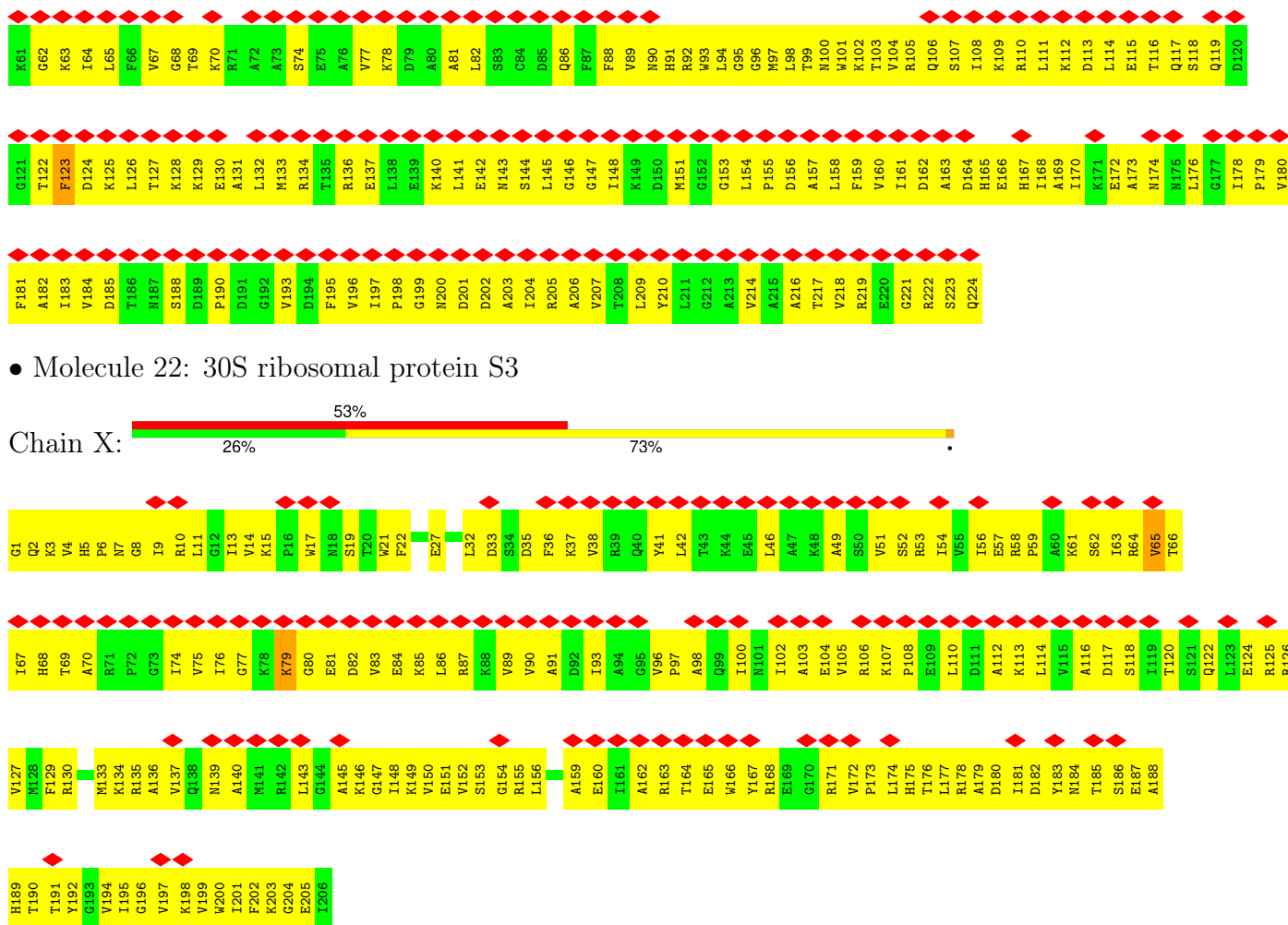


• Molecule 20: 50S ribosomal protein L17

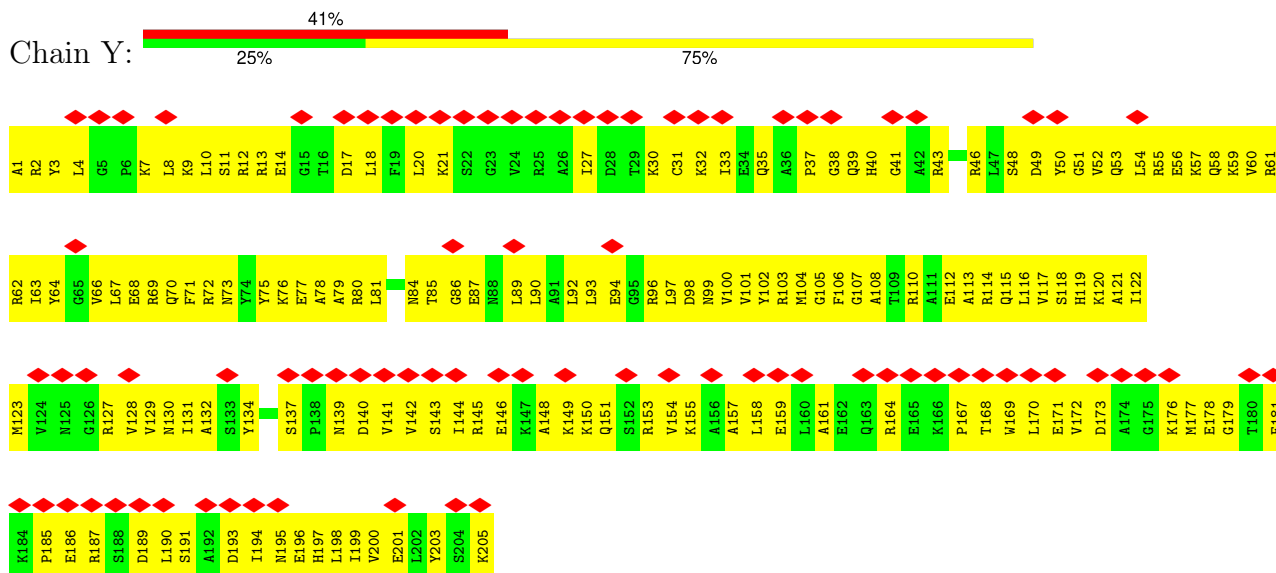


• Molecule 21: 30S ribosomal protein S2

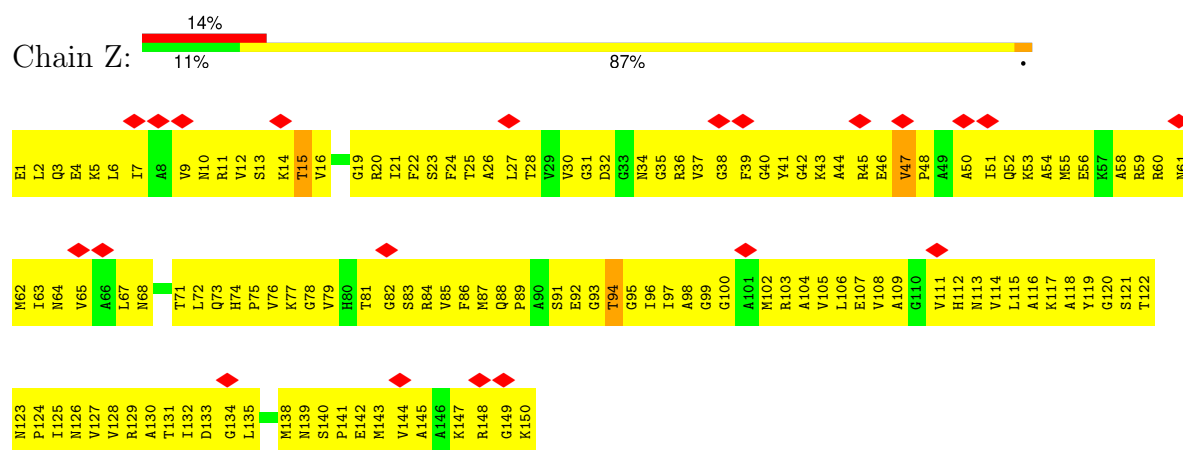




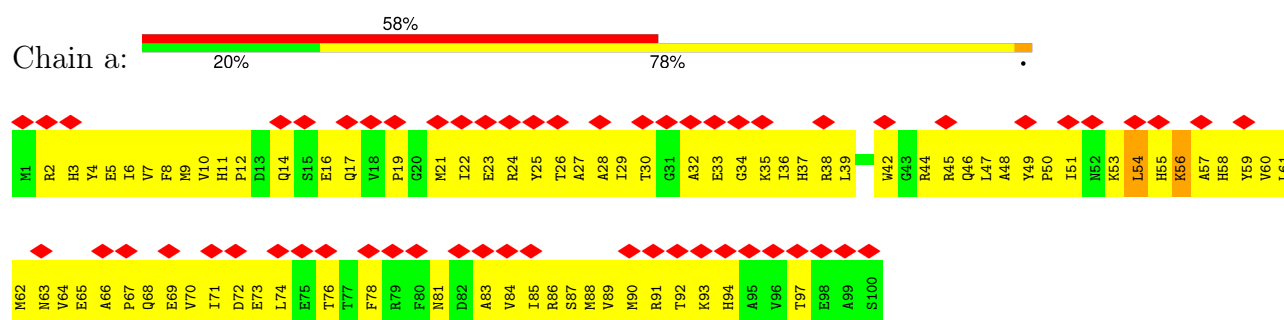
• Molecule 23: 30S ribosomal protein S4



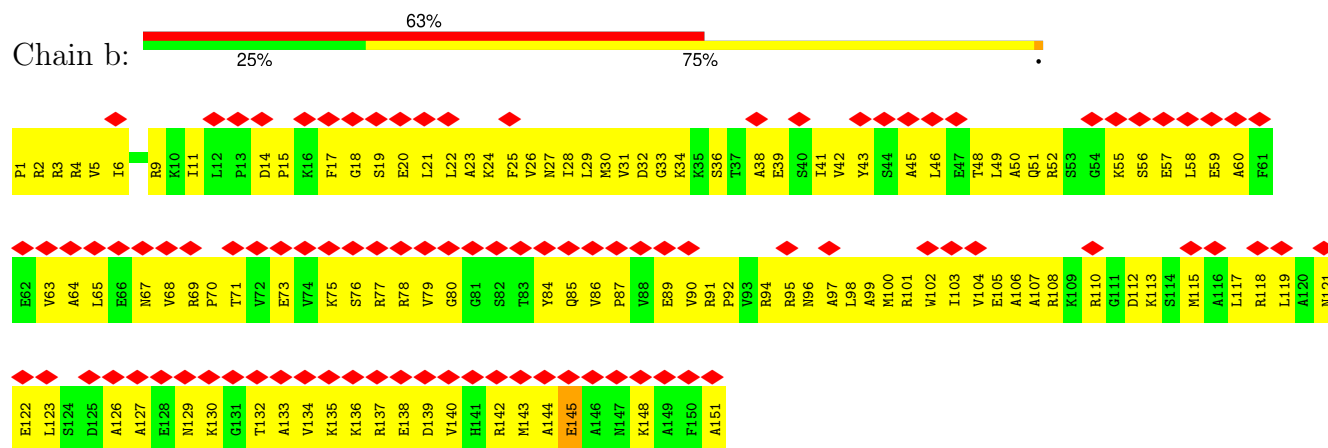
• Molecule 24: 30S ribosomal protein S5



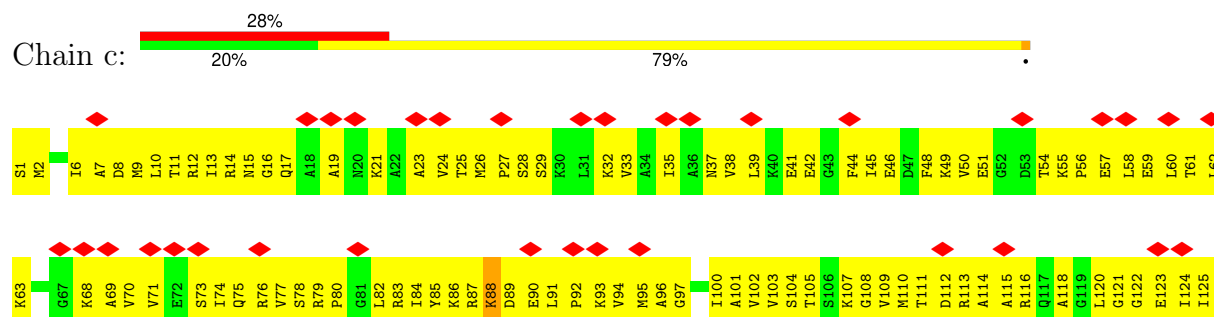
• Molecule 25: 30S ribosomal protein S6



• Molecule 26: 30S ribosomal protein S7



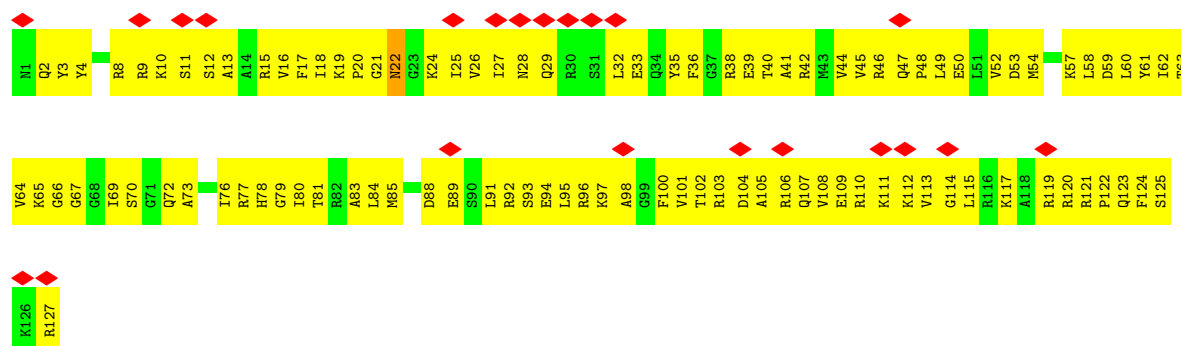
• Molecule 27: 30S ribosomal protein S8



C126
Y127
V128
A129

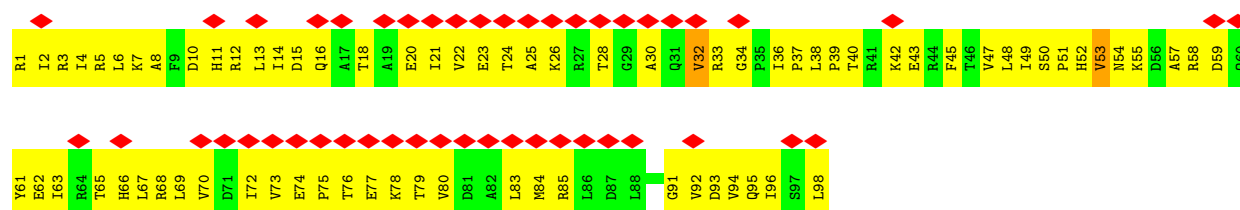
• Molecule 28: 30S ribosomal protein S9

Chain d: 

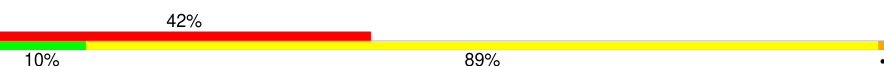


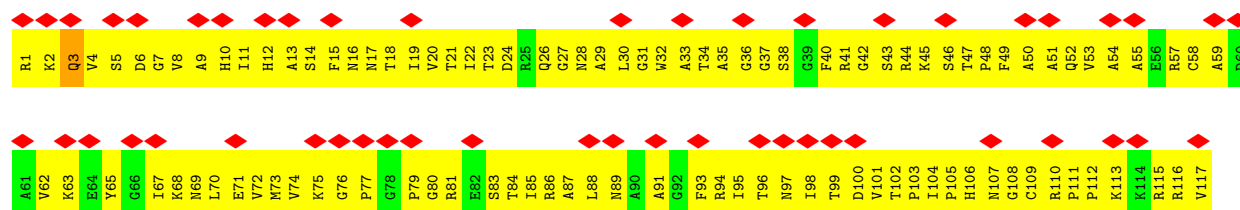
• Molecule 29: 30S ribosomal protein S10

Chain e: 



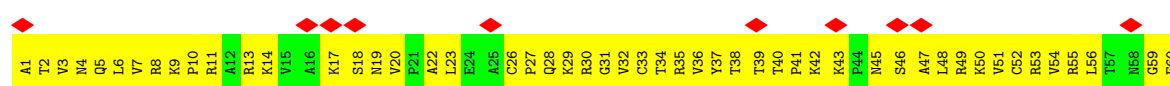
• Molecule 30: 30S ribosomal protein S11

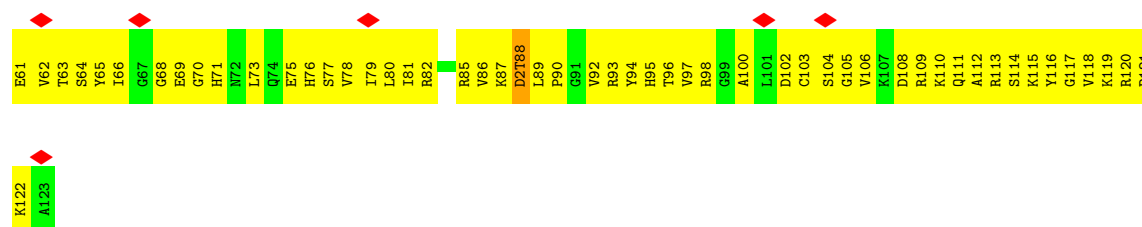
Chain f: 



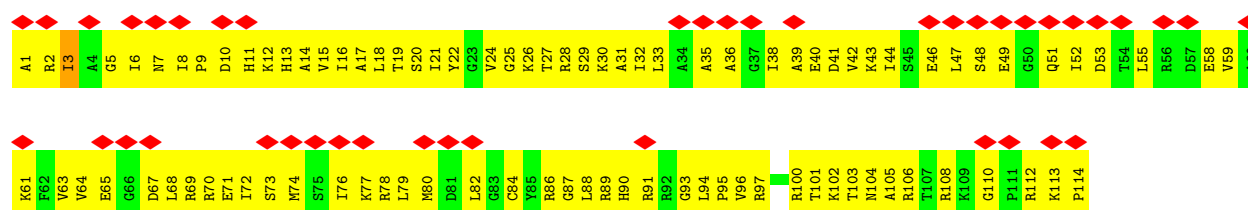
• Molecule 31: 30S ribosomal protein S12

Chain g: 





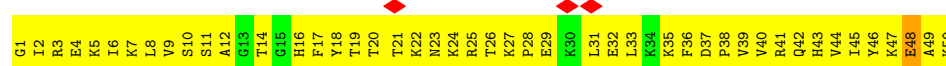
• Molecule 32: 30S ribosomal protein S13



• Molecule 33: 50S ribosomal protein L32



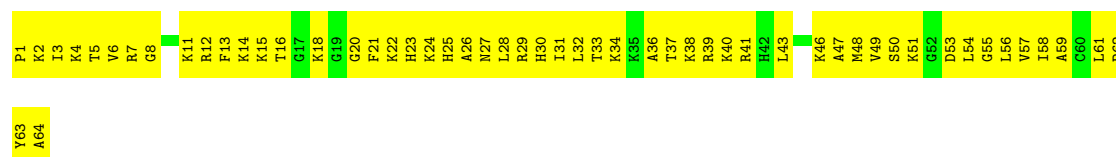
• Molecule 34: 50S ribosomal protein L33



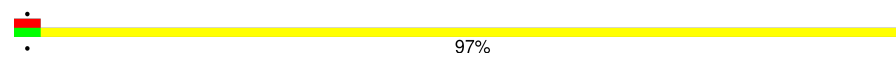
• Molecule 35: 50S ribosomal protein L34



• Molecule 36: 50S ribosomal protein L35



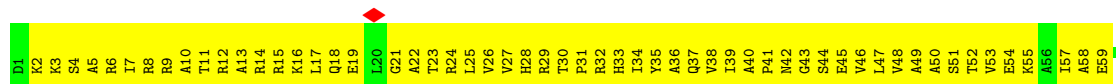
• Molecule 37: 50S ribosomal protein L36

Chain m:  97%



• Molecule 38: 50S ribosomal protein L18

Chain n:  91%



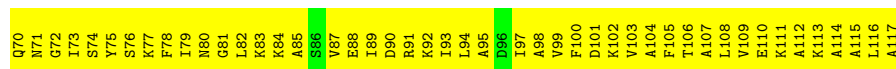
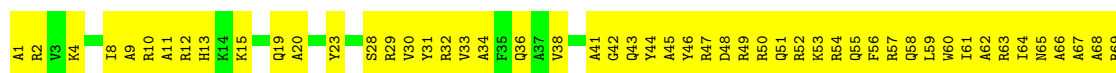
• Molecule 39: 50S ribosomal protein L19

Chain o:  82%

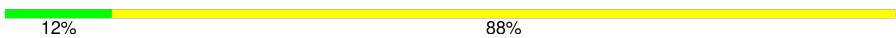


• Molecule 40: 50S ribosomal protein L20

Chain p:  83%

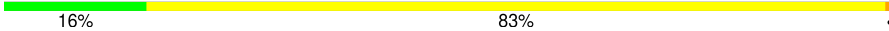


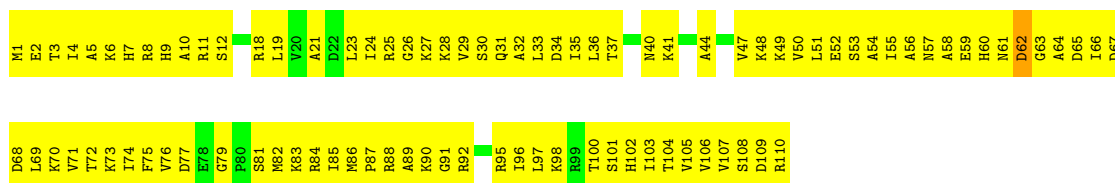
• Molecule 41: 50S ribosomal protein L21

Chain q:  88%

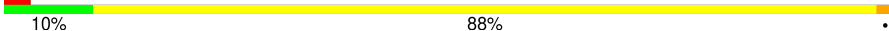


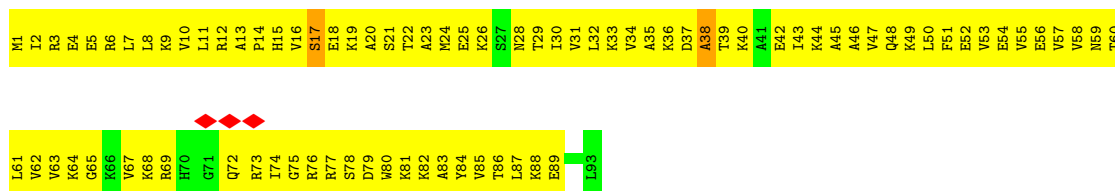
• Molecule 42: 50S ribosomal protein L22

Chain r:  16% 83%

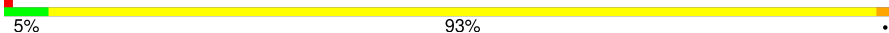


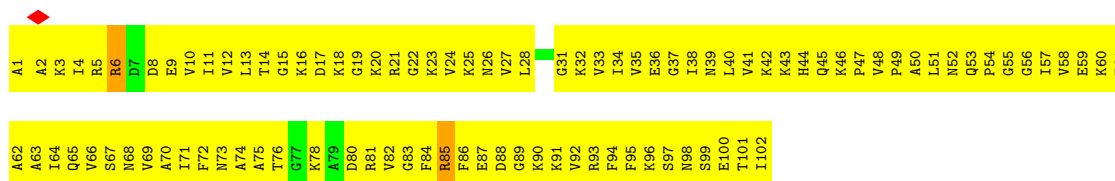
• Molecule 43: 50S ribosomal protein L23

Chain s:  10% 88%

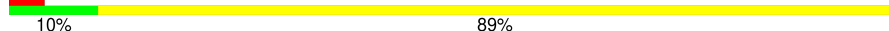


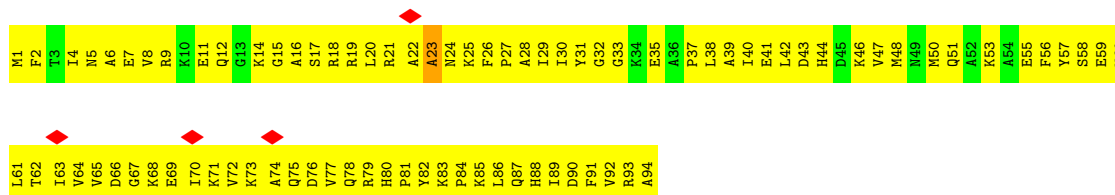
• Molecule 44: 50S ribosomal protein L24

Chain t:  5% 93%

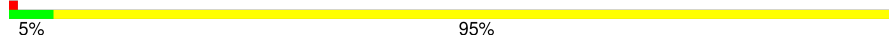


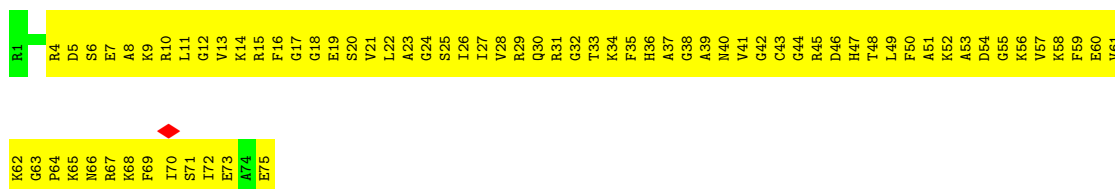
• Molecule 45: 50S ribosomal protein L25

Chain u:  10% 89%



• Molecule 46: 50S ribosomal protein L27

Chain v:  5% 95%



A1	K2	T3	I4	K5	I6	T7	Q8	T9	R10	R11	A12	I13	G14	R15	L16	P17	K18	H19	K20	A21	T22	L23		L26	G27	L28	R29	R30	I31	G32	R33	T34	V35	E36	R37	E38	D39	L40	P41	I42	I43	R44	G45	M46	N47	N48	A49	V50	S51	F52	F53	V54	K55	V56	E57	T58
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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	34852	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	TFS TALOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1200	Depositor
Maximum defocus (nm)	2200	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.073	Depositor
Minimum map value	-0.210	Depositor
Average map value	0.005	Depositor
Map value standard deviation	0.058	Depositor
Recommended contour level	0.1	Depositor
Map size ($\text{\AA}$)	430.0, 430.0, 430.0	wwPDB
Map dimensions	500, 500, 500	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.86, 0.86, 0.86	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: OMU, UR3, 5MU, 5MC, OMC, 3TD, 6MZ, MA6, 2MG, 4OC, D2T, 4D4, PSU, OMG, G7M, 2MA, 1MG

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	0.24	0/817	0.46	0/1088
2	B	0.20	0/722	0.51	0/964
3	C	0.81	2/659 (0.3%)	1.06	3/884 (0.3%)
4	D	0.28	0/658	0.64	0/881
5	E	0.18	0/463	0.56	0/621
6	F	0.16	0/653	0.43	0/877
7	G	0.20	0/671	0.52	1/888 (0.1%)
8	H	0.16	0/472	0.72	2/627 (0.3%)
9	I	0.33	6/105898 (0.0%)	0.46	4/165194 (0.0%)
10	J	0.33	0/2828	0.57	0/4410
11	K	0.80	6/2122 (0.3%)	0.99	7/2852 (0.2%)
12	L	0.49	1/1586 (0.1%)	0.84	2/2134 (0.1%)
13	M	0.35	0/1571	0.69	2/2113 (0.1%)
14	N	0.40	0/1435	0.68	1/1926 (0.1%)
15	O	0.34	0/1343	0.70	0/1816
16	R	0.54	1/1152 (0.1%)	0.83	0/1551
17	S	0.60	1/947 (0.1%)	0.93	2/1268 (0.2%)
18	T	0.42	0/1062	0.83	4/1413 (0.3%)
19	U	0.43	0/1081	0.73	2/1443 (0.1%)
20	V	0.39	0/973	0.78	0/1301
21	W	0.14	0/1784	0.40	0/2403
22	X	0.13	0/1652	0.36	0/2225
23	Y	0.13	0/1665	0.32	0/2227
24	Z	0.38	1/1118 (0.1%)	0.63	0/1504
25	a	0.21	0/835	0.50	0/1128
26	b	0.15	0/1196	0.41	0/1602
27	c	0.16	0/989	0.41	0/1326
28	d	0.16	0/1034	0.42	0/1375
29	e	0.19	0/797	0.50	0/1077
30	f	0.19	0/893	0.55	0/1205
31	g	0.18	0/960	0.41	0/1286

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
32	h	0.17	0/893	0.43	0/1193
33	i	0.33	0/450	0.57	0/599
34	j	0.26	0/416	0.59	0/554
35	k	0.36	0/380	0.66	0/498
36	l	0.36	0/513	0.66	0/676
37	m	0.32	0/303	0.67	0/397
38	n	0.30	0/902	0.68	0/1209
39	o	0.42	0/929	0.75	0/1242
40	p	0.40	0/960	0.71	0/1278
41	q	0.36	0/829	0.67	0/1107
42	r	0.35	0/864	0.57	0/1156
43	s	0.46	0/745	0.75	1/994 (0.1%)
44	t	0.38	0/788	0.84	0/1051
45	u	0.33	0/766	0.71	0/1025
46	v	0.38	0/576	0.74	0/762
47	w	0.37	0/635	0.70	2/848 (0.2%)
48	x	0.50	0/502	0.82	0/667
49	y	0.34	0/453	0.76	0/605
All	All	0.35	18/151940 (0.0%)	0.52	33/227470 (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
1	A	0	1
4	D	0	2
11	K	0	5
12	L	0	3
16	R	0	1
20	V	0	1
39	o	0	1
44	t	0	1
All	All	0	15

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	I	226	G	C2-N3	23.37	1.79	1.32
9	I	226	G	N3-C4	23.05	1.81	1.35
11	K	26	GLY	C-N	20.00	1.59	1.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	I	226	G	N1-C2	17.98	1.73	1.37
9	I	226	G	C6-N1	17.44	1.74	1.39

The worst 5 of 33 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	62	GLY	CA-C-N	18.33	149.36	122.41
3	C	62	GLY	C-N-CA	18.33	149.36	122.41
8	H	8	ASN	CA-C-N	-10.05	106.15	121.83
8	H	8	ASN	C-N-CA	-10.05	106.15	121.83
11	K	26	GLY	N-CA-C	9.88	120.76	111.67

There are no chirality outliers.

5 of 15 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	60	ARG	Sidechain
4	D	12	ASP	Peptide
4	D	13	LYS	Peptide
11	K	63	ILE	Mainchain
11	K	81	GLU	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	805	0	847	335	0
2	B	714	0	737	214	0
3	C	649	0	665	246	0
4	D	649	0	694	312	0
5	E	456	0	481	191	0
6	F	638	0	668	84	0
7	G	665	0	717	142	0
8	H	465	0	493	199	0
9	I	95284	625	47973	16244	0
10	J	2529	0	1281	760	0
11	K	2083	0	2157	881	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
12	L	1565	0	1616	851	0
13	M	1552	0	1619	578	0
14	N	1411	0	1447	634	0
15	O	1323	0	1374	662	0
16	R	1129	0	1162	429	0
17	S	938	0	1012	464	0
18	T	1053	0	1129	357	0
19	U	1075	0	1155	384	0
20	V	960	0	1000	411	0
21	W	1753	0	1783	395	0
22	X	1625	0	1699	298	0
23	Y	1643	0	1710	323	0
24	Z	1105	0	1151	267	0
25	a	817	0	808	243	0
26	b	1182	0	1240	240	0
27	c	979	0	1034	211	0
28	d	1022	0	1073	240	0
29	e	787	0	831	217	0
30	f	877	0	890	394	0
31	g	957	0	1020	239	0
32	h	884	0	944	237	0
33	i	444	0	461	85	0
34	j	409	0	443	206	0
35	k	377	0	418	51	0
36	l	504	0	574	127	0
37	m	302	0	343	210	0
38	n	892	0	926	383	0
39	o	917	0	965	513	0
40	p	947	0	1022	367	0
41	q	816	0	839	337	0
42	r	857	0	922	195	0
43	s	739	0	807	336	0
44	t	780	0	834	454	0
45	u	753	0	780	382	0
46	v	569	0	584	356	0
47	w	625	0	655	200	0
48	x	501	0	534	226	0
49	y	449	0	491	231	0
All	All	140455	625	94008	28535	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 123.

The worst 5 of 28535 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)
3:C:63:GLN:N	9:I:226:G:C4	1.68	1.57
9:I:226:G:N1	9:I:226:G:C6	1.74	1.56
9:I:226:G:N1	9:I:226:G:C2	1.73	1.54
10:J:76:U:C2	10:J:98:A:N7	1.82	1.47
9:I:226:G:C4	9:I:226:G:N3	1.81	1.44

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	
1	A	98/100 (98%)	90 (92%)	8 (8%)	0	100	100
2	B	86/88 (98%)	78 (91%)	6 (7%)	2 (2%)	5	29
3	C	80/82 (98%)	68 (85%)	10 (12%)	2 (2%)	4	27
4	D	78/80 (98%)	66 (85%)	10 (13%)	2 (3%)	4	26
5	E	53/55 (96%)	47 (89%)	5 (9%)	1 (2%)	6	33
6	F	77/79 (98%)	74 (96%)	2 (3%)	1 (1%)	9	41
7	G	83/85 (98%)	77 (93%)	6 (7%)	0	100	100
8	H	54/56 (96%)	51 (94%)	3 (6%)	0	100	100
11	K	269/271 (99%)	216 (80%)	52 (19%)	1 (0%)	30	66
12	L	207/209 (99%)	171 (83%)	30 (14%)	6 (3%)	3	24
13	M	199/201 (99%)	172 (86%)	25 (13%)	2 (1%)	12	47
14	N	175/177 (99%)	147 (84%)	27 (15%)	1 (1%)	21	58
15	O	174/176 (99%)	133 (76%)	38 (22%)	3 (2%)	7	35
16	R	140/142 (99%)	113 (81%)	24 (17%)	3 (2%)	5	30

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
17	S	120/122 (98%)	97 (81%)	20 (17%)	3 (2%)	4	27
18	T	142/144 (99%)	118 (83%)	23 (16%)	1 (1%)	18	55
19	U	133/136 (98%)	119 (90%)	14 (10%)	0	100	100
20	V	118/120 (98%)	100 (85%)	18 (15%)	0	100	100
21	W	222/224 (99%)	211 (95%)	10 (4%)	1 (0%)	24	62
22	X	204/206 (99%)	190 (93%)	12 (6%)	2 (1%)	12	47
23	Y	203/205 (99%)	200 (98%)	3 (2%)	0	100	100
24	Z	148/150 (99%)	127 (86%)	19 (13%)	2 (1%)	9	39
25	a	98/100 (98%)	85 (87%)	11 (11%)	2 (2%)	6	32
26	b	149/151 (99%)	143 (96%)	5 (3%)	1 (1%)	18	55
27	c	127/129 (98%)	116 (91%)	10 (8%)	1 (1%)	16	52
28	d	125/127 (98%)	118 (94%)	6 (5%)	1 (1%)	16	52
29	e	96/98 (98%)	81 (84%)	13 (14%)	2 (2%)	5	30
30	f	115/117 (98%)	94 (82%)	20 (17%)	1 (1%)	14	49
31	g	120/123 (98%)	107 (89%)	13 (11%)	0	100	100
32	h	112/114 (98%)	105 (94%)	6 (5%)	1 (1%)	14	49
33	i	54/56 (96%)	45 (83%)	7 (13%)	2 (4%)	2	20
34	j	48/50 (96%)	41 (85%)	6 (12%)	1 (2%)	5	30
35	k	44/46 (96%)	36 (82%)	7 (16%)	1 (2%)	5	29
36	l	62/64 (97%)	53 (86%)	9 (14%)	0	100	100
37	m	36/38 (95%)	31 (86%)	5 (14%)	0	100	100
38	n	114/116 (98%)	96 (84%)	18 (16%)	0	100	100
39	o	112/114 (98%)	93 (83%)	18 (16%)	1 (1%)	14	49
40	p	115/117 (98%)	97 (84%)	18 (16%)	0	100	100
41	q	101/103 (98%)	79 (78%)	22 (22%)	0	100	100
42	r	108/110 (98%)	94 (87%)	13 (12%)	1 (1%)	14	49
43	s	91/93 (98%)	71 (78%)	19 (21%)	1 (1%)	11	45
44	t	100/102 (98%)	72 (72%)	27 (27%)	1 (1%)	12	47
45	u	92/94 (98%)	80 (87%)	11 (12%)	1 (1%)	11	45
46	v	73/75 (97%)	65 (89%)	8 (11%)	0	100	100
47	w	75/77 (97%)	63 (84%)	12 (16%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
48	x	60/62 (97%)	49 (82%)	10 (17%)	1 (2%)	7	35
49	y	56/58 (97%)	46 (82%)	9 (16%)	1 (2%)	6	34
All	All	5346/5442 (98%)	4625 (86%)	668 (12%)	53 (1%)	15	47

5 of 53 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	C	43	ALA
11	K	101	ARG
12	L	104	VAL
15	O	91	VAL
16	R	81	ILE

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	83/83 (100%)	83 (100%)	0	100	100
2	B	76/76 (100%)	76 (100%)	0	100	100
3	C	65/65 (100%)	65 (100%)	0	100	100
4	D	74/74 (100%)	74 (100%)	0	100	100
5	E	48/48 (100%)	48 (100%)	0	100	100
6	F	70/70 (100%)	70 (100%)	0	100	100
7	G	65/65 (100%)	65 (100%)	0	100	100
8	H	48/48 (100%)	48 (100%)	0	100	100
11	K	216/216 (100%)	216 (100%)	0	100	100
12	L	164/164 (100%)	164 (100%)	0	100	100
13	M	165/165 (100%)	165 (100%)	0	100	100
14	N	148/148 (100%)	148 (100%)	0	100	100
15	O	137/137 (100%)	137 (100%)	0	100	100
16	R	116/116 (100%)	116 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
17	S	103/103 (100%)	103 (100%)	0	100	100
18	T	103/103 (100%)	103 (100%)	0	100	100
19	U	108/108 (100%)	107 (99%)	1 (1%)	70	75
20	V	100/100 (100%)	100 (100%)	0	100	100
21	W	186/186 (100%)	186 (100%)	0	100	100
22	X	170/170 (100%)	170 (100%)	0	100	100
23	Y	172/172 (100%)	172 (100%)	0	100	100
24	Z	113/113 (100%)	113 (100%)	0	100	100
25	a	87/87 (100%)	87 (100%)	0	100	100
26	b	124/124 (100%)	124 (100%)	0	100	100
27	c	104/104 (100%)	104 (100%)	0	100	100
28	d	105/105 (100%)	105 (100%)	0	100	100
29	e	86/86 (100%)	86 (100%)	0	100	100
30	f	90/90 (100%)	90 (100%)	0	100	100
31	g	102/102 (100%)	102 (100%)	0	100	100
32	h	92/92 (100%)	92 (100%)	0	100	100
33	i	47/47 (100%)	47 (100%)	0	100	100
34	j	45/45 (100%)	45 (100%)	0	100	100
35	k	38/38 (100%)	38 (100%)	0	100	100
36	l	51/51 (100%)	51 (100%)	0	100	100
37	m	34/34 (100%)	34 (100%)	0	100	100
38	n	86/86 (100%)	86 (100%)	0	100	100
39	o	99/99 (100%)	99 (100%)	0	100	100
40	p	89/89 (100%)	89 (100%)	0	100	100
41	q	84/84 (100%)	84 (100%)	0	100	100
42	r	93/93 (100%)	93 (100%)	0	100	100
43	s	80/80 (100%)	80 (100%)	0	100	100
44	t	83/83 (100%)	83 (100%)	0	100	100
45	u	78/78 (100%)	78 (100%)	0	100	100
46	v	56/57 (98%)	56 (100%)	0	100	100
47	w	67/67 (100%)	67 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
48	x	54/54 (100%)	54 (100%)	0	100	100
49	y	48/48 (100%)	48 (100%)	0	100	100
All	All	4452/4453 (100%)	4451 (100%)	1 (0%)	100	100

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

Mol	Chain	Res	Type
19	U	13	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 80 such sidechains are listed below:

Mol	Chain	Res	Type
32	h	90	HIS
44	t	44	HIS
33	i	41	HIS
40	p	58	GLN
45	u	78	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
10	J	117/118 (99%)	54 (46%)	2 (1%)
9	I	4435/4438 (99%)	1494 (33%)	119 (2%)
All	All	4552/4556 (99%)	1548 (34%)	121 (2%)

5 of 1548 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
9	I	3	U
9	I	4	U
9	I	8	G
9	I	27	A
9	I	31	A

5 of 121 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
9	I	2363	A
9	I	4174	U

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Mol	Chain	Res	Type
9	I	2832	G
9	I	4096	A
9	I	4325	G

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

35 non-standard protein/DNA/RNA residues are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
9	2MA	I	3997	9	22,25,26	3.78	7 (31%)	32,37,40	3.23	11 (34%)
9	4OC	I	1401	9	20,23,24	3.16	8 (40%)	25,32,35	0.84	0
9	5MC	I	966	9	19,22,23	4.12	9 (47%)	26,32,35	1.02	1 (3%)
9	PSU	I	3405	9	18,21,22	4.57	8 (44%)	21,30,33	2.12	5 (23%)
9	G7M	I	3563	9	23,26,27	2.78	8 (34%)	34,39,42	1.89	8 (23%)
19	4D4	U	81	19	9,11,12	2.74	3 (33%)	7,13,15	0.82	0
9	PSU	I	3411	9	18,21,22	4.61	8 (44%)	21,30,33	1.94	6 (28%)
9	PSU	I	4074	9	18,21,22	4.37	9 (50%)	21,30,33	2.18	5 (23%)
9	OMC	I	3992	9	19,22,23	3.34	8 (42%)	25,31,34	0.81	0
9	G7M	I	526	9	23,26,27	2.85	9 (39%)	34,39,42	1.74	8 (23%)
31	D2T	g	88	31	8,9,10	2.93	1 (12%)	6,11,13	2.42	2 (33%)
9	5MC	I	3456	9	19,22,23	3.89	9 (47%)	26,32,35	1.11	1 (3%)
9	UR3	I	4402	9	19,22,23	2.86	8 (42%)	26,32,35	1.61	4 (15%)
9	PSU	I	2449	9	18,21,22	4.36	8 (44%)	21,30,33	2.23	5 (23%)
9	2MG	I	3329	9	23,26,27	2.94	9 (39%)	33,38,41	2.27	9 (27%)
9	PSU	I	4099	9	18,21,22	4.39	9 (50%)	21,30,33	2.07	7 (33%)
9	MA6	I	4423	9	23,26,27	1.35	4 (17%)	33,38,41	4.77	13 (39%)
9	3TD	I	3409	9	19,22,23	4.14	6 (31%)	23,32,35	1.74	3 (13%)
9	1MG	I	2245	9	23,26,27	2.77	8 (34%)	33,39,42	1.53	5 (15%)
9	PSU	I	3998	9	18,21,22	4.50	9 (50%)	21,30,33	2.19	5 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
9	2MG	I	3939	9	23,26,27	2.81	9 (39%)	33,38,41	2.35	9 (27%)
9	2MG	I	4420	9	23,26,27	2.98	9 (39%)	33,38,41	2.18	10 (30%)
9	PSU	I	515	9	18,21,22	4.58	8 (44%)	21,30,33	2.18	6 (28%)
9	2MG	I	1206	9	23,26,27	3.01	8 (34%)	33,38,41	2.16	10 (30%)
9	6MZ	I	3112	9	22,25,26	2.59	4 (18%)	29,36,39	2.30	9 (31%)
9	5MU	I	2247	9	19,22,23	4.64	6 (31%)	27,32,35	3.79	10 (37%)
9	5MU	I	3433	9	19,22,23	4.60	6 (31%)	27,32,35	3.84	10 (37%)
9	6MZ	I	3524	9	22,25,26	2.51	5 (22%)	29,36,39	2.45	12 (41%)
9	OMG	I	3745	9	23,26,27	2.58	9 (39%)	32,38,41	1.97	9 (28%)
9	OMU	I	4046	9	19,22,23	2.86	6 (31%)	25,31,34	1.83	4 (16%)
9	2MG	I	965	9	23,26,27	3.01	9 (39%)	33,38,41	2.15	10 (30%)
9	PSU	I	2246	9	18,21,22	4.61	8 (44%)	21,30,33	2.51	6 (28%)
9	PSU	I	3951	9	18,21,22	4.33	8 (44%)	21,30,33	2.54	7 (33%)
9	5MC	I	1406	9	19,22,23	3.99	9 (47%)	26,32,35	1.00	1 (3%)
9	MA6	I	4422	9	23,26,27	1.35	4 (17%)	33,38,41	4.72	13 (39%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	2MA	I	3997	9	-	4/7/25/26	0/3/3/3
9	4OC	I	1401	9	-	1/9/29/30	0/2/2/2
9	5MC	I	966	9	-	1/7/25/26	0/2/2/2
9	PSU	I	3405	9	-	0/7/25/26	0/2/2/2
9	G7M	I	3563	9	-	1/7/25/26	0/3/3/3
19	4D4	U	81	19	-	5/11/12/14	-
9	PSU	I	3411	9	-	1/7/25/26	0/2/2/2
9	PSU	I	4074	9	-	2/7/25/26	0/2/2/2
9	OMC	I	3992	9	-	4/9/27/28	0/2/2/2
9	G7M	I	526	9	-	3/7/25/26	0/3/3/3
31	D2T	g	88	31	-	3/7/12/14	-
9	5MC	I	3456	9	-	4/7/25/26	0/2/2/2
9	UR3	I	4402	9	-	4/7/25/26	0/2/2/2
9	PSU	I	2449	9	-	2/7/25/26	0/2/2/2
9	2MG	I	3329	9	-	1/9/27/28	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
9	PSU	I	4099	9	-	5/7/25/26	0/2/2/2
9	MA6	I	4423	9	-	1/11/29/30	0/3/3/3
9	3TD	I	3409	9	-	1/7/25/26	0/2/2/2
9	1MG	I	2245	9	-	2/7/25/26	0/3/3/3
9	PSU	I	3998	9	-	2/7/25/26	0/2/2/2
9	2MG	I	3939	9	-	2/9/27/28	0/3/3/3
9	2MG	I	4420	9	-	1/9/27/28	0/3/3/3
9	PSU	I	515	9	-	0/7/25/26	0/2/2/2
9	2MG	I	1206	9	-	1/9/27/28	0/3/3/3
9	6MZ	I	3112	9	-	6/9/27/28	0/3/3/3
9	5MU	I	2247	9	-	2/7/25/26	0/2/2/2
9	5MU	I	3433	9	-	1/7/25/26	0/2/2/2
9	6MZ	I	3524	9	-	4/9/27/28	0/3/3/3
9	OMG	I	3745	9	-	0/9/27/28	0/3/3/3
9	OMU	I	4046	9	-	4/9/27/28	0/2/2/2
9	2MG	I	965	9	-	2/9/27/28	0/3/3/3
9	PSU	I	2246	9	-	3/7/25/26	0/2/2/2
9	PSU	I	3951	9	-	2/7/25/26	0/2/2/2
9	5MC	I	1406	9	-	4/7/25/26	0/2/2/2
9	MA6	I	4422	9	-	4/11/29/30	0/3/3/3

The worst 5 of 256 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
9	I	3409	3TD	C6-C5	12.54	1.49	1.35
9	I	3411	PSU	C6-C5	12.31	1.48	1.35
9	I	2246	PSU	C6-C5	12.21	1.48	1.35
9	I	515	PSU	C6-C5	11.97	1.48	1.35
9	I	3405	PSU	C6-C5	11.97	1.48	1.35

The worst 5 of 224 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	I	4422	MA6	N1-C6-N6	-17.74	95.25	116.86
9	I	4423	MA6	N1-C6-N6	-17.65	95.36	116.86
9	I	3433	5MU	C5-C4-N3	12.34	126.05	115.32
9	I	4422	MA6	C5-C6-N6	12.20	144.65	125.33
9	I	2247	5MU	C5-C4-N3	12.12	125.86	115.32

There are no chirality outliers.

5 of 83 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
19	U	81	4D4	O-C-CA-CB
19	U	81	4D4	N-CA-CB-CG
19	U	81	4D4	NE-CD-CG-CB
31	g	88	D2T	CA-CB-SB-CB1
31	g	88	D2T	CG-CB-SB-CB1

There are no ring outliers.

34 monomers are involved in 182 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	I	3997	2MA	5	0
9	I	1401	4OC	7	0
9	I	966	5MC	9	0
9	I	3405	PSU	6	0
9	I	3563	G7M	1	0
9	I	3411	PSU	7	0
9	I	4074	PSU	5	0
9	I	3992	OMC	7	0
9	I	526	G7M	6	0
31	g	88	D2T	6	0
9	I	3456	5MC	6	0
9	I	4402	UR3	3	0
9	I	2449	PSU	9	0
9	I	3329	2MG	3	0
9	I	4099	PSU	6	0
9	I	4423	MA6	8	0
9	I	3409	3TD	7	0
9	I	2245	1MG	12	0
9	I	3998	PSU	3	0
9	I	3939	2MG	8	0
9	I	4420	2MG	7	0
9	I	515	PSU	3	0
9	I	1206	2MG	5	0
9	I	3112	6MZ	5	0
9	I	2247	5MU	1	0
9	I	3433	5MU	4	0
9	I	3524	6MZ	4	0
9	I	3745	OMG	6	0
9	I	4046	OMU	6	0
9	I	965	2MG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
9	I	2246	PSU	3	0
9	I	3951	PSU	7	0
9	I	1406	5MC	6	0
9	I	4422	MA6	8	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

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
9	I	2

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	I	1500:U	O3'	1501:G	P	17.28
1	I	2385:C	O3'	2386:A	P	15.64

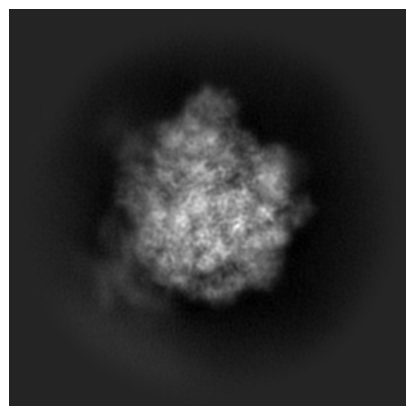
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26666. These allow visual inspection of the internal detail of the map and identification of artifacts.

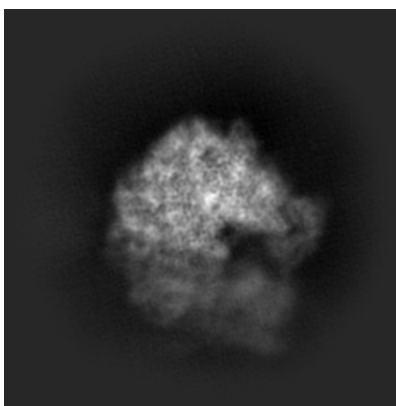
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

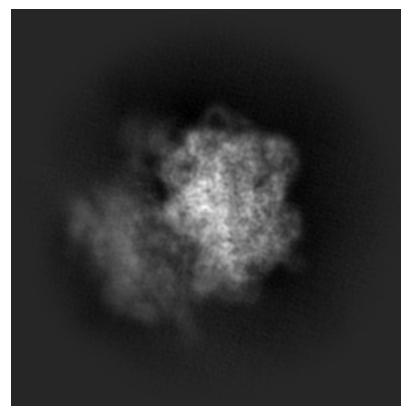
6.1.1 Primary map



X

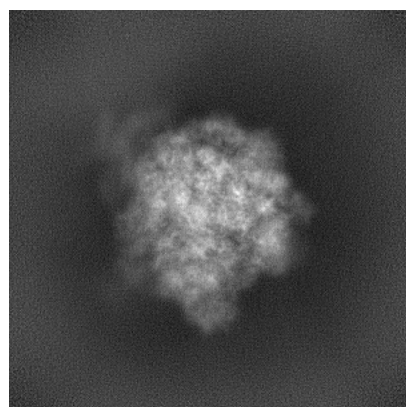


Y

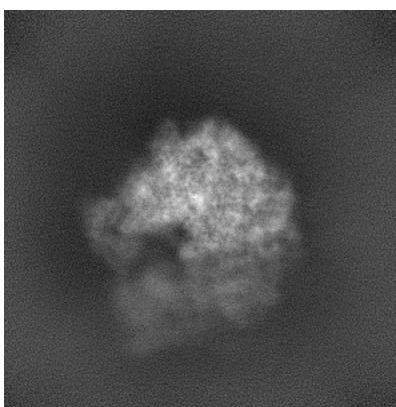


Z

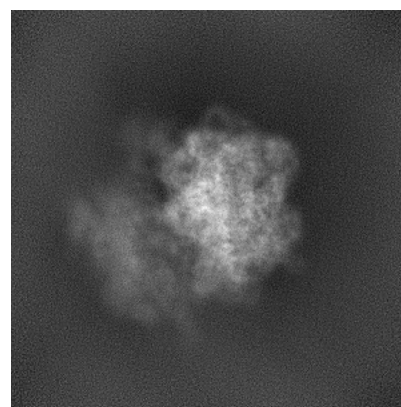
6.1.2 Raw map



X



Y

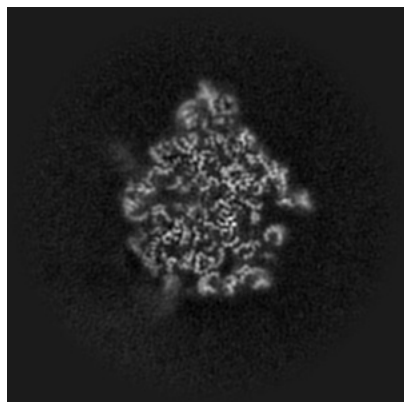


Z

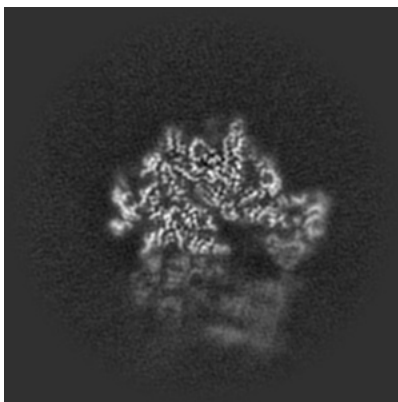
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

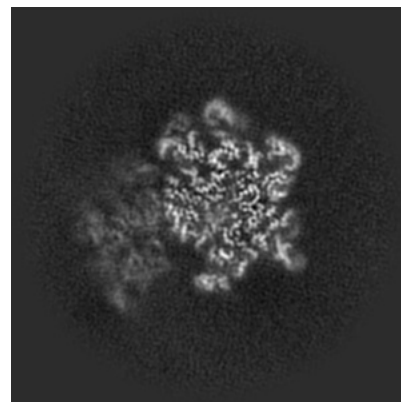
6.2.1 Primary map



X Index: 250

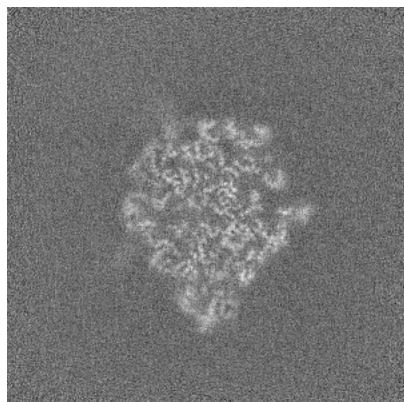


Y Index: 250

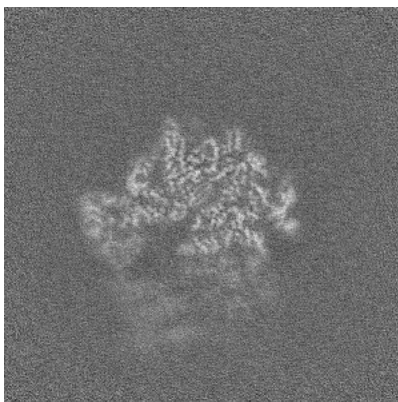


Z Index: 250

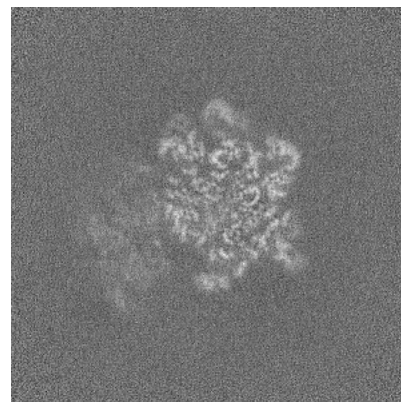
6.2.2 Raw map



X Index: 250



Y Index: 250

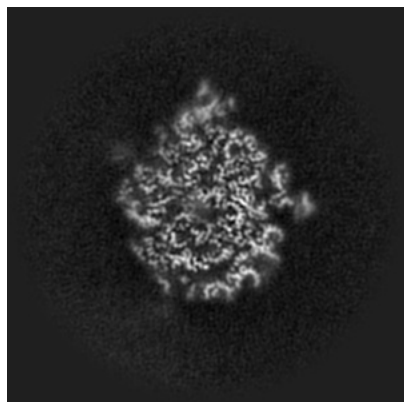


Z Index: 250

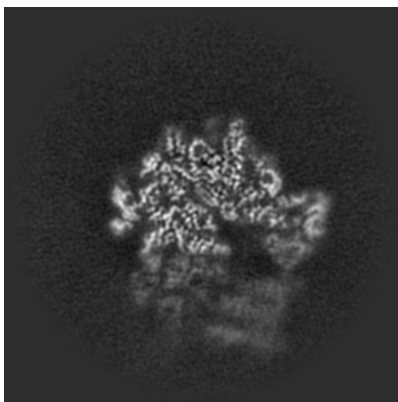
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

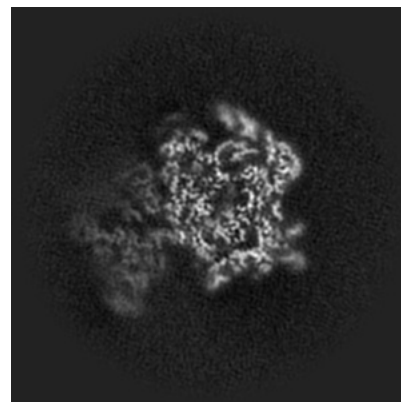
6.3.1 Primary map



X Index: 259

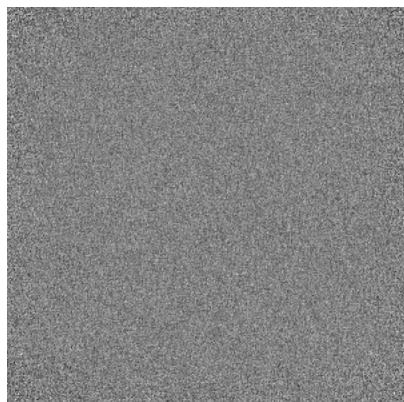


Y Index: 249

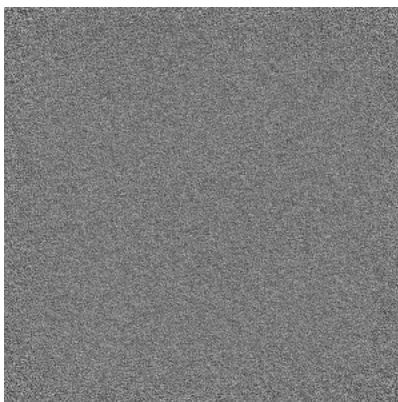


Z Index: 239

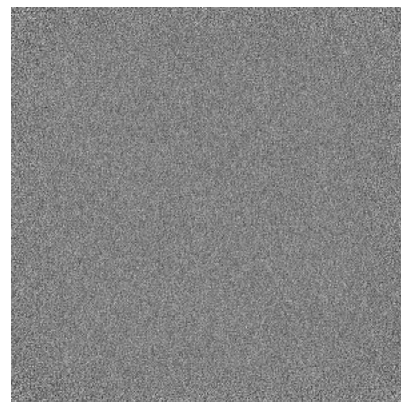
6.3.2 Raw map



X Index: 0



Y Index: 0

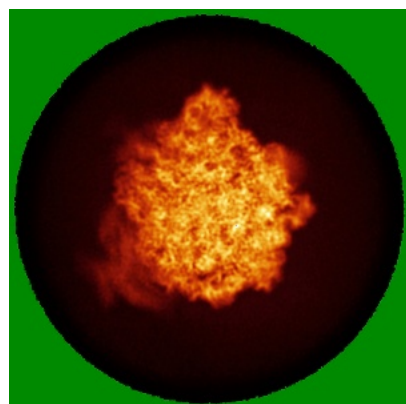


Z Index: 0

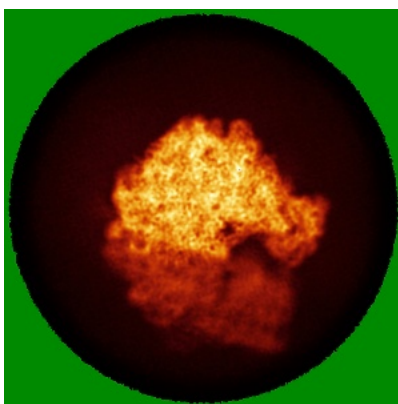
The images above show the largest variance slices of the map in three orthogonal directions.

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

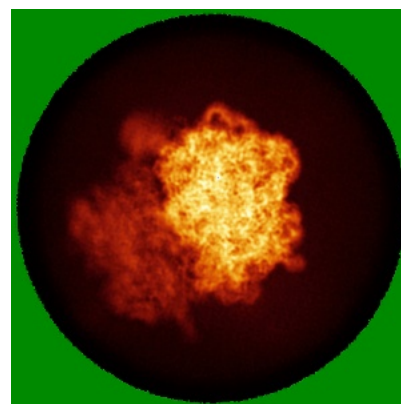
6.4.1 Primary map



X

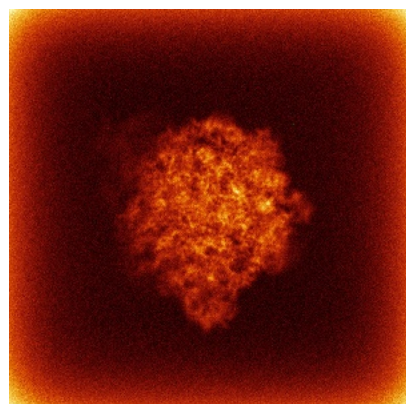


Y

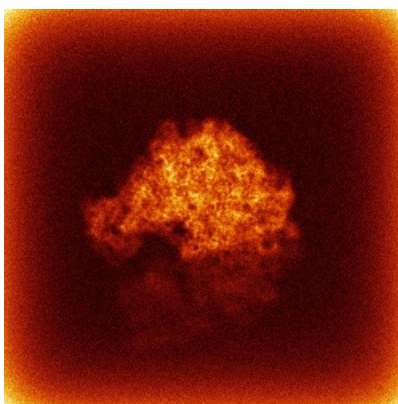


Z

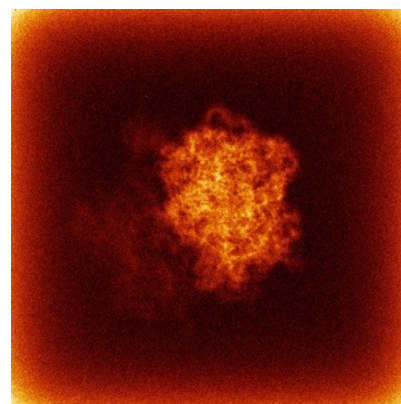
6.4.2 Raw map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



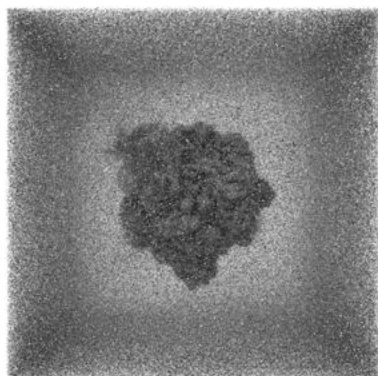
Y



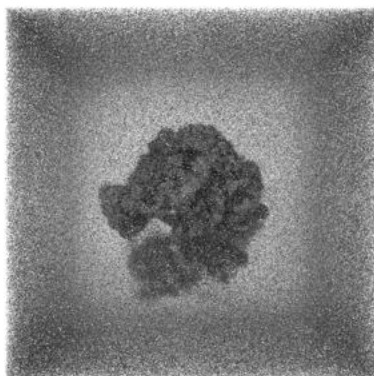
Z

The images above show the 3D surface view of the map at the recommended contour level 0.1. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

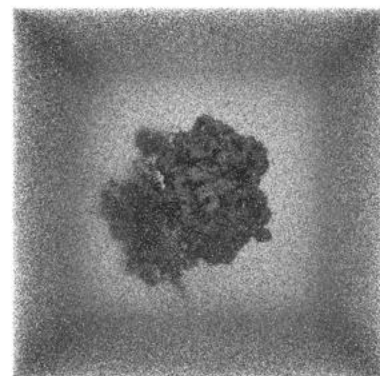
6.5.2 Raw map



X



Y



Z

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

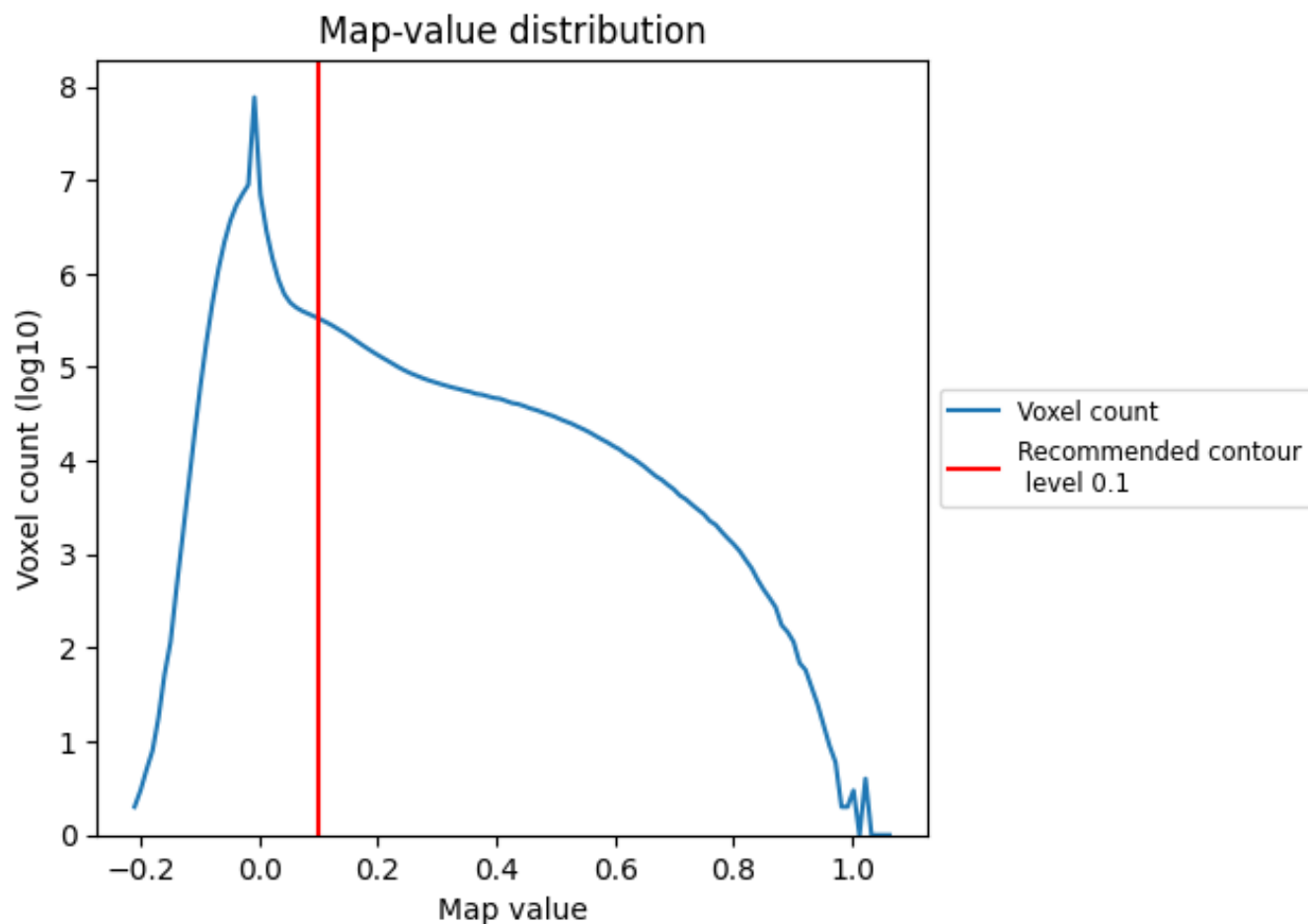
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

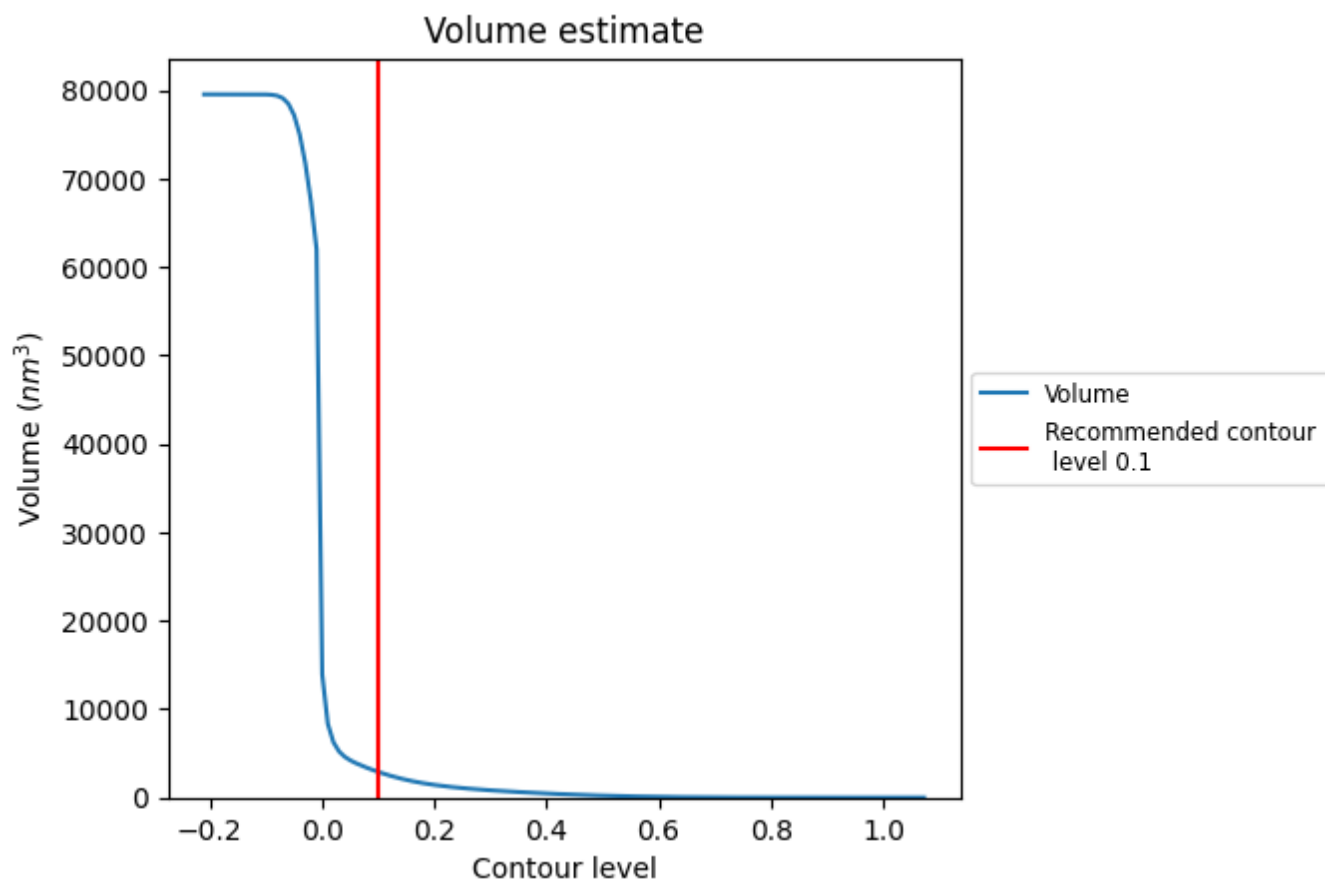
This section contains the results of statistical analysis of the map.

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.

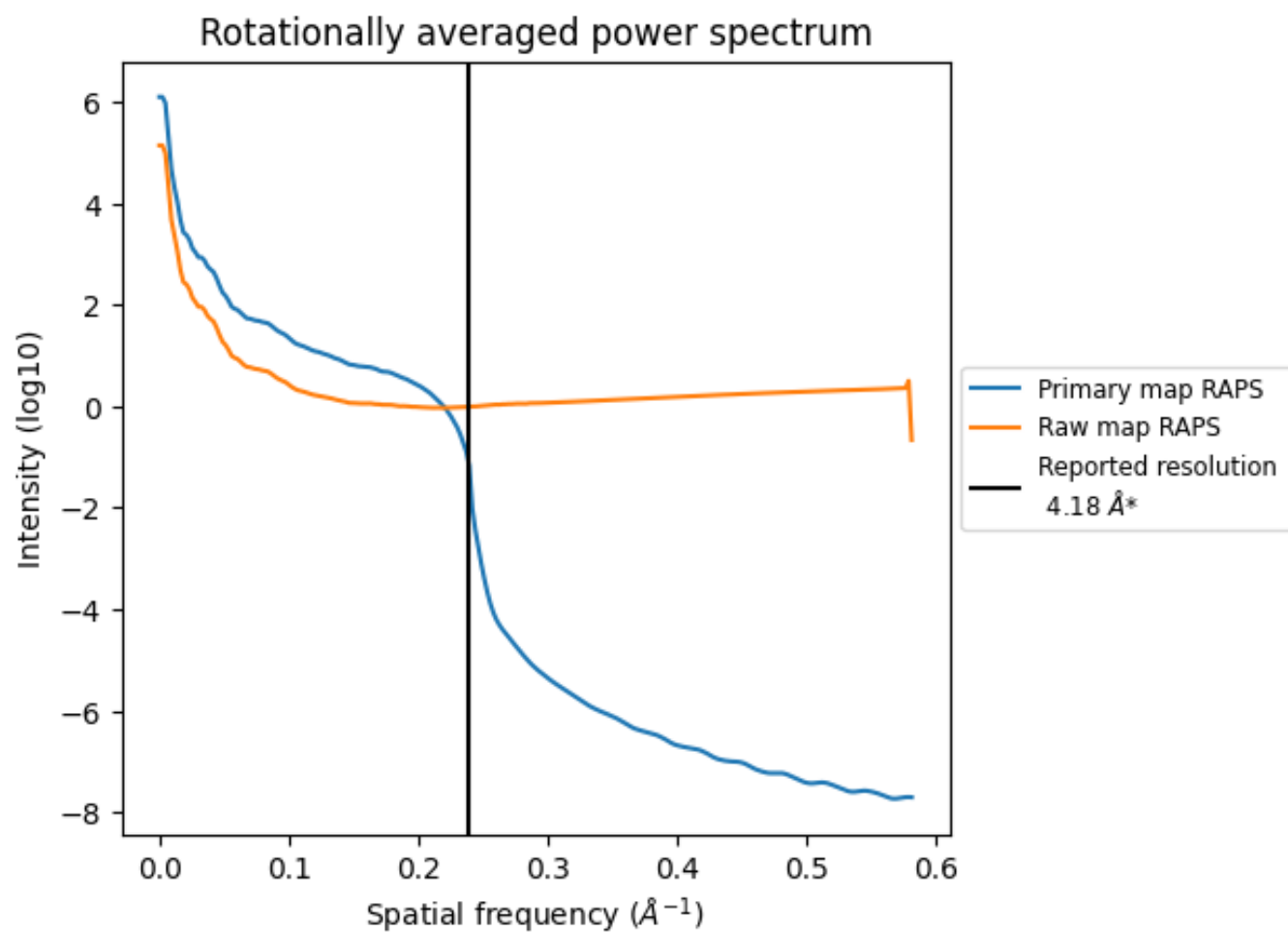
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 2909 nm³; this corresponds to an approximate mass of 2628 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 ⓘ

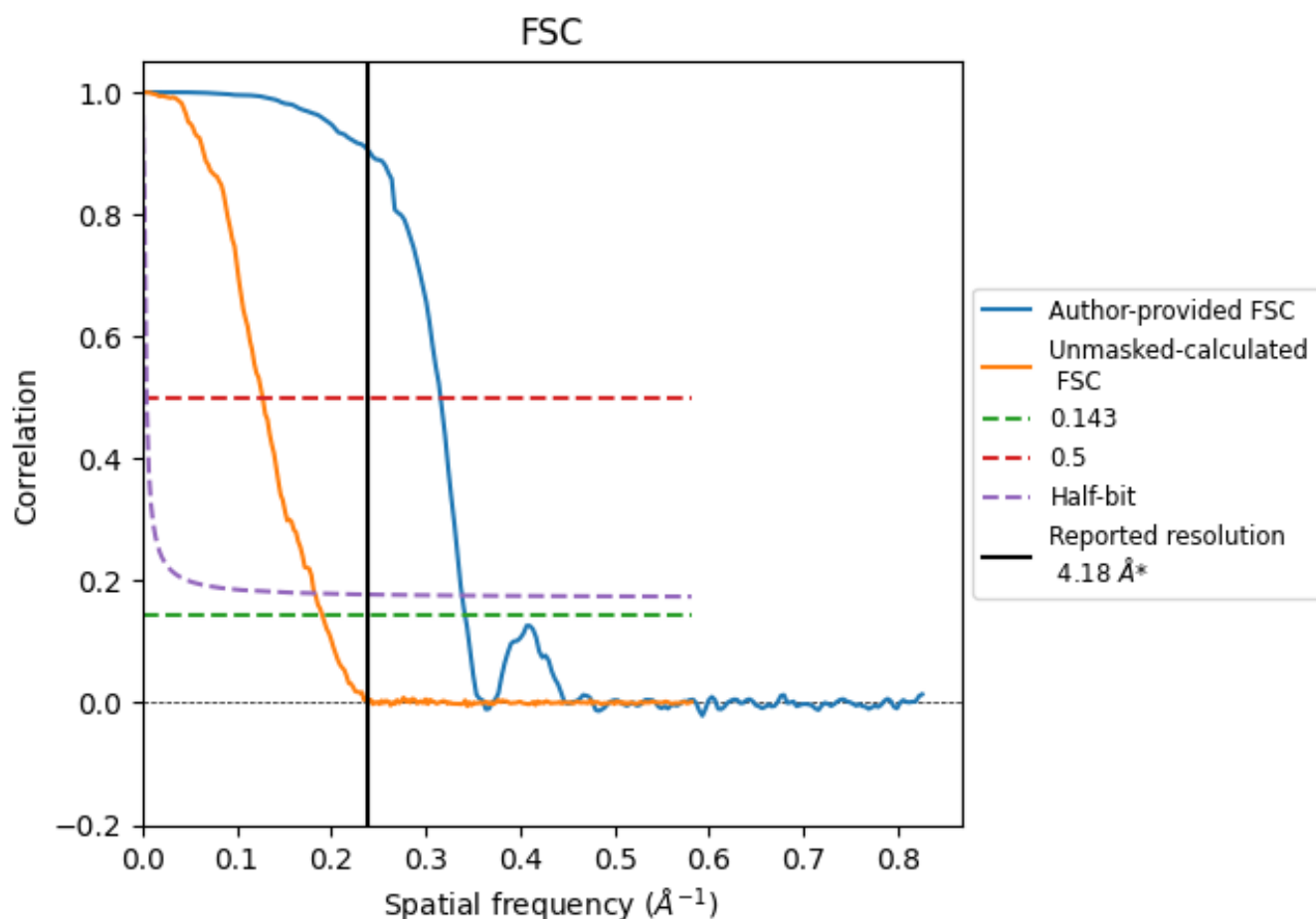


*Reported resolution corresponds to spatial frequency of 0.239 $\text{\AA}^{-1}$

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.239 Å⁻¹

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	4.18	-	-
Author-provided FSC curve	2.93	3.17	2.96
Unmasked-calculated*	5.24	7.85	5.47

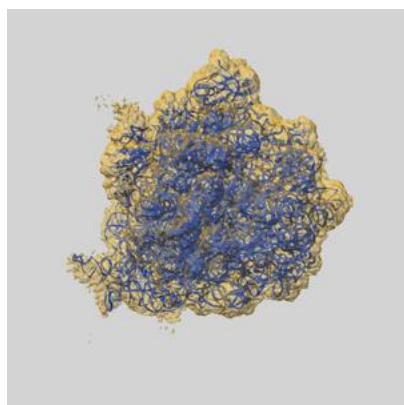
*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from author-provided FSC intersecting FSC 0.143 CUT-OFF 2.93 differs from the reported value 4.18 by more than 10 %

The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 5.24 differs from the reported value 4.18 by more than 10 %

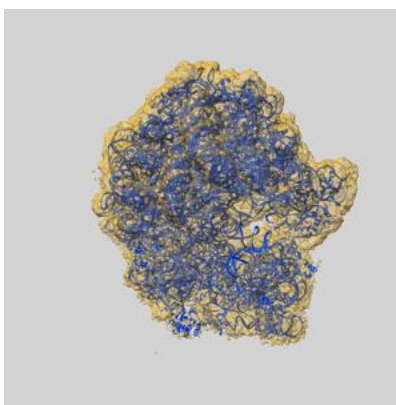
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-26666 and PDB model 7UPH. Per-residue inclusion information can be found in section 3 on page 14.

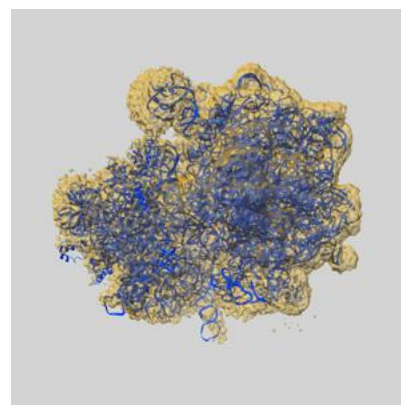
9.1 Map-model overlay [i](#)



X



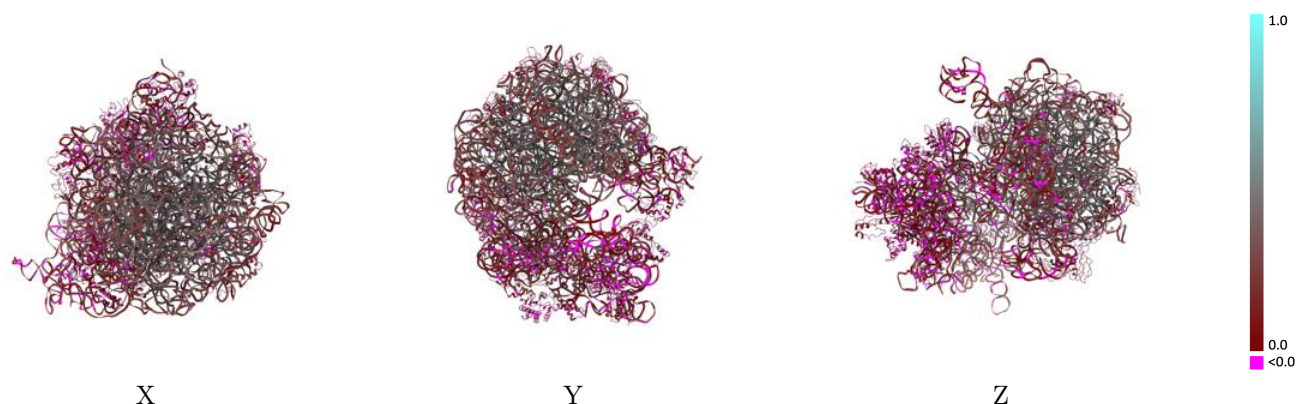
Y



Z

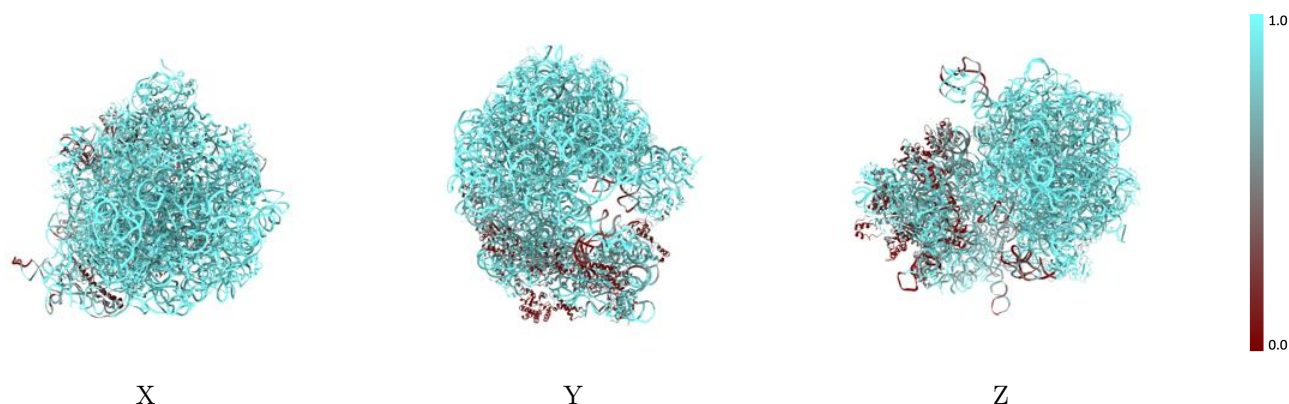
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



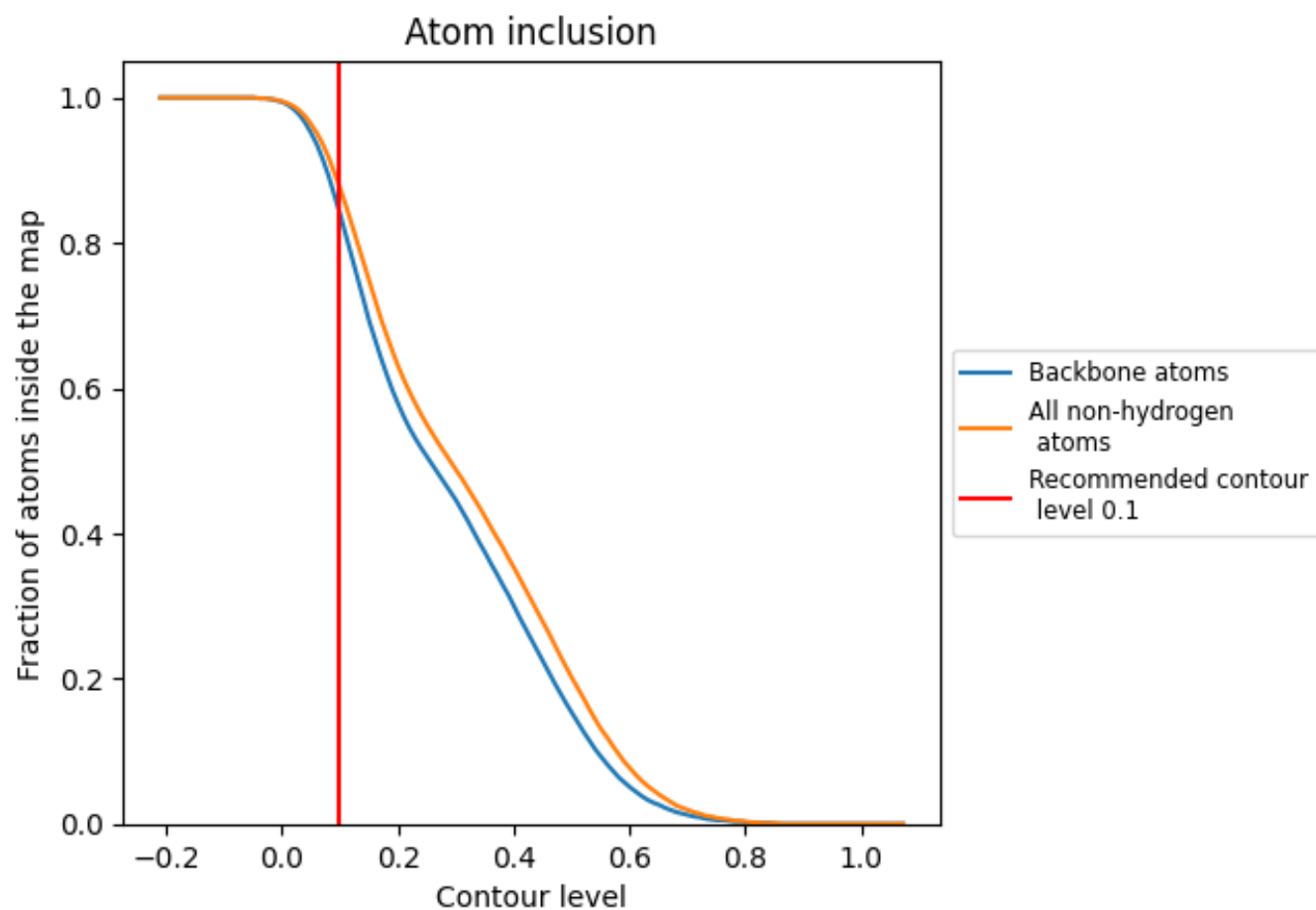
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.1).

9.4 Atom inclusion ⓘ



At the recommended contour level, 84% of all backbone atoms, 88% 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 (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8770	 0.2400
A	 0.7610	 0.0710
B	 0.4930	 0.1220
C	 0.7030	 0.1010
D	 0.4990	 0.1100
E	 0.5420	 0.1030
F	 0.2240	 0.0820
G	 0.2720	 0.0490
H	 0.2070	 0.0660
I	 0.9290	 0.2690
J	 0.9840	 0.1870
K	 0.9670	 0.2620
L	 0.9720	 0.2320
M	 0.9860	 0.2770
N	 0.9030	 0.1100
O	 0.9240	 0.1440
R	 0.9800	 0.2660
S	 0.9100	 0.1770
T	 0.9770	 0.2820
U	 0.9750	 0.2430
V	 0.9650	 0.2620
W	 0.1530	 0.0940
X	 0.4000	 0.0950
Y	 0.4910	 0.0880
Z	 0.7220	 0.1790
a	 0.3730	 0.1020
b	 0.3680	 0.1200
c	 0.6360	 0.1640
d	 0.7370	 0.0920
e	 0.4960	 0.1210
f	 0.4700	 0.0950
g	 0.7480	 0.2300
h	 0.5410	 0.1190
i	 0.9860	 0.3590
j	 0.8650	 0.1050



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Chain	Atom inclusion	Q-score
k	 0.9920	 0.4020
l	 0.9880	 0.3730
m	 0.9110	 0.1150
n	 0.9290	 0.1460
o	 0.8480	 0.1160
p	 0.9780	 0.2860
q	 0.9770	 0.2520
r	 0.9900	 0.3640
s	 0.9500	 0.2180
t	 0.9600	 0.1780
u	 0.9380	 0.1380
v	 0.9300	 0.1640
w	 0.9770	 0.2910
x	 0.9960	 0.2010
y	 0.9950	 0.2680