



wwPDB EM Validation Summary Report ⓘ

Mar 23, 2026 – 02:00 AM UTC

PDB ID : 9BDL / pdb_00009bdl
EMDB ID : EMD-44461
Title : 80S ribosome with angiogenin
Authors : Loveland, A.B.; Korostelev, A.A.
Deposited on : 2024-04-12
Resolution : 2.80 Å(reported)

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:

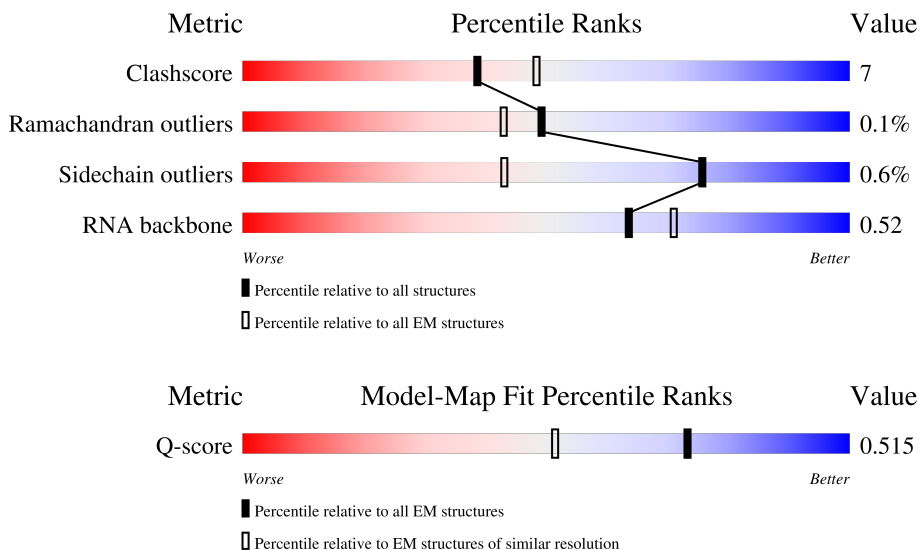
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

The reported resolution of this entry is 2.80 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.




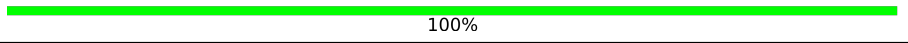
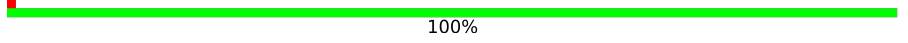
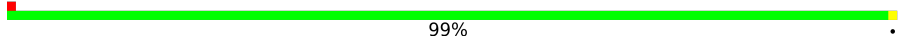
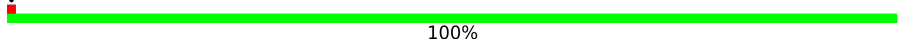

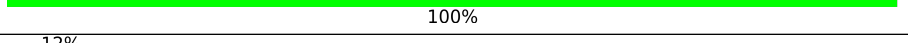
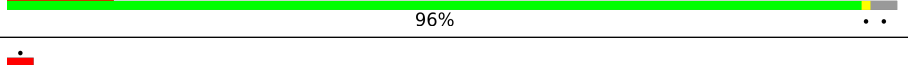
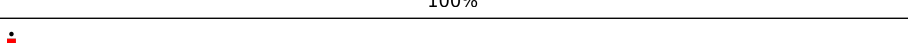
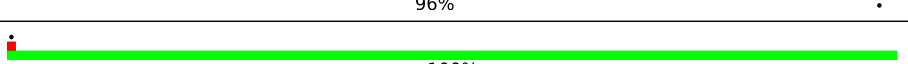
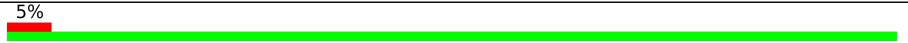
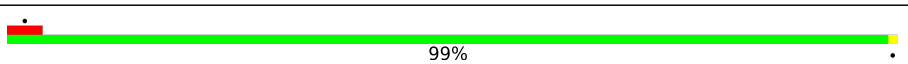
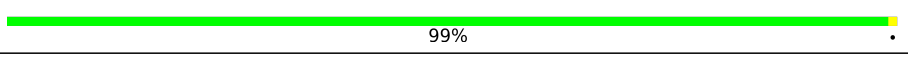
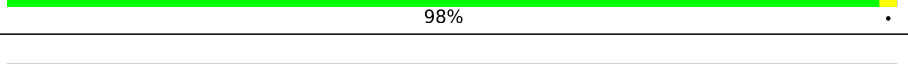
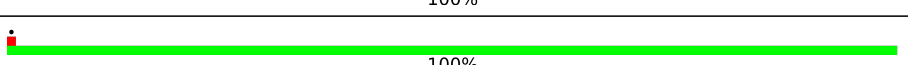
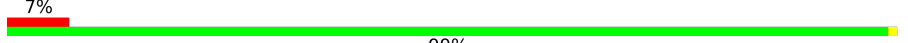
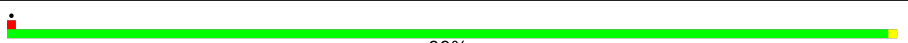
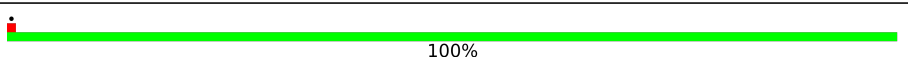
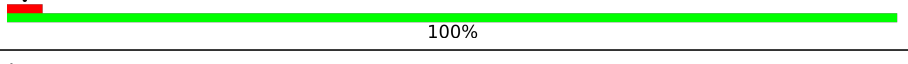
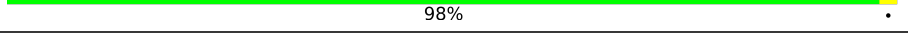

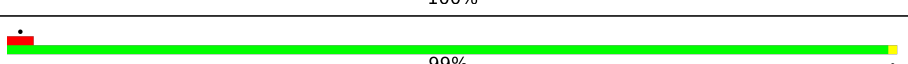



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	11806 (2.30 - 3.30)

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	A18S	1869	
2	A28S	3601	
3	A58S	156	

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Mol	Chain	Length	Quality of chain
4	A5S	120	 88% 12%
5	AL02	248	 100%
6	AL03	394	 100%
7	AL04	362	 99%
8	AL05	293	 100%
9	AL06	251	 85% 14%
10	AL07	225	 100%
11	AL08	240	 12% 96%
12	AL09	190	 100%
13	AL10	213	 96%
14	AL11	170	 100%
15	AL13	210	 5% 100%
16	AL14	138	 99%
17	AL15	203	 99%
18	AL16	199	 98%
19	AL17	153	 100%
20	AL18	187	 100%
21	AL19	180	 7% 99%
22	AL20	176	 99%
23	AL21	159	 100%
24	AL22	99	 100%
25	AL23	131	 98%
26	AL24	121	 52% 48%
27	AL25	118	 100%
28	AL26	134	 99%

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Mol	Chain	Length	Quality of chain
29	AL27	135	100%
30	AL28	147	100%
31	AL29	245	7% 42% 58%
32	AL30	98	100%
33	AL31	107	99%
34	AL32	128	100%
35	AL33	109	100%
36	AL34	114	5% 98%
37	AL35	122	98%
38	AL36	102	99%
39	AL37	86	100%
40	AL38	69	100%
41	AL39	50	98%
42	AL40	52	100%
43	AL41	25	100%
44	AL42	104	98%
45	AL43	91	99%
46	ALNW	124	99%
47	ANG	147	7% 66% 15% 18%
48	ARAC	313	52% 99%
49	AS00	217	10% 100%
50	AS01	213	100%
51	AS02	221	7% 100%
52	AS03	228	18% 100%
53	AS04	262	5% 100%

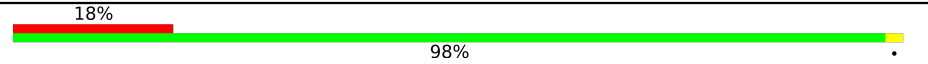

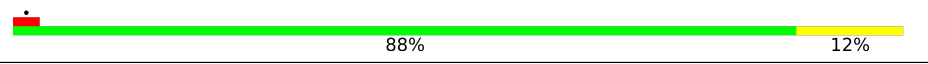

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Mol	Chain	Length	Quality of chain
54	AS05	191	5% 97%
55	AS06	237	23% 98%
56	AS07	189	32% 98%
57	AS08	206	11% 99%
58	AS09	185	8% 100%
59	AS10	96	27% 99%
60	AS11	151	8% 95% 5%
61	AS12	117	85% 84% 15%
62	AS13	149	100%
63	AS14	135	6% 100%
64	AS15	120	12% 99%
65	AS16	142	8% 100%
66	AS17	132	33% 100%
67	AS18	144	8% 99%
68	AS19	141	9% 99%
69	AS20	100	34% 99%
70	AS21	83	11% 100%
71	AS22	129	99%
72	AS23	141	100%
73	AS24	124	6% 100%
74	AS25	75	9% 99%
75	AS26	101	99%
76	AS27	83	12% 99%
77	AS28	62	10% 100%
78	AS29	55	22% 100%

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Mol	Chain	Length	Quality of chain
79	AS30	55	 <p>18% 98% 14%</p>
80	ETRN	77	 <p>9% 86% 14%</p>
80	PTRN	77	 <p>11% 88% 12%</p>
81	MRNA	28	 <p>11% 18% 18% 64%</p>

2 Entry composition

There are 81 unique types of molecules in this entry. The entry contains 212927 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 18S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	A18S	1691	36103	16115	6485	11813	1690	0	0

- Molecule 2 is a RNA chain called 28S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
2	A28S	3557	76286	33975	13979	24775	3557	0	0

- Molecule 3 is a RNA chain called 5.8S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
3	A58S	151	3208	1432	564	1062	150	0	0

- Molecule 4 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
4	A5S	120	2558	1141	456	842	119	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A5S	2	U	N	conflict	GB X06789.1
A5S	36	C	N	conflict	GB X06789.1
A5S	102	U	N	conflict	GB X06789.1
A5S	112	U	N	conflict	GB X06789.1
A5S	114	U	N	conflict	GB X06789.1
A5S	119	U	C	conflict	GB X06789.1
A5S	120	U	N	conflict	GB X06789.1

- Molecule 5 is a protein called 60S ribosomal protein L8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	AL02	248	1898	1189	389	314	6	0	0

- Molecule 6 is a protein called 60S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	AL03	394	3172	2020	597	542	13	0	0

- Molecule 7 is a protein called 60S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	AL04	362	2883	1812	577	480	14	0	0

- Molecule 8 is a protein called Large ribosomal subunit protein uL18.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	AL05	293	2391	1512	438	427	14	0	0

- Molecule 9 is a protein called 60S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	AL06	216	1729	1115	329	282	3	0	0

- Molecule 10 is a protein called 60S ribosomal protein L7.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	AL07	225	1875	1205	358	303	9	0	0

- Molecule 11 is a protein called Large ribosomal subunit protein eL8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
11	AL08	233	1879	1199	361	315	4	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AL08	244	GLY	CYS	conflict	UNP G1STW0

- Molecule 12 is a protein called 60S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
12	AL09	190	1516	954	284	272	6	0	0

- Molecule 13 is a protein called Ribosomal protein L10.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
13	AL10	205	1664	1056	321	274	13	0	0

- Molecule 14 is a protein called 60S ribosomal protein L11.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
14	AL11	170	1361	861	254	240	6	0	0

- Molecule 15 is a protein called Large ribosomal subunit protein eL13.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
15	AL13	210	1702	1065	354	279	4	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AL13	74	ARG	HIS	conflict	UNP G1TKB3
AL13	190	ARG	HIS	conflict	UNP G1TKB3

- Molecule 16 is a protein called Large ribosomal subunit protein eL14.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
16	AL14	138	1137	727	221	182	7	0	0

- Molecule 17 is a protein called 60S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
17	AL15	203	1701	1072	359	266	4	0	0

- Molecule 18 is a protein called Large ribosomal subunit protein uL13.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
18	AL16	199	1630	1051	319	255	5	0	0

- Molecule 19 is a protein called 60S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
19	AL17	153	1242	777	241	215	9	0	0

- Molecule 20 is a protein called Large ribosomal subunit protein eL18.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
20	AL18	187	1515	946	315	250	4	0	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AL18	4	ASP	ASN	conflict	UNP G1TFE0
AL18	14	ARG	TRP	conflict	UNP G1TFE0
AL18	53	MET	LEU	conflict	UNP G1TFE0
AL18	58	ARG	TRP	conflict	UNP G1TFE0
AL18	75	ARG	GLN	conflict	UNP G1TFE0
AL18	80	ALA	PRO	conflict	UNP G1TFE0
AL18	86	VAL	ILE	conflict	UNP G1TFE0
AL18	104	ARG	HIS	conflict	UNP G1TFE0
AL18	110	ARG	CYS	conflict	UNP G1TFE0
AL18	137	VAL	GLY	conflict	UNP G1TFE0
AL18	157	GLY	ARG	conflict	UNP G1TFE0
AL18	181	ARG	TRP	conflict	UNP G1TFE0

- Molecule 21 is a protein called 60S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
21	AL19	180	1508	933	328	238	9	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AL19	38	ARG	HIS	conflict	UNP G1TYL6
AL19	151	ARG	HIS	conflict	UNP G1TYL6

- Molecule 22 is a protein called 60S ribosomal protein L18a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
22	AL20	176	1462	930	285	236	11	0	0

There are 23 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AL20	1	MET	THR	conflict	UNP G1TTY7
AL20	18	PRO	-	insertion	UNP G1TTY7
AL20	19	THR	-	insertion	UNP G1TTY7
AL20	20	PRO	SER	conflict	UNP G1TTY7
AL20	22	CYS	SER	conflict	UNP G1TTY7
AL20	23	ARG	PRO	conflict	UNP G1TTY7
AL20	24	THR	ALA	conflict	UNP G1TTY7
AL20	49	SER	LEU	conflict	UNP G1TTY7
AL20	50	GLN	GLU	conflict	UNP G1TTY7
AL20	95	ARG	HIS	conflict	UNP G1TTY7
AL20	101	THR	ILE	conflict	UNP G1TTY7
AL20	102	THR	MET	conflict	UNP G1TTY7
AL20	104	GLY	SER	conflict	UNP G1TTY7
AL20	126	ILE	VAL	conflict	UNP G1TTY7
AL20	132	ILE	MET	conflict	UNP G1TTY7
AL20	135	SER	ALA	conflict	UNP G1TTY7
AL20	136	LYS	ARG	conflict	UNP G1TTY7
AL20	138	ARG	PRO	conflict	UNP G1TTY7
AL20	149	LYS	ARG	conflict	UNP G1TTY7
AL20	151	LYS	ARG	conflict	UNP G1TTY7
AL20	168	THR	TYR	conflict	UNP G1TTY7
AL20	169	THR	ALA	conflict	UNP G1TTY7
AL20	176	PHE	-	insertion	UNP G1TTY7

- Molecule 23 is a protein called 60S ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
23	AL21	159	1298	823	252	217	6	0	0

- Molecule 24 is a protein called Large ribosomal subunit protein eL22.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
24	AL22	99	809	519	141	147	2	0	0

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AL22	18	LEU	VAL	conflict	UNP G1TSG1
AL22	32	GLY	ARG	conflict	UNP G1TSG1
AL22	36	ALA	GLU	conflict	UNP G1TSG1
AL22	39	PHE	SER	conflict	UNP G1TSG1
AL22	54	GLY	ARG	conflict	UNP G1TSG1
AL22	60	VAL	ALA	conflict	UNP G1TSG1
AL22	62	SER	THR	conflict	UNP G1TSG1
AL22	63	LEU	ILE	conflict	UNP G1TSG1
AL22	97	ARG	HIS	conflict	UNP G1TSG1
AL22	106	THR	SER	conflict	UNP G1TSG1

- Molecule 25 is a protein called 60S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
25	AL23	131	979	618	184	172	5	0	0

- Molecule 26 is a protein called Ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
26	AL24	63	528	337	103	85	3	0	0

- Molecule 27 is a protein called Large ribosomal subunit protein uL23.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
27	AL25	118	967	618	181	167	1	0	0

- Molecule 28 is a protein called 60S ribosomal protein L26.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
28	AL26	134	1115	700	226	186	3	0	0

- Molecule 29 is a protein called 60S ribosomal protein L27.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
29	AL27	135	1107	714	208	182	3	0	0

- Molecule 30 is a protein called 60S ribosomal protein L27a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
30	AL28	147	1162	734	239	185	4	0	0

- Molecule 31 is a protein called Large ribosomal subunit protein eL29.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
31	AL29	104	848	527	189	129	3	0	0

- Molecule 32 is a protein called 60S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
32	AL30	98	761	481	134	140	6	0	0

- Molecule 33 is a protein called 60S ribosomal protein L31.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
33	AL31	107	888	560	171	155	2	0	0

- Molecule 34 is a protein called Large ribosomal subunit protein eL32.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
34	AL32	128	1053	667	216	165	5	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AL32	3	ALA	SER	conflict	UNP G1TUN8
AL32	13	VAL	ILE	conflict	UNP G1TUN8
AL32	16	ARG	TRP	conflict	UNP G1TUN8
AL32	81	ASN	SER	conflict	UNP G1TUN8

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Chain	Residue	Modelled	Actual	Comment	Reference
AL32	98	GLU	LYS	conflict	UNP G1TUN8
AL32	108	ARG	CYS	conflict	UNP G1TUN8
AL32	115	ALA	VAL	conflict	UNP G1TUN8

- Molecule 35 is a protein called 60S ribosomal protein L35a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
35	AL33	109	876	555	174	143	4	0	0

- Molecule 36 is a protein called 60S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
36	AL34	114	906	566	187	147	6	0	0

- Molecule 37 is a protein called 60S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
37	AL35	122	1013	640	204	168	1	0	0

- Molecule 38 is a protein called 60S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
38	AL36	102	830	520	176	129	5	0	0

- Molecule 39 is a protein called 60S ribosomal protein L37.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
39	AL37	86	705	434	155	111	5	0	0

- Molecule 40 is a protein called Large ribosomal subunit protein eL38.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
40	AL38	69	569	366	103	99	1	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AL38	24	LYS	ASN	conflict	UNP G1U001

- Molecule 41 is a protein called 60S ribosomal protein L39.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
41	AL39	50	447	286	96	64	1	0	0

- Molecule 42 is a protein called Large ribosomal subunit protein eL40.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
42	AL40	52	429	266	90	67	6	0	0

- Molecule 43 is a protein called eL41.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
43	AL41	25	239	145	64	27	3	0	0

- Molecule 44 is a protein called eL42.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
44	AL42	104	851	533	174	138	6	0	0

- Molecule 45 is a protein called 60S ribosomal protein L37a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
45	AL43	91	708	445	136	120	7	0	0

- Molecule 46 is a protein called 60S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
46	ALNW	124	994	616	205	167	6	0	0

- Molecule 47 is a protein called Angiogenin.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
47	ANG	120	955	591	186	171	7	0	0

- Molecule 48 is a protein called Receptor of activated protein C kinase 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
48	ARAC	313	2436	1535	424	465	12	0	0

- Molecule 49 is a protein called 40S_SA_C domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
49	AS00	217	1710	1086	300	316	8	0	0

- Molecule 50 is a protein called 40S ribosomal protein S3a.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
50	AS01	213	1729	1098	309	308	14	0	0

- Molecule 51 is a protein called 40S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
51	AS02	221	1716	1111	295	301	9	0	0

- Molecule 52 is a protein called 40S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
52	AS03	228	1768	1126	318	316	8	0	0

- Molecule 53 is a protein called 40S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
53	AS04	262	2076	1324	386	358	8	0	0

- Molecule 54 is a protein called Ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
54	AS05	185	1471	921	277	266	7	0	0

- Molecule 55 is a protein called 40S ribosomal protein S6.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
55	AS06	237	1922	1199	387	329	7	0	0

- Molecule 56 is a protein called Small ribosomal subunit protein eS7.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
56	AS07	185	1488	952	271	264	1	0	0

- Molecule 57 is a protein called 40S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
57	AS08	206	1686	1058	332	291	5	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AS08	47	ARG	GLY	conflict	UNP G1TJW1

- Molecule 58 is a protein called 40S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
58	AS09	185	1525	969	306	248	2	0	0

- Molecule 59 is a protein called 40S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
59	AS10	96	810	530	143	131	6	0	0

- Molecule 60 is a protein called 40S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
60	AS11	143	1175	749	222	198	6	0	0

- Molecule 61 is a protein called 40S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
61	AS12	99	758	475	139	137	7	0	0

- Molecule 62 is a protein called 40S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
62	AS13	149	1202	770	228	203	1	0	0

- Molecule 63 is a protein called Small ribosomal subunit protein uS11.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
63	AS14	135	1004	614	196	188	6	0	0

- Molecule 64 is a protein called 40S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
64	AS15	120	997	635	187	168	7	0	0

- Molecule 65 is a protein called Small ribosomal subunit protein uS9.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
65	AS16	142	1128	717	213	195	3	0	0

- Molecule 66 is a protein called 40S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
66	AS17	132	1068	670	199	195	4	0	0

- Molecule 67 is a protein called 40S ribosomal protein S18.

Mol	Chain	Residues	Atoms					AltConf	Trace
67	AS18	144	Total	C	N	O	S	0	0
			1190	746	241	202	1		

- Molecule 68 is a protein called Small ribosomal subunit protein eS19.

Mol	Chain	Residues	Atoms					AltConf	Trace
68	AS19	141	Total	C	N	O	S	0	0
			1097	688	211	195	3		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AS19	119	GLY	TRP	conflict	UNP G1TN62

- Molecule 69 is a protein called 40S ribosomal protein S20.

Mol	Chain	Residues	Atoms					AltConf	Trace
69	AS20	100	Total	C	N	O	S	0	0
			795	498	152	141	4		

- Molecule 70 is a protein called eS21.

Mol	Chain	Residues	Atoms					AltConf	Trace
70	AS21	83	Total	C	N	O	S	0	0
			636	393	117	121	5		

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AS21	3	ASN	SER	conflict	UNP G1TM82
AS21	4	ASP	ASN	conflict	UNP G1TM82
AS21	33	GLN	PRO	conflict	UNP G1TM82
AS21	50	PHE	SER	conflict	UNP G1TM82
AS21	75	ALA	SER	conflict	UNP G1TM82
AS21	76	ASP	HIS	conflict	UNP G1TM82
AS21	81	LYS	GLN	conflict	UNP G1TM82

- Molecule 71 is a protein called 40S ribosomal protein S15a.

Mol	Chain	Residues	Atoms					AltConf	Trace
71	AS22	129	Total	C	N	O	S	0	0
			1034	659	193	176	6		

- Molecule 72 is a protein called 40S ribosomal protein S23.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
72	AS23	141	1098	693	219	183	3	0	0

- Molecule 73 is a protein called 40S ribosomal protein S24.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
73	AS24	124	1011	640	198	168	5	0	0

- Molecule 74 is a protein called 40S ribosomal protein S25.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
74	AS25	75	598	382	111	104	1	0	0

- Molecule 75 is a protein called eS26.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
75	AS26	101	814	507	170	132	5	0	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
AS26	28	ARG	CYS	conflict	UNP G1TFE8
AS26	56	ALA	VAL	conflict	UNP G1TFE8

- Molecule 76 is a protein called 40S ribosomal protein S27.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
76	AS27	83	651	408	121	115	7	0	0

- Molecule 77 is a protein called 40S ribosomal protein S28.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
77	AS28	62	488	297	97	92	2	0	0

- Molecule 78 is a protein called 40S ribosomal protein S29.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
78	AS29	55	459	286	94	74	5	0	0

- Molecule 79 is a protein called 40S ribosomal protein S30.

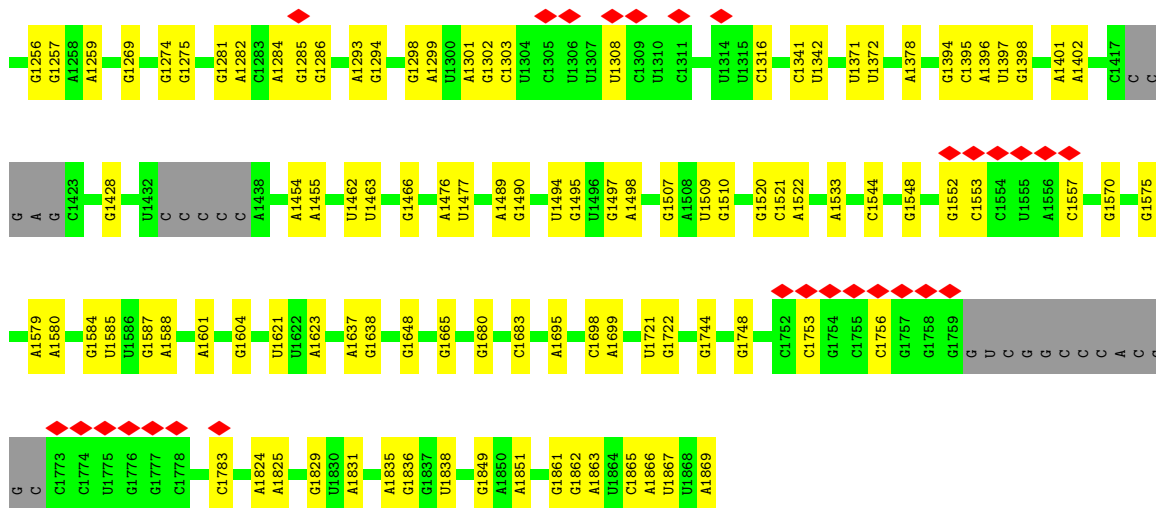
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
79	AS30	55	443	274	97	71	1	0	0

- Molecule 80 is a RNA chain called tRNAfMet.

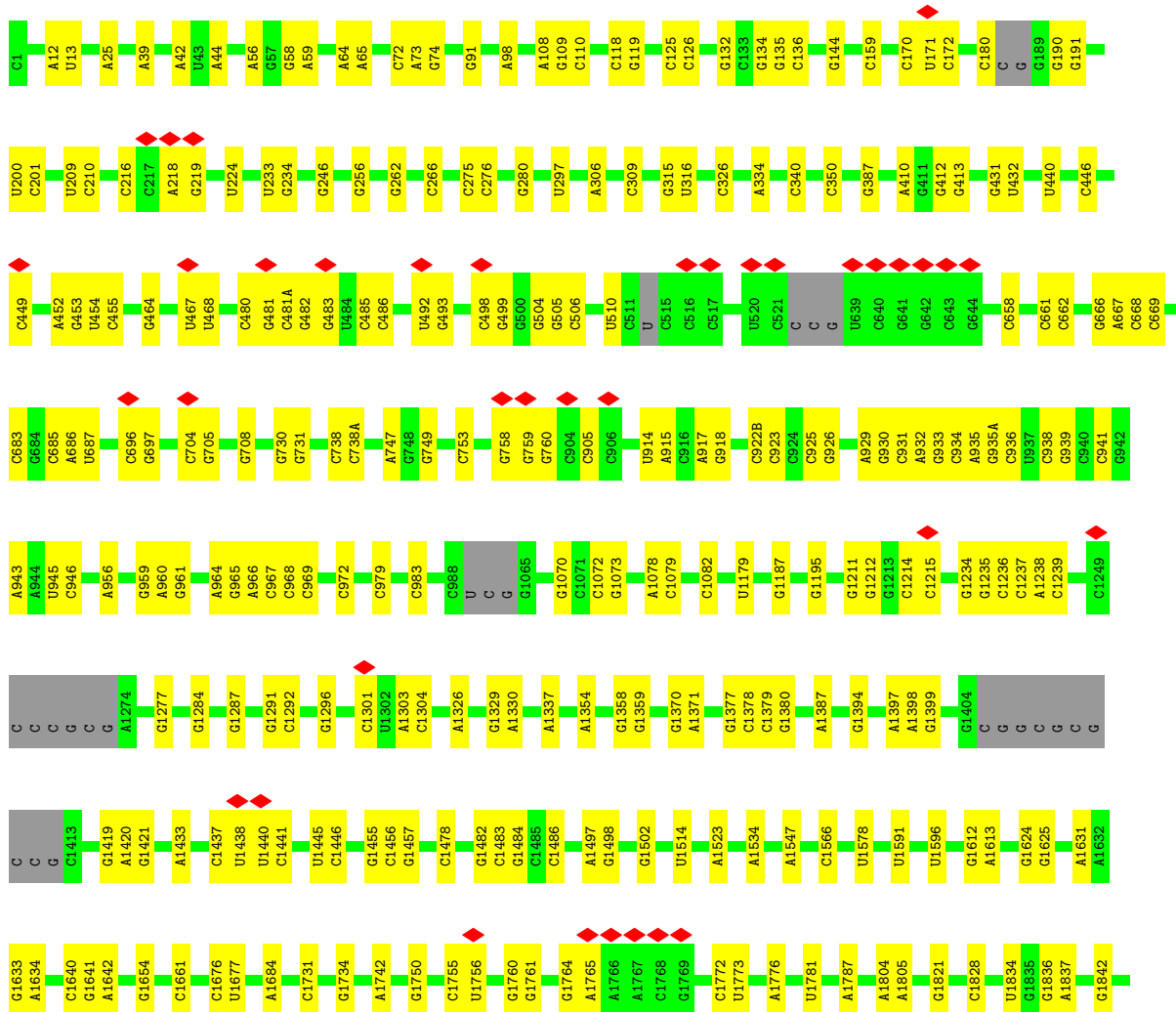
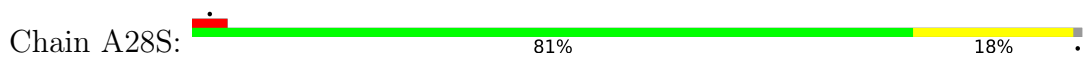
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
80	ETRN	77	1640	732	297	535	76	0	0
80	PTRN	77	1640	732	297	535	76	0	0

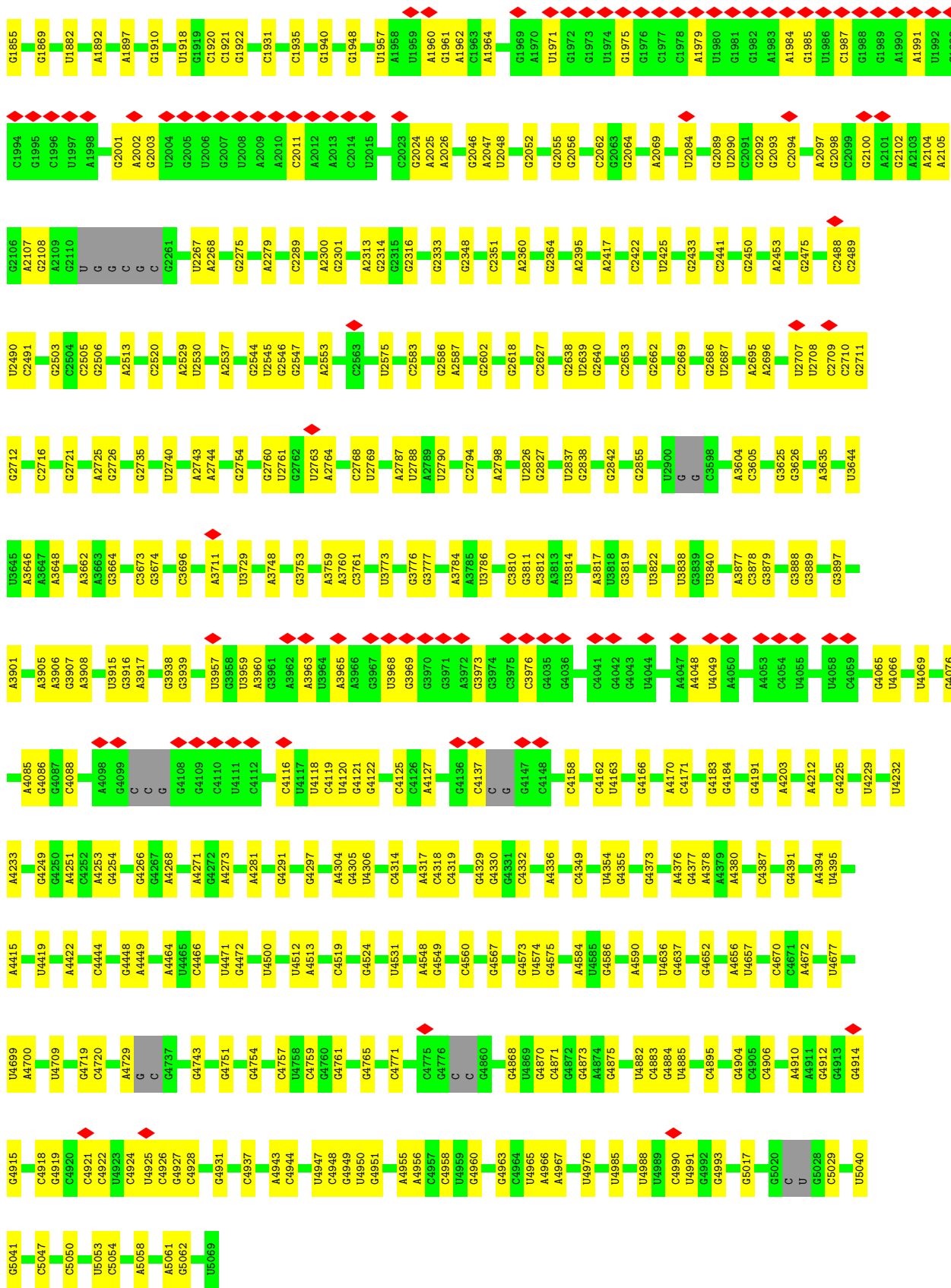
- Molecule 81 is a RNA chain called Model mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
81	MRNA	10	207	93	33	71	10	0	0

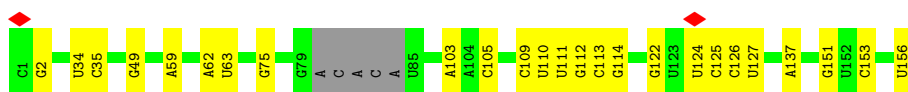
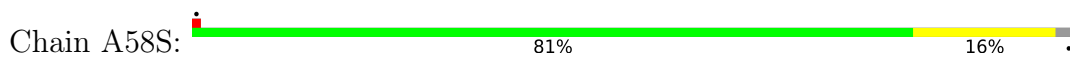


• Molecule 2: 28S rRNA





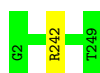
• Molecule 3: 5.8S rRNA



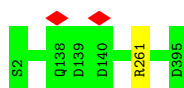
• Molecule 4: 5S rRNA



• Molecule 5: 60S ribosomal protein L8



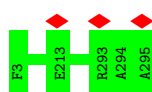
• Molecule 6: 60S ribosomal protein L3



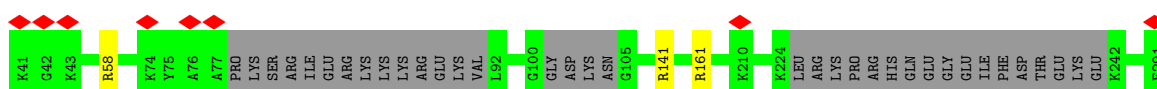
• Molecule 7: 60S ribosomal protein L4



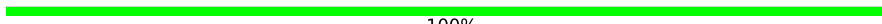
• Molecule 8: Large ribosomal subunit protein uL18



• Molecule 9: 60S ribosomal protein L6



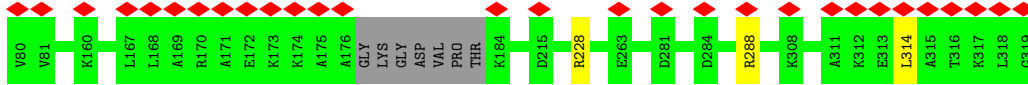
- Molecule 10: 60S ribosomal protein L7

Chain AL07:  100%



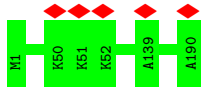
- Molecule 11: Large ribosomal subunit protein eL8

Chain AL08:  12%  96%



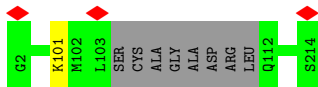
- Molecule 12: 60S ribosomal protein L9

Chain AL09:  100%



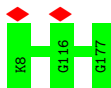
- Molecule 13: Ribosomal protein L10

Chain AL10:  96%



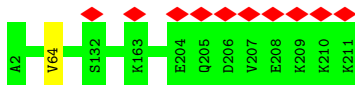
- Molecule 14: 60S ribosomal protein L11

Chain AL11:  100%



- Molecule 15: Large ribosomal subunit protein eL13

Chain AL13:  5%  100%



- Molecule 16: Large ribosomal subunit protein eL14

Chain AL14:  99%



- Molecule 17: 60S ribosomal protein L15

Chain AL15: 99%



- Molecule 18: Large ribosomal subunit protein uL13

Chain AL16: 98%



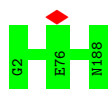
- Molecule 19: 60S ribosomal protein L17

Chain AL17: 100%

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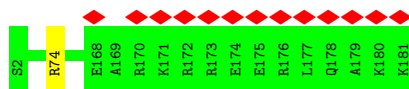
- Molecule 20: Large ribosomal subunit protein eL18

Chain AL18: 100%



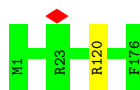
- Molecule 21: 60S ribosomal protein L19

Chain AL19: 99%



- Molecule 22: 60S ribosomal protein L18a

Chain AL20: 99%

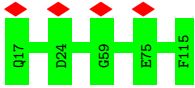


- Molecule 23: 60S ribosomal protein L21

Chain AL21: 100%



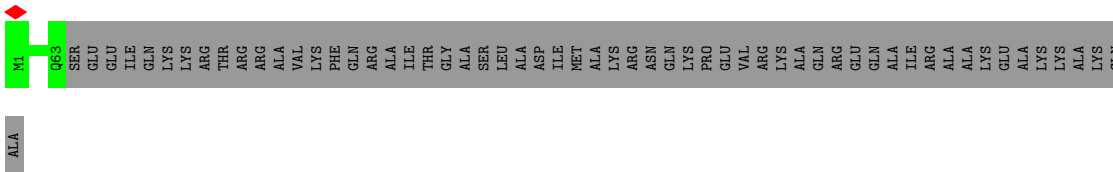
- Molecule 24: Large ribosomal subunit protein eL22



- Molecule 25: 60S ribosomal protein L23



- Molecule 26: Ribosomal protein L24



- Molecule 27: Large ribosomal subunit protein uL23



- Molecule 28: 60S ribosomal protein L26

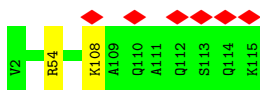


- Molecule 29: 60S ribosomal protein L27

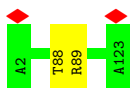


There are no outlier residues recorded for this chain.

- Molecule 30: 60S ribosomal protein L27a



- Molecule 37: 60S ribosomal protein L35



- Molecule 38: 60S ribosomal protein L36



- Molecule 39: 60S ribosomal protein L37



There are no outlier residues recorded for this chain.

- Molecule 40: Large ribosomal subunit protein eL38



- Molecule 41: 60S ribosomal protein L39



- Molecule 42: Large ribosomal subunit protein eL40



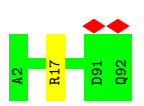
- Molecule 43: eL41



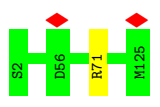
• Molecule 44: eL42



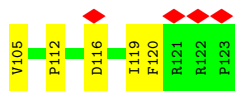
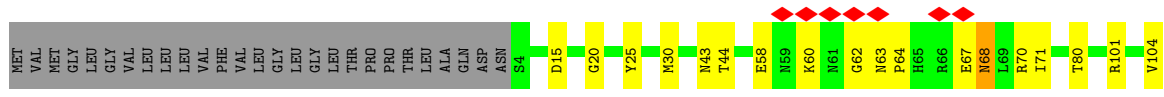
• Molecule 45: 60S ribosomal protein L37a



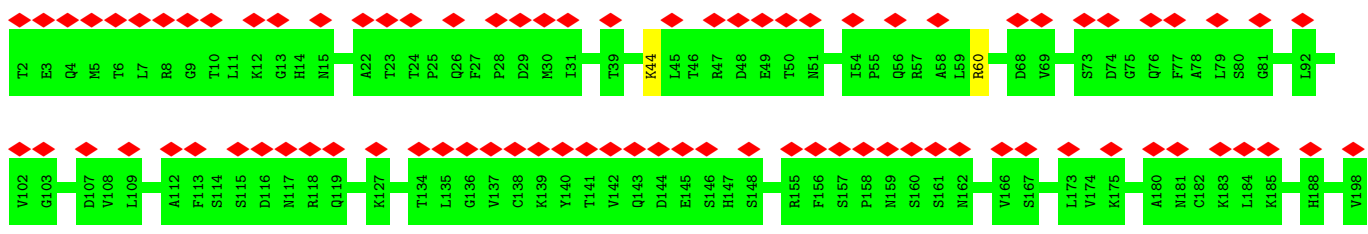
• Molecule 46: 60S ribosomal protein L28

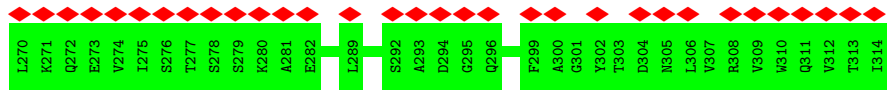
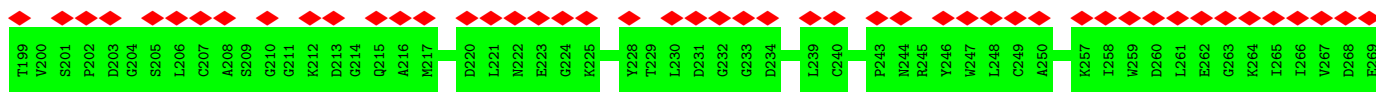


• Molecule 47: Angiogenin

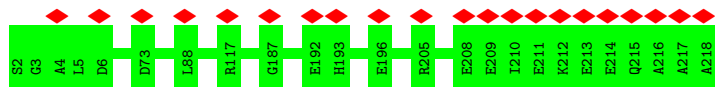


• Molecule 48: Receptor of activated protein C kinase 1

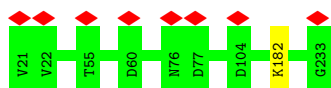




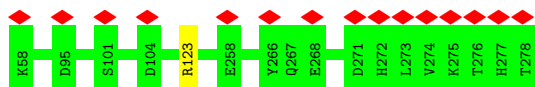
• Molecule 49: 40S_SA_C domain-containing protein



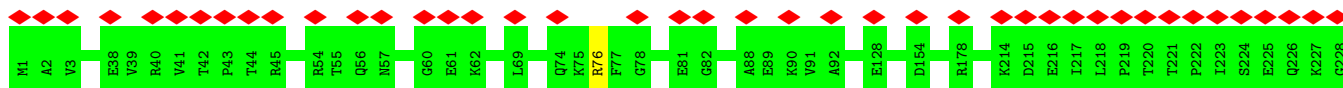
• Molecule 50: 40S ribosomal protein S3a



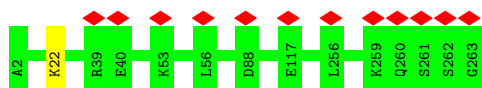
• Molecule 51: 40S ribosomal protein S2



• Molecule 52: 40S ribosomal protein S3

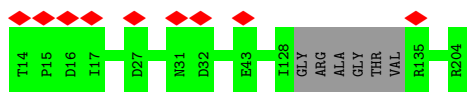


• Molecule 53: 40S ribosomal protein S4

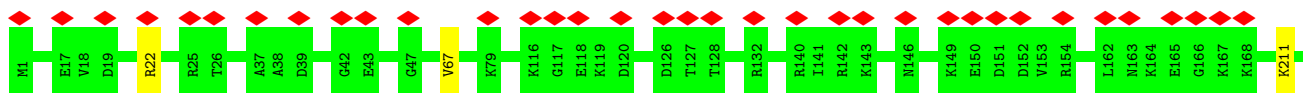


• Molecule 54: Ribosomal protein S5

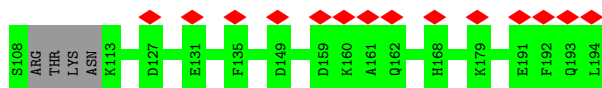
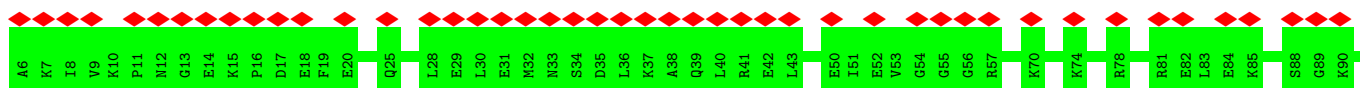




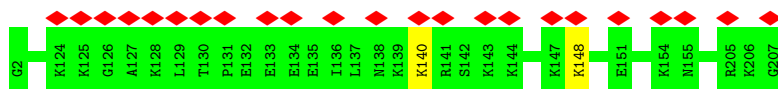
- Molecule 55: 40S ribosomal protein S6



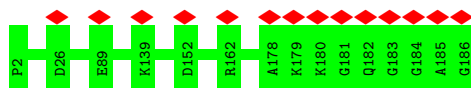
- Molecule 56: Small ribosomal subunit protein eS7



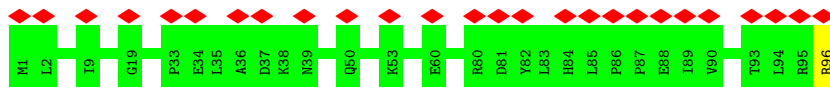
- Molecule 57: 40S ribosomal protein S8



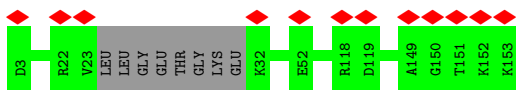
- Molecule 58: 40S ribosomal protein S9



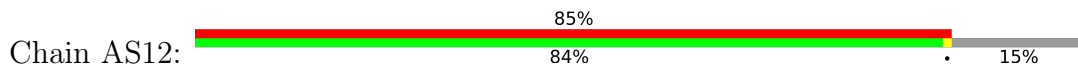
- Molecule 59: 40S ribosomal protein S10



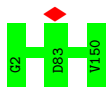
- Molecule 60: 40S ribosomal protein S11



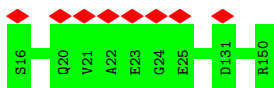
- Molecule 61: 40S ribosomal protein S12



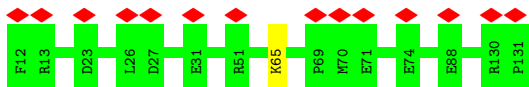
- Molecule 62: 40S ribosomal protein S13



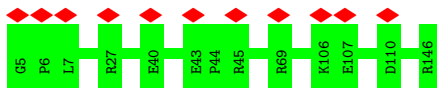
- Molecule 63: Small ribosomal subunit protein uS11



- Molecule 64: 40S ribosomal protein S15

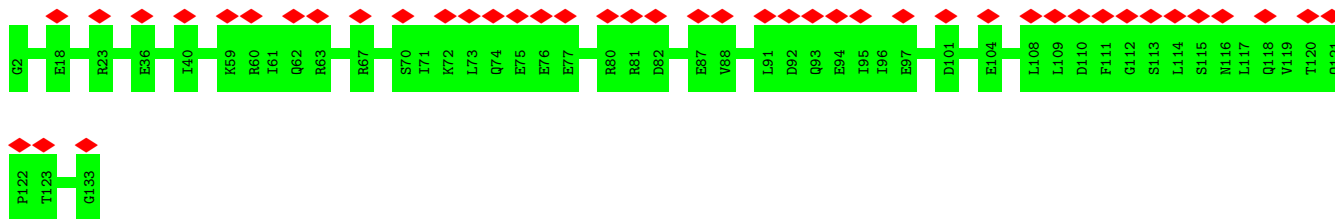


- Molecule 65: Small ribosomal subunit protein uS9

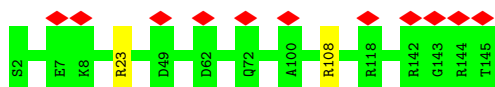


- Molecule 66: 40S ribosomal protein S17

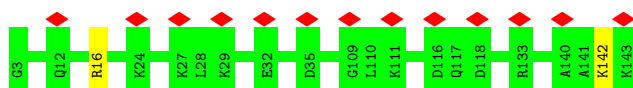




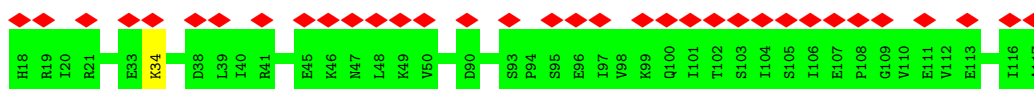
- Molecule 67: 40S ribosomal protein S18



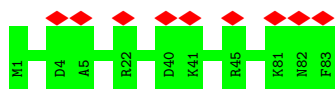
- Molecule 68: Small ribosomal subunit protein eS19



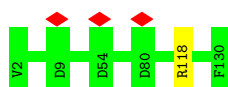
- Molecule 69: 40S ribosomal protein S20



- Molecule 70: eS21



- Molecule 71: 40S ribosomal protein S15a

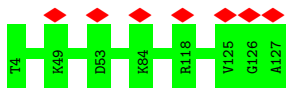


- Molecule 72: 40S ribosomal protein S23

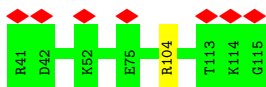




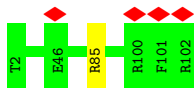
- Molecule 73: 40S ribosomal protein S24



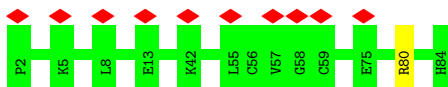
- Molecule 74: 40S ribosomal protein S25



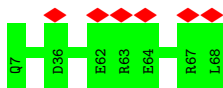
- Molecule 75: eS26



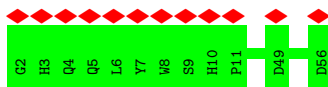
- Molecule 76: 40S ribosomal protein S27



- Molecule 77: 40S ribosomal protein S28



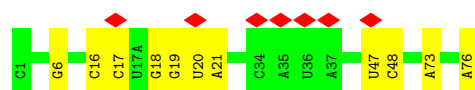
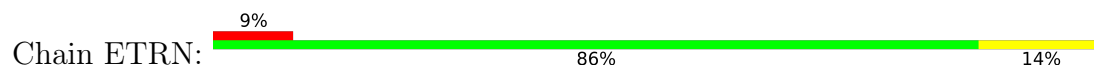
- Molecule 78: 40S ribosomal protein S29



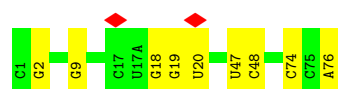
- Molecule 79: 40S ribosomal protein S30



- Molecule 80: tRNA^{fMet}



- Molecule 80: tRNA^{fMet}



- Molecule 81: Model mRNA



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	101572	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	FEI TALOS ARCTICA	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	30	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	30.079	Depositor
Minimum map value	-10.342	Depositor
Average map value	0.006	Depositor
Map value standard deviation	0.932	Depositor
Recommended contour level	3.5	Depositor
Map size (\AA)	528.96, 528.96, 528.96	wwPDB
Map dimensions	608, 608, 608	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.87000006, 0.87000006, 0.87000006	Depositor

5 Model quality

5.1 Standard geometry

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	A18S	0.28	0/40369	0.25	0/62910
2	A28S	0.39	0/85336	0.29	0/133087
3	A58S	0.39	0/3581	0.24	0/5577
4	A5S	0.34	0/2858	0.21	0/4455
5	AL02	0.42	0/1936	0.49	0/2596
6	AL03	0.37	0/3240	0.44	0/4339
7	AL04	0.38	0/2937	0.46	0/3946
8	AL05	0.30	0/2437	0.40	0/3264
9	AL06	0.28	0/1762	0.44	0/2362
10	AL07	0.38	0/1911	0.45	0/2549
11	AL08	0.31	0/1910	0.46	0/2569
12	AL09	0.33	0/1535	0.48	0/2063
13	AL10	0.35	0/1702	0.43	0/2272
14	AL11	0.28	0/1384	0.44	0/1850
15	AL13	0.34	0/1733	0.47	0/2316
16	AL14	0.31	0/1158	0.46	0/1547
17	AL15	0.44	0/1746	0.51	0/2338
18	AL16	0.38	0/1662	0.46	1/2222 (0.0%)
19	AL17	0.40	0/1268	0.47	0/1700
20	AL18	0.38	0/1539	0.49	0/2054
21	AL19	0.34	0/1524	0.45	0/2013
22	AL20	0.37	0/1501	0.45	0/2012
23	AL21	0.36	0/1326	0.46	0/1770
24	AL22	0.27	0/823	0.49	0/1104
25	AL23	0.38	0/993	0.43	0/1332
26	AL24	0.35	0/541	0.46	0/720
27	AL25	0.34	0/984	0.40	0/1323
28	AL26	0.35	0/1132	0.45	0/1504
29	AL27	0.33	0/1130	0.42	0/1507
30	AL28	0.40	0/1191	0.44	0/1590
31	AL29	0.29	0/861	0.43	0/1138
32	AL30	0.34	0/771	0.36	0/1034
33	AL31	0.36	0/903	0.43	0/1216
34	AL32	0.42	0/1071	0.52	0/1429

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
35	AL33	0.41	0/895	0.50	0/1198
36	AL34	0.36	0/916	0.47	0/1220
37	AL35	0.32	0/1021	0.42	0/1348
38	AL36	0.31	0/841	0.47	0/1112
39	AL37	0.44	0/720	0.50	0/952
40	AL38	0.28	0/575	0.47	0/761
41	AL39	0.37	0/459	0.47	0/608
42	AL40	0.32	0/435	0.49	0/575
43	AL41	0.32	0/240	0.49	0/305
44	AL42	0.36	0/864	0.45	0/1140
45	AL43	0.39	0/718	0.40	0/953
46	ALNW	0.35	0/1010	0.44	0/1354
47	ANG	0.30	0/979	0.52	0/1322
48	ARAC	0.23	1/2493 (0.0%)	0.47	0/3394
49	AS00	0.24	0/1747	0.45	0/2374
50	AS01	0.26	0/1756	0.43	0/2350
51	AS02	0.26	0/1753	0.44	0/2369
52	AS03	0.23	0/1796	0.50	0/2417
53	AS04	0.24	0/2118	0.43	0/2849
54	AS05	0.25	0/1492	0.49	0/2005
55	AS06	0.21	0/1944	0.45	1/2586 (0.0%)
56	AS07	0.22	0/1510	0.50	0/2022
57	AS08	0.27	0/1715	0.46	0/2287
58	AS09	0.23	0/1550	0.44	0/2069
59	AS10	0.24	0/834	0.55	0/1125
60	AS11	0.31	0/1195	0.44	0/1597
61	AS12	0.22	0/765	0.53	0/1029
62	AS13	0.28	0/1226	0.43	0/1649
63	AS14	0.30	0/1017	0.45	0/1365
64	AS15	0.23	0/1017	0.48	0/1358
65	AS16	0.24	0/1146	0.45	0/1534
66	AS17	0.23	0/1082	0.53	0/1452
67	AS18	0.25	0/1208	0.52	0/1618
68	AS19	0.24	0/1115	0.45	0/1493
69	AS20	0.23	0/805	0.44	0/1081
70	AS21	0.23	0/643	0.45	0/860
71	AS22	0.28	0/1051	0.40	0/1406
72	AS23	0.30	0/1116	0.49	0/1490
73	AS24	0.23	0/1028	0.40	0/1366
74	AS25	0.24	0/604	0.44	0/810
75	AS26	0.31	0/828	0.46	0/1109
76	AS27	0.23	0/665	0.40	0/891
77	AS28	0.25	0/490	0.42	0/656

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
78	AS29	0.23	0/470	0.50	0/623
79	AS30	0.22	0/447	0.42	0/587
80	ETRN	0.18	0/1832	0.23	0/2855
80	PTRN	0.23	0/1832	0.24	0/2855
81	MRNA	0.25	0/229	0.39	0/353
All	All	0.34	1/228947 (0.0%)	0.35	2/336440 (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
33	AL31	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
48	ARAC	44	LYS	CA-CB	5.13	1.62	1.53

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
55	AS06	67	VAL	N-CA-C	-5.81	108.19	113.71
18	AL16	185	VAL	N-CA-C	-5.32	108.10	113.20

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
33	AL31	95	ASP	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	A18S	36103	0	0	0	0
2	A28S	76286	0	0	0	0
3	A58S	3208	0	0	0	0
4	A5S	2558	0	1296	3	0
5	AL02	1898	0	0	0	0
6	AL03	3172	0	0	0	0
7	AL04	2883	0	0	0	0
8	AL05	2391	0	0	0	0
9	AL06	1729	0	0	0	0
10	AL07	1875	0	0	0	0
11	AL08	1879	0	0	0	0
12	AL09	1516	0	0	0	0
13	AL10	1664	0	0	0	0
14	AL11	1361	0	0	0	0
15	AL13	1702	0	0	0	0
16	AL14	1137	0	0	0	0
17	AL15	1701	0	0	0	0
18	AL16	1630	0	0	0	0
19	AL17	1242	0	0	0	0
20	AL18	1515	0	0	0	0
21	AL19	1508	0	0	0	0
22	AL20	1462	0	0	0	0
23	AL21	1298	0	0	0	0
24	AL22	809	0	0	0	0
25	AL23	979	0	0	0	0
26	AL24	528	0	0	0	0
27	AL25	967	0	0	0	0
28	AL26	1115	0	0	0	0
29	AL27	1107	0	0	0	0
30	AL28	1162	0	0	0	0
31	AL29	848	0	0	0	0
32	AL30	761	0	0	0	0
33	AL31	888	0	0	0	0
34	AL32	1053	0	0	0	0
35	AL33	876	0	0	0	0
36	AL34	906	0	0	0	0
37	AL35	1013	0	0	0	0
38	AL36	830	0	0	0	0
39	AL37	705	0	0	0	0
40	AL38	569	0	0	0	0
41	AL39	447	0	0	0	0
42	AL40	429	0	0	0	0
43	AL41	239	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
44	AL42	851	0	0	0	0
45	AL43	708	0	0	0	0
46	ALNW	994	0	0	0	0
47	ANG	955	0	922	12	0
48	ARAC	2436	0	0	0	0
49	AS00	1710	0	0	0	0
50	AS01	1729	0	0	0	0
51	AS02	1716	0	0	0	0
52	AS03	1768	0	0	0	0
53	AS04	2076	0	0	0	0
54	AS05	1471	0	0	0	0
55	AS06	1922	0	0	0	0
56	AS07	1488	0	0	0	0
57	AS08	1686	0	0	0	0
58	AS09	1525	0	0	0	0
59	AS10	810	0	0	0	0
60	AS11	1175	0	0	0	0
61	AS12	758	0	0	0	0
62	AS13	1202	0	0	0	0
63	AS14	1004	0	0	0	0
64	AS15	997	0	0	0	0
65	AS16	1128	0	0	0	0
66	AS17	1068	0	0	0	0
67	AS18	1190	0	0	0	0
68	AS19	1097	0	0	0	0
69	AS20	795	0	0	0	0
70	AS21	636	0	0	0	0
71	AS22	1034	0	0	0	0
72	AS23	1098	0	0	0	0
73	AS24	1011	0	0	0	0
74	AS25	598	0	0	0	0
75	AS26	814	0	0	0	0
76	AS27	651	0	0	0	0
77	AS28	488	0	0	0	0
78	AS29	459	0	0	0	0
79	AS30	443	0	0	0	0
80	ETRN	1640	0	0	0	0
80	PTRN	1640	0	0	0	0
81	MRNA	207	0	0	0	0
All	All	212927	0	2218	15	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 7.

The worst 5 of 15 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)
47:ANG:20:GLY:HA3	47:ANG:25:TYR:HB2	1.65	0.79
47:ANG:30:MET:HE1	47:ANG:43:ASN:HB3	1.81	0.62
47:ANG:105:VAL:HG11	47:ANG:112:PRO:HB3	1.86	0.58
47:ANG:71:ILE:HG22	47:ANG:104:VAL:HB	1.88	0.55
47:ANG:104:VAL:HG13	47:ANG:116:ASP:HB3	1.93	0.51

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	
5	AL02	246/248 (99%)	233 (95%)	13 (5%)	0	100	100
6	AL03	392/394 (100%)	378 (96%)	14 (4%)	0	100	100
7	AL04	360/362 (99%)	346 (96%)	14 (4%)	0	100	100
8	AL05	291/293 (99%)	279 (96%)	12 (4%)	0	100	100
9	AL06	208/251 (83%)	198 (95%)	10 (5%)	0	100	100
10	AL07	223/225 (99%)	218 (98%)	4 (2%)	1 (0%)	30	60
11	AL08	229/240 (95%)	219 (96%)	10 (4%)	0	100	100
12	AL09	188/190 (99%)	177 (94%)	11 (6%)	0	100	100
13	AL10	201/213 (94%)	191 (95%)	10 (5%)	0	100	100
14	AL11	168/170 (99%)	161 (96%)	7 (4%)	0	100	100
15	AL13	208/210 (99%)	197 (95%)	10 (5%)	1 (0%)	24	55
16	AL14	136/138 (99%)	128 (94%)	8 (6%)	0	100	100
17	AL15	201/203 (99%)	188 (94%)	12 (6%)	1 (0%)	24	55
18	AL16	197/199 (99%)	195 (99%)	2 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
19	AL17	151/153 (99%)	148 (98%)	3 (2%)	0	100	100
20	AL18	185/187 (99%)	172 (93%)	13 (7%)	0	100	100
21	AL19	178/180 (99%)	174 (98%)	4 (2%)	0	100	100
22	AL20	174/176 (99%)	167 (96%)	7 (4%)	0	100	100
23	AL21	157/159 (99%)	151 (96%)	6 (4%)	0	100	100
24	AL22	97/99 (98%)	91 (94%)	6 (6%)	0	100	100
25	AL23	129/131 (98%)	126 (98%)	3 (2%)	0	100	100
26	AL24	61/121 (50%)	59 (97%)	2 (3%)	0	100	100
27	AL25	116/118 (98%)	113 (97%)	3 (3%)	0	100	100
28	AL26	132/134 (98%)	130 (98%)	2 (2%)	0	100	100
29	AL27	133/135 (98%)	126 (95%)	7 (5%)	0	100	100
30	AL28	145/147 (99%)	137 (94%)	8 (6%)	0	100	100
31	AL29	100/245 (41%)	94 (94%)	6 (6%)	0	100	100
32	AL30	96/98 (98%)	93 (97%)	3 (3%)	0	100	100
33	AL31	105/107 (98%)	97 (92%)	7 (7%)	1 (1%)	12	38
34	AL32	126/128 (98%)	122 (97%)	4 (3%)	0	100	100
35	AL33	107/109 (98%)	101 (94%)	6 (6%)	0	100	100
36	AL34	112/114 (98%)	109 (97%)	3 (3%)	0	100	100
37	AL35	120/122 (98%)	116 (97%)	2 (2%)	2 (2%)	7	25
38	AL36	100/102 (98%)	95 (95%)	5 (5%)	0	100	100
39	AL37	84/86 (98%)	83 (99%)	1 (1%)	0	100	100
40	AL38	67/69 (97%)	65 (97%)	2 (3%)	0	100	100
41	AL39	48/50 (96%)	44 (92%)	4 (8%)	0	100	100
42	AL40	50/52 (96%)	50 (100%)	0	0	100	100
43	AL41	23/25 (92%)	23 (100%)	0	0	100	100
44	AL42	102/104 (98%)	99 (97%)	3 (3%)	0	100	100
45	AL43	89/91 (98%)	84 (94%)	5 (6%)	0	100	100
46	ALNW	122/124 (98%)	119 (98%)	3 (2%)	0	100	100
47	ANG	118/147 (80%)	113 (96%)	4 (3%)	1 (1%)	16	44
48	ARAC	311/313 (99%)	293 (94%)	18 (6%)	0	100	100
49	AS00	215/217 (99%)	201 (94%)	14 (6%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
50	AS01	211/213 (99%)	200 (95%)	11 (5%)	0	100	100
51	AS02	219/221 (99%)	214 (98%)	5 (2%)	0	100	100
52	AS03	226/228 (99%)	220 (97%)	6 (3%)	0	100	100
53	AS04	260/262 (99%)	253 (97%)	7 (3%)	0	100	100
54	AS05	181/191 (95%)	165 (91%)	16 (9%)	0	100	100
55	AS06	235/237 (99%)	227 (97%)	8 (3%)	0	100	100
56	AS07	181/189 (96%)	175 (97%)	6 (3%)	0	100	100
57	AS08	204/206 (99%)	193 (95%)	11 (5%)	0	100	100
58	AS09	183/185 (99%)	180 (98%)	3 (2%)	0	100	100
59	AS10	94/96 (98%)	85 (90%)	9 (10%)	0	100	100
60	AS11	139/151 (92%)	130 (94%)	9 (6%)	0	100	100
61	AS12	97/117 (83%)	93 (96%)	4 (4%)	0	100	100
62	AS13	147/149 (99%)	142 (97%)	5 (3%)	0	100	100
63	AS14	133/135 (98%)	127 (96%)	6 (4%)	0	100	100
64	AS15	118/120 (98%)	112 (95%)	6 (5%)	0	100	100
65	AS16	140/142 (99%)	135 (96%)	5 (4%)	0	100	100
66	AS17	130/132 (98%)	118 (91%)	12 (9%)	0	100	100
67	AS18	142/144 (99%)	131 (92%)	11 (8%)	0	100	100
68	AS19	139/141 (99%)	133 (96%)	6 (4%)	0	100	100
69	AS20	98/100 (98%)	93 (95%)	5 (5%)	0	100	100
70	AS21	81/83 (98%)	80 (99%)	1 (1%)	0	100	100
71	AS22	127/129 (98%)	122 (96%)	5 (4%)	0	100	100
72	AS23	139/141 (99%)	130 (94%)	9 (6%)	0	100	100
73	AS24	122/124 (98%)	121 (99%)	1 (1%)	0	100	100
74	AS25	73/75 (97%)	73 (100%)	0	0	100	100
75	AS26	99/101 (98%)	94 (95%)	5 (5%)	0	100	100
76	AS27	81/83 (98%)	78 (96%)	3 (4%)	0	100	100
77	AS28	60/62 (97%)	59 (98%)	1 (2%)	0	100	100
78	AS29	53/55 (96%)	51 (96%)	2 (4%)	0	100	100
79	AS30	53/55 (96%)	52 (98%)	1 (2%)	0	100	100
All	All	11166/11646 (96%)	10689 (96%)	470 (4%)	7 (0%)	49	77

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
10	AL07	236	GLU
15	AL13	64	VAL
17	AL15	79	ALA
37	AL35	89	ARG
47	ANG	68	ASN

5.3.2 Protein sidechains [i](#)

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
5	AL02	190/190 (100%)	189 (100%)	1 (0%)	81 93
6	AL03	342/342 (100%)	341 (100%)	1 (0%)	86 95
7	AL04	302/302 (100%)	299 (99%)	3 (1%)	68 88
8	AL05	247/247 (100%)	247 (100%)	0	100 100
9	AL06	190/223 (85%)	187 (98%)	3 (2%)	55 83
10	AL07	196/196 (100%)	196 (100%)	0	100 100
11	AL08	200/205 (98%)	197 (98%)	3 (2%)	57 84
12	AL09	169/169 (100%)	169 (100%)	0	100 100
13	AL10	175/180 (97%)	174 (99%)	1 (1%)	78 92
14	AL11	143/143 (100%)	143 (100%)	0	100 100
15	AL13	175/175 (100%)	175 (100%)	0	100 100
16	AL14	117/117 (100%)	116 (99%)	1 (1%)	70 89
17	AL15	171/171 (100%)	170 (99%)	1 (1%)	78 92
18	AL16	171/171 (100%)	169 (99%)	2 (1%)	63 87
19	AL17	134/134 (100%)	134 (100%)	0	100 100
20	AL18	164/164 (100%)	164 (100%)	0	100 100
21	AL19	159/159 (100%)	158 (99%)	1 (1%)	78 92
22	AL20	157/157 (100%)	156 (99%)	1 (1%)	78 92
23	AL21	139/139 (100%)	139 (100%)	0	100 100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
24	AL22	89/89 (100%)	89 (100%)	0	100	100
25	AL23	101/101 (100%)	99 (98%)	2 (2%)	48	80
26	AL24	55/100 (55%)	55 (100%)	0	100	100
27	AL25	106/106 (100%)	106 (100%)	0	100	100
28	AL26	124/124 (100%)	122 (98%)	2 (2%)	55	83
29	AL27	117/117 (100%)	117 (100%)	0	100	100
30	AL28	119/119 (100%)	119 (100%)	0	100	100
31	AL29	84/184 (46%)	83 (99%)	1 (1%)	63	87
32	AL30	84/84 (100%)	84 (100%)	0	100	100
33	AL31	98/98 (100%)	98 (100%)	0	100	100
34	AL32	114/114 (100%)	114 (100%)	0	100	100
35	AL33	88/88 (100%)	88 (100%)	0	100	100
36	AL34	98/98 (100%)	96 (98%)	2 (2%)	48	80
37	AL35	109/109 (100%)	109 (100%)	0	100	100
38	AL36	86/86 (100%)	85 (99%)	1 (1%)	63	87
39	AL37	73/73 (100%)	73 (100%)	0	100	100
40	AL38	64/64 (100%)	64 (100%)	0	100	100
41	AL39	47/47 (100%)	46 (98%)	1 (2%)	47	79
42	AL40	48/48 (100%)	48 (100%)	0	100	100
43	AL41	24/24 (100%)	24 (100%)	0	100	100
44	AL42	92/92 (100%)	90 (98%)	2 (2%)	45	78
45	AL43	74/74 (100%)	73 (99%)	1 (1%)	59	85
46	ALNW	108/108 (100%)	107 (99%)	1 (1%)	70	89
47	ANG	105/129 (81%)	103 (98%)	2 (2%)	50	81
48	ARAC	272/272 (100%)	271 (100%)	1 (0%)	84	94
49	AS00	180/181 (99%)	180 (100%)	0	100	100
50	AS01	194/194 (100%)	193 (100%)	1 (0%)	81	93
51	AS02	187/187 (100%)	186 (100%)	1 (0%)	81	93
52	AS03	190/190 (100%)	189 (100%)	1 (0%)	81	93
53	AS04	224/224 (100%)	223 (100%)	1 (0%)	84	94
54	AS05	158/161 (98%)	158 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
55	AS06	206/207 (100%)	203 (98%)	3 (2%)	57	84
56	AS07	165/169 (98%)	165 (100%)	0	100	100
57	AS08	178/178 (100%)	176 (99%)	2 (1%)	65	87
58	AS09	161/161 (100%)	161 (100%)	0	100	100
59	AS10	87/87 (100%)	86 (99%)	1 (1%)	65	87
60	AS11	130/136 (96%)	130 (100%)	0	100	100
61	AS12	83/99 (84%)	82 (99%)	1 (1%)	63	87
62	AS13	130/130 (100%)	130 (100%)	0	100	100
63	AS14	104/105 (99%)	104 (100%)	0	100	100
64	AS15	109/109 (100%)	108 (99%)	1 (1%)	70	89
65	AS16	117/117 (100%)	117 (100%)	0	100	100
66	AS17	119/119 (100%)	119 (100%)	0	100	100
67	AS18	125/125 (100%)	123 (98%)	2 (2%)	55	83
68	AS19	111/111 (100%)	109 (98%)	2 (2%)	51	82
69	AS20	92/92 (100%)	91 (99%)	1 (1%)	65	87
70	AS21	67/67 (100%)	67 (100%)	0	100	100
71	AS22	112/112 (100%)	111 (99%)	1 (1%)	70	89
72	AS23	113/113 (100%)	113 (100%)	0	100	100
73	AS24	107/107 (100%)	107 (100%)	0	100	100
74	AS25	66/66 (100%)	65 (98%)	1 (2%)	57	84
75	AS26	88/88 (100%)	87 (99%)	1 (1%)	65	87
76	AS27	75/75 (100%)	74 (99%)	1 (1%)	61	86
77	AS28	55/55 (100%)	55 (100%)	0	100	100
78	AS29	48/48 (100%)	48 (100%)	0	100	100
79	AS30	46/46 (100%)	45 (98%)	1 (2%)	45	78
All	All	9747/9991 (98%)	9691 (99%)	56 (1%)	76	92

5 of 56 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
44	AL42	87	ARG
79	AS30	99	LYS
51	AS02	123	ARG

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Mol	Chain	Res	Type
76	AS27	80	ARG
68	AS19	142	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A18S	1680/1869 (89%)	293 (17%)	14 (0%)
2	A28S	3534/3601 (98%)	629 (17%)	44 (1%)
3	A58S	149/156 (95%)	24 (16%)	1 (0%)
4	A5S	119/120 (99%)	10 (8%)	0
80	ETR N	76/77 (98%)	11 (14%)	0
80	PTR N	76/77 (98%)	9 (11%)	1 (1%)
81	MRNA	9/28 (32%)	4 (44%)	1 (11%)
All	All	5643/5928 (95%)	980 (17%)	61 (1%)

5 of 980 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A18S	2	A
1	A18S	3	C
1	A18S	17	C
1	A18S	25	A
1	A18S	26	U

5 of 61 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
2	A28S	1238	A
2	A28S	4884	G
2	A28S	1633	G
2	A28S	4719	G
81	MRNA	6	U

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

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
2	A28S	12

The worst 5 of 12 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A28S	1219:G	O3'	1233:G	P	20.06
1	A28S	3976:C	O3'	4035:G	P	14.96
1	A28S	1696:C	O3'	1720:C	P	14.73
1	A28S	1364:U	O3'	1368:A	P	13.71
1	A28S	1180:C	O3'	1183:C	P	8.80

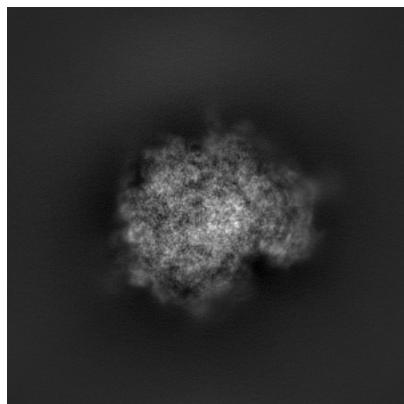
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-44461. These allow visual inspection of the internal detail of the map and identification of artifacts.

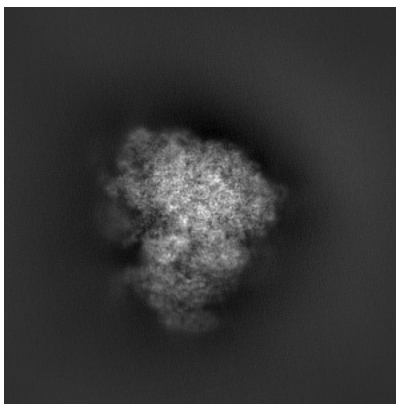
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

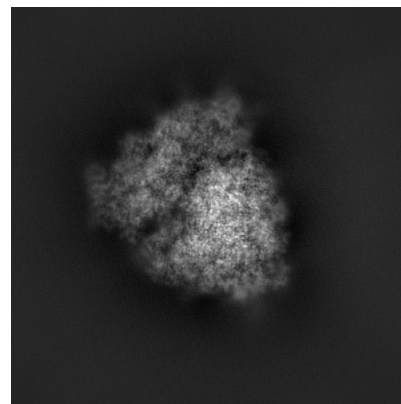
6.1.1 Primary map



X

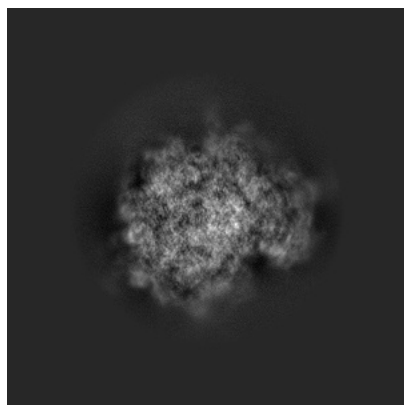


Y

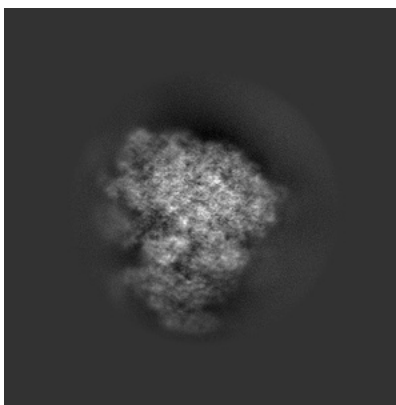


Z

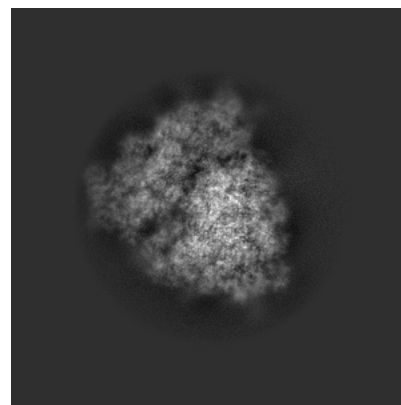
6.1.2 Raw map



X



Y

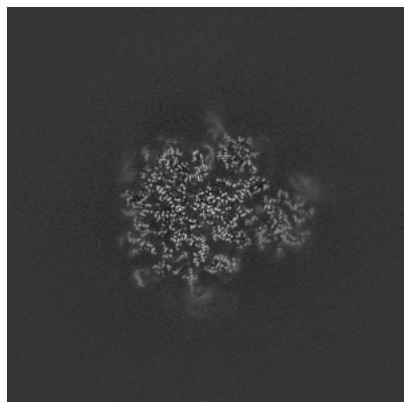


Z

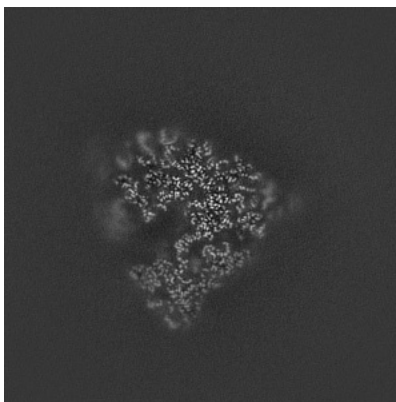
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

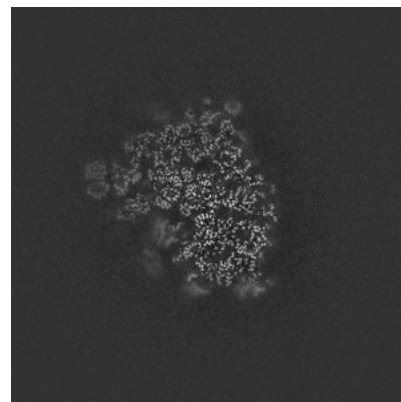
6.2.1 Primary map



X Index: 304

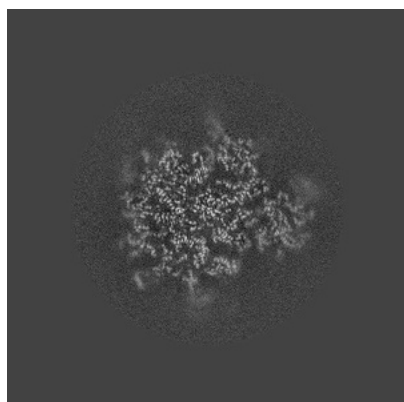


Y Index: 304

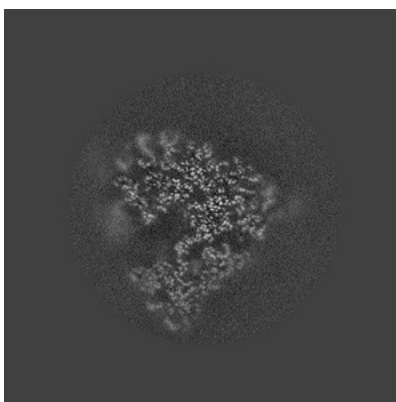


Z Index: 304

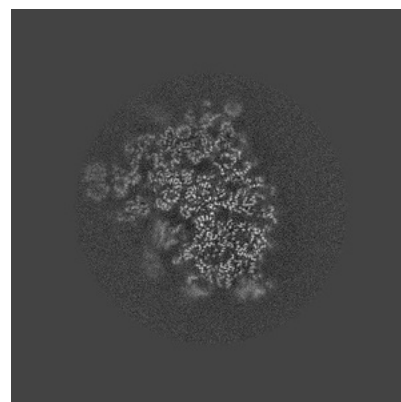
6.2.2 Raw map



X Index: 304



Y Index: 304

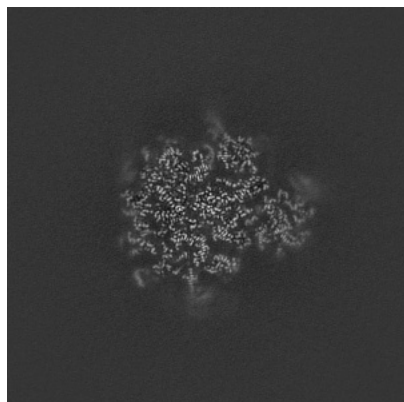


Z Index: 304

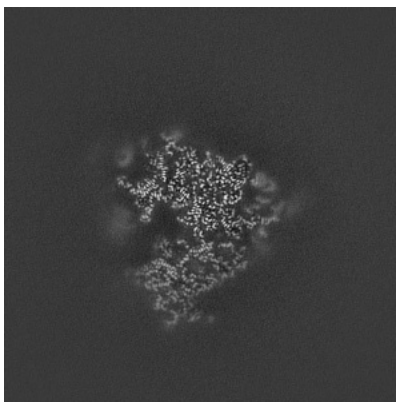
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

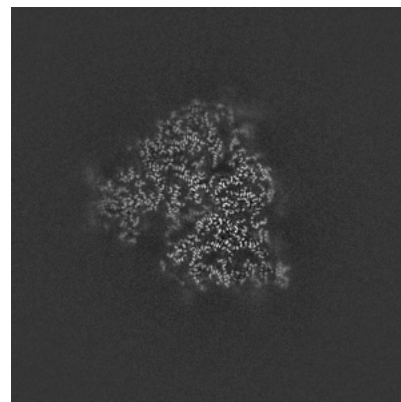
6.3.1 Primary map



X Index: 304

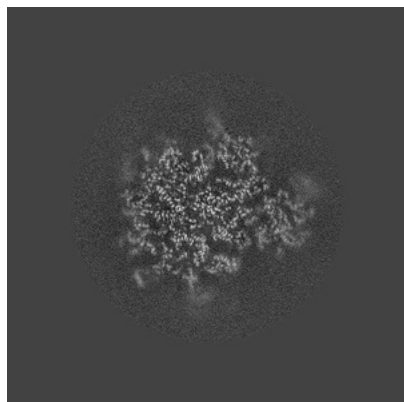


Y Index: 314

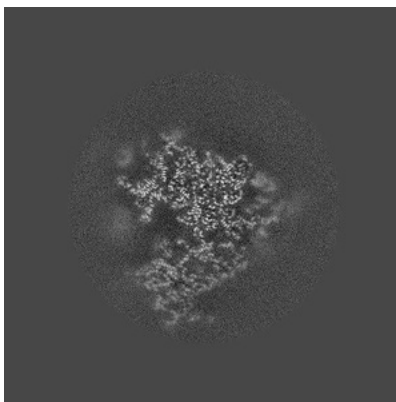


Z Index: 274

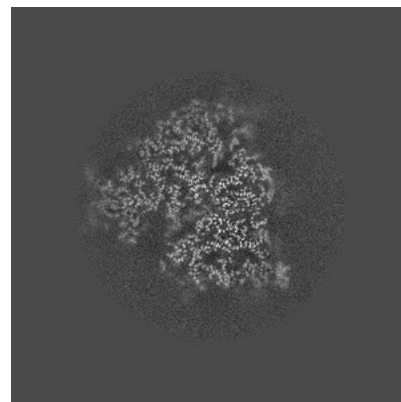
6.3.2 Raw map



X Index: 304



Y Index: 314

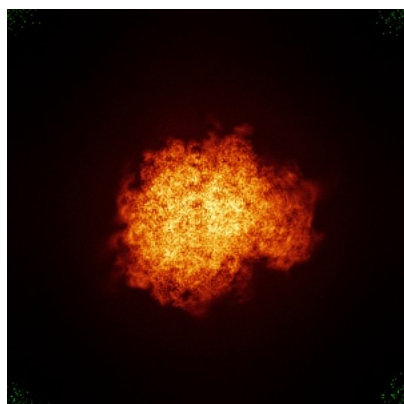


Z Index: 274

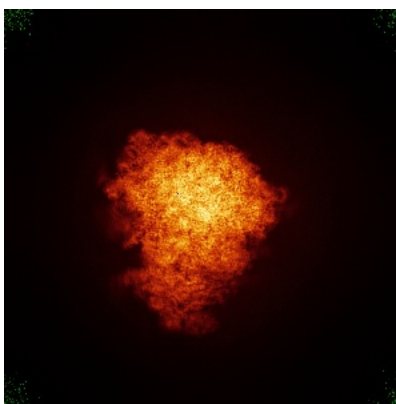
The images above show the largest variance slices of the map in three orthogonal directions.

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

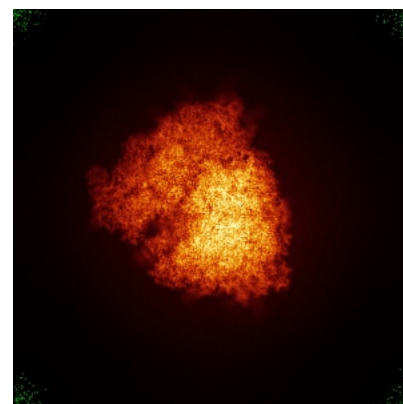
6.4.1 Primary map



X

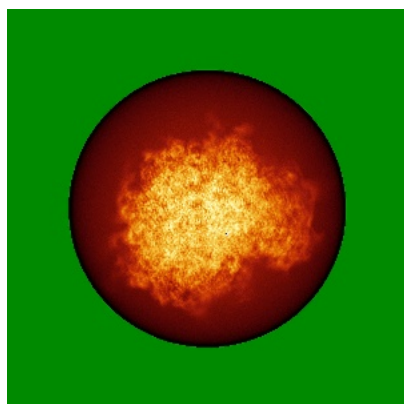


Y

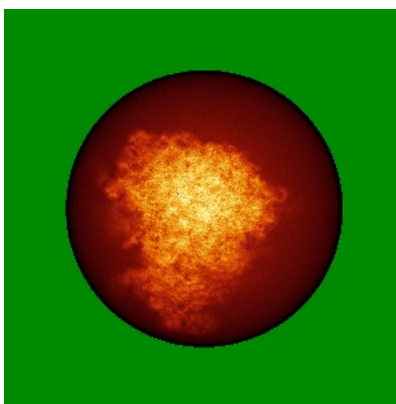


Z

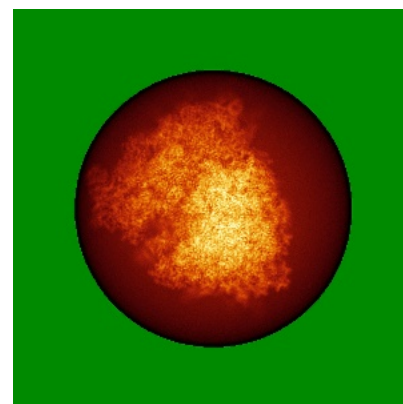
6.4.2 Raw map



X



Y



Z

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

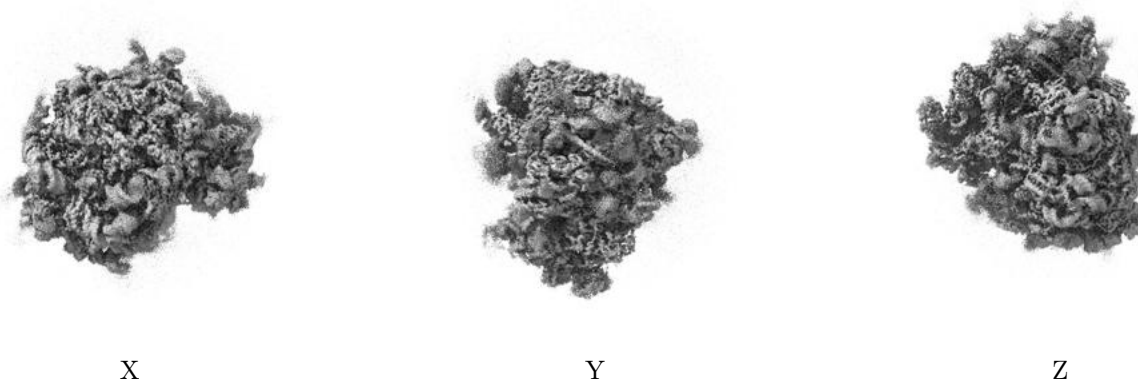
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 3.5. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

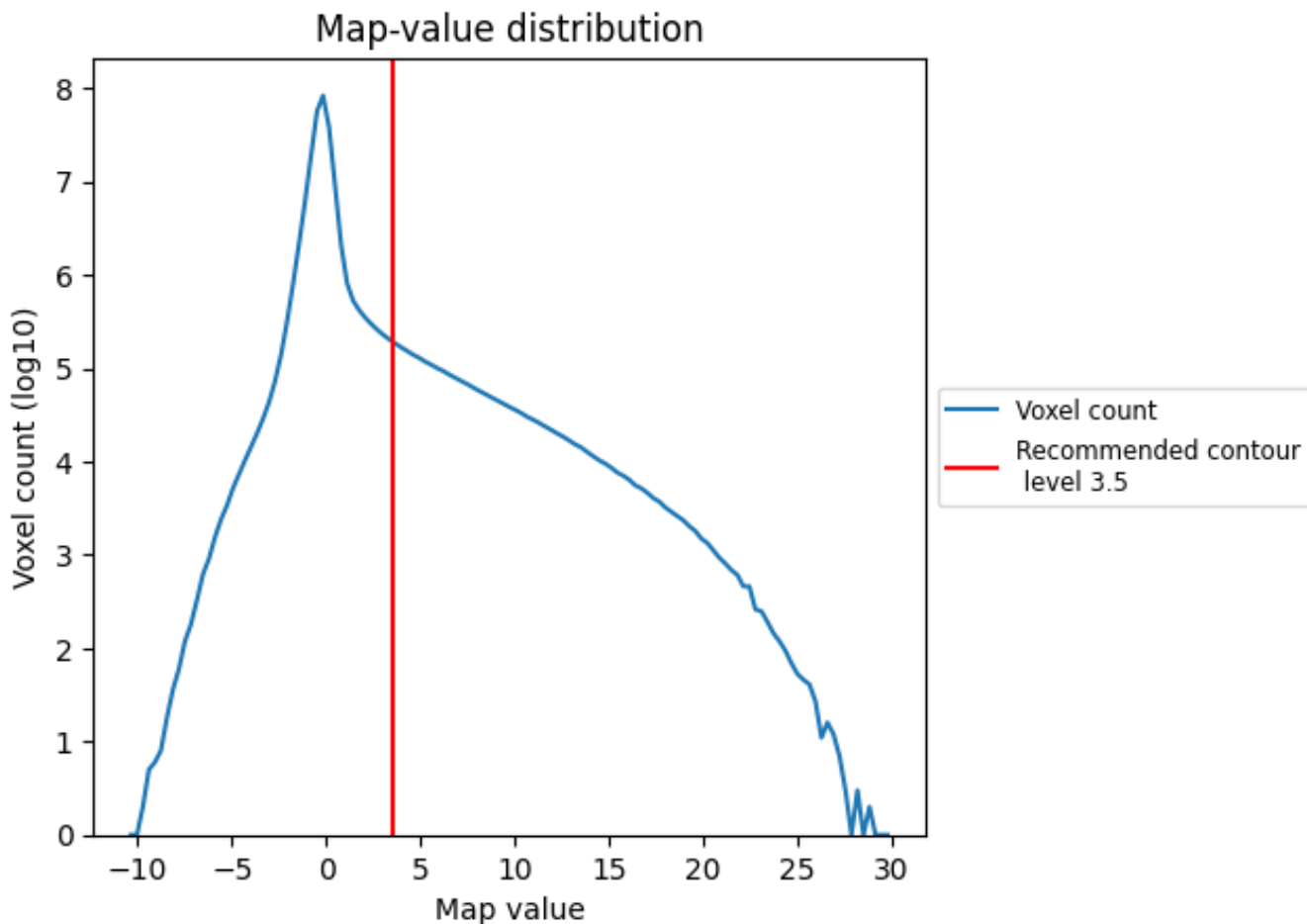
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

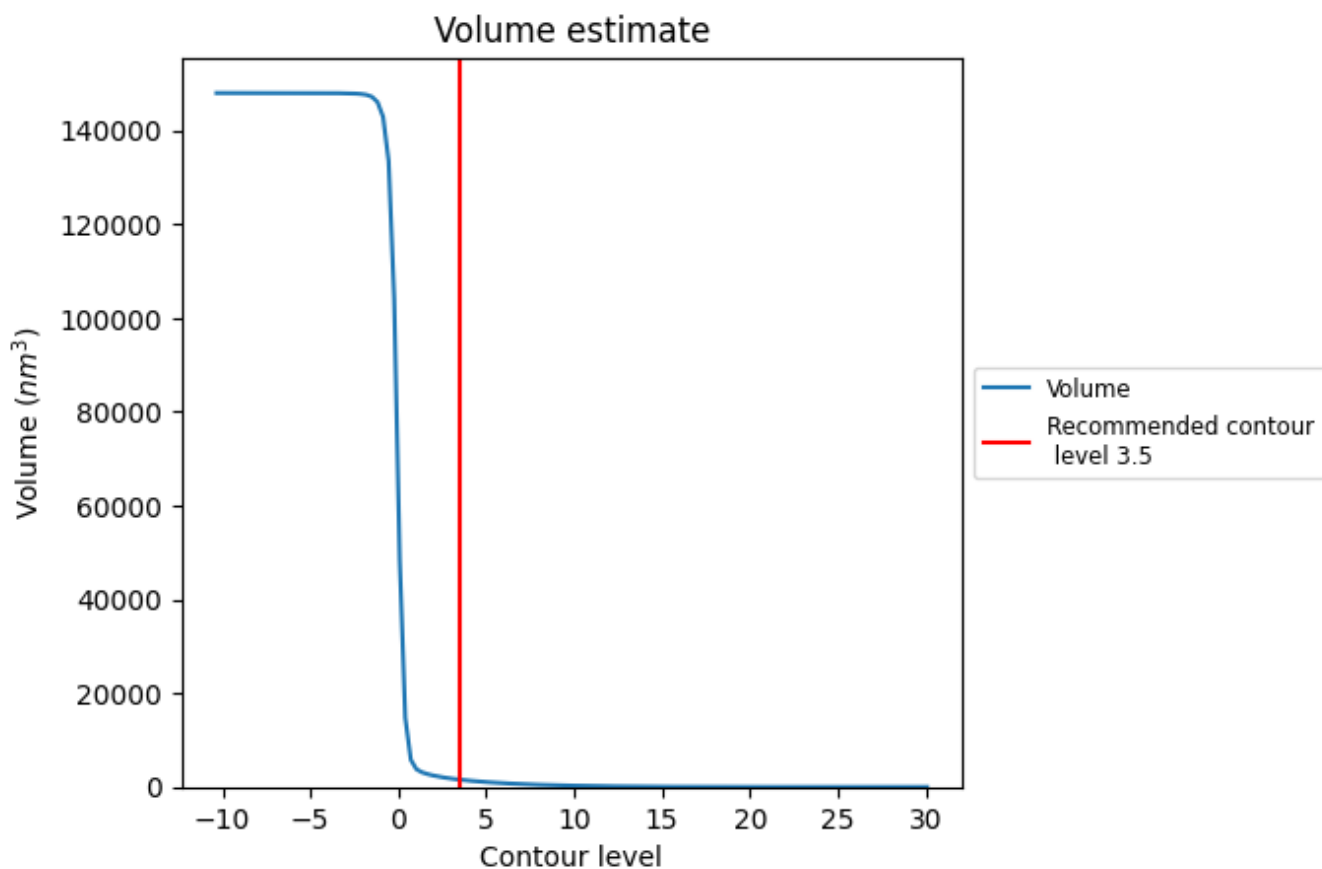
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

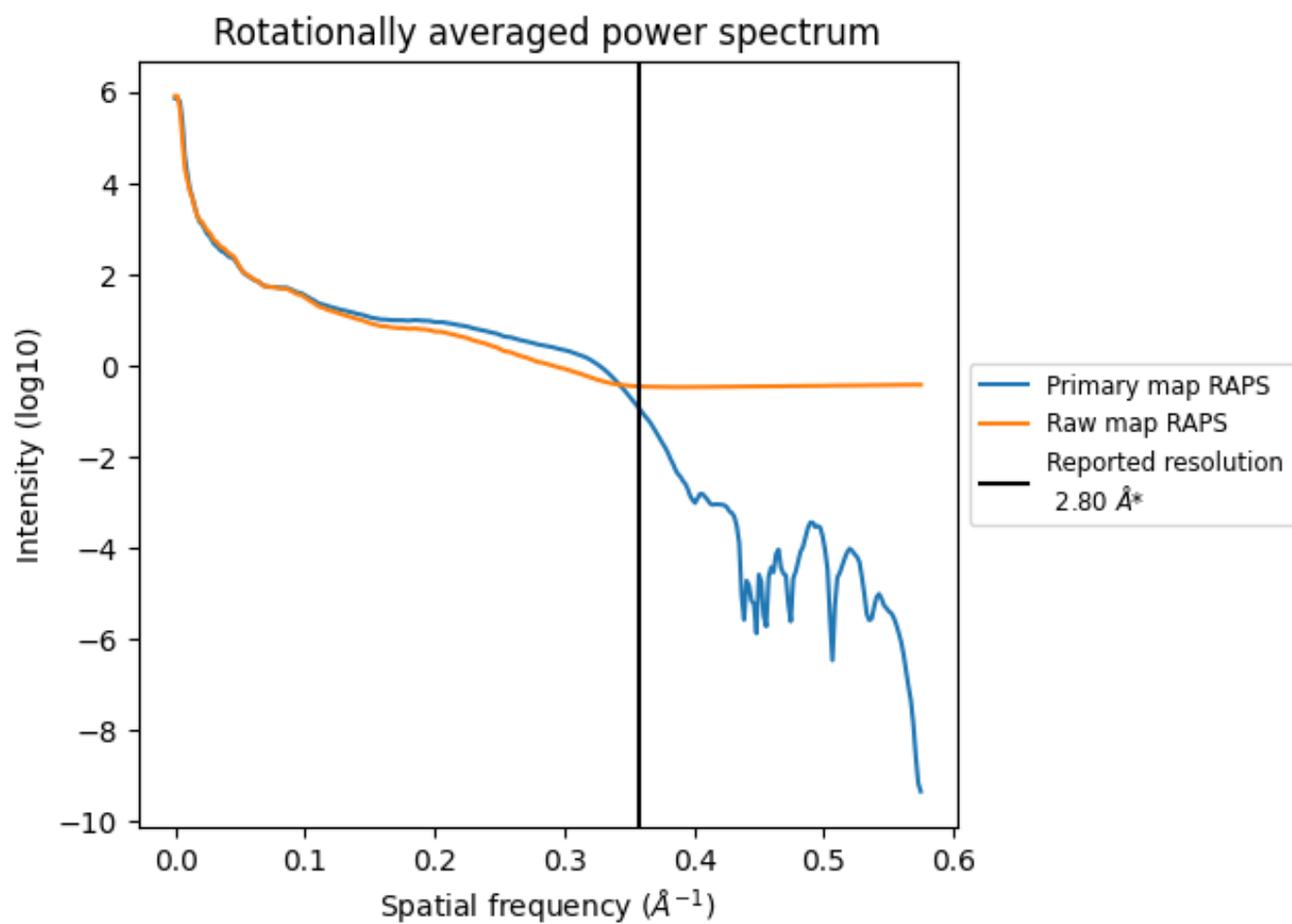
7.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 1555 nm^3 ; this corresponds to an approximate mass of 1404 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 [\(i\)](#)

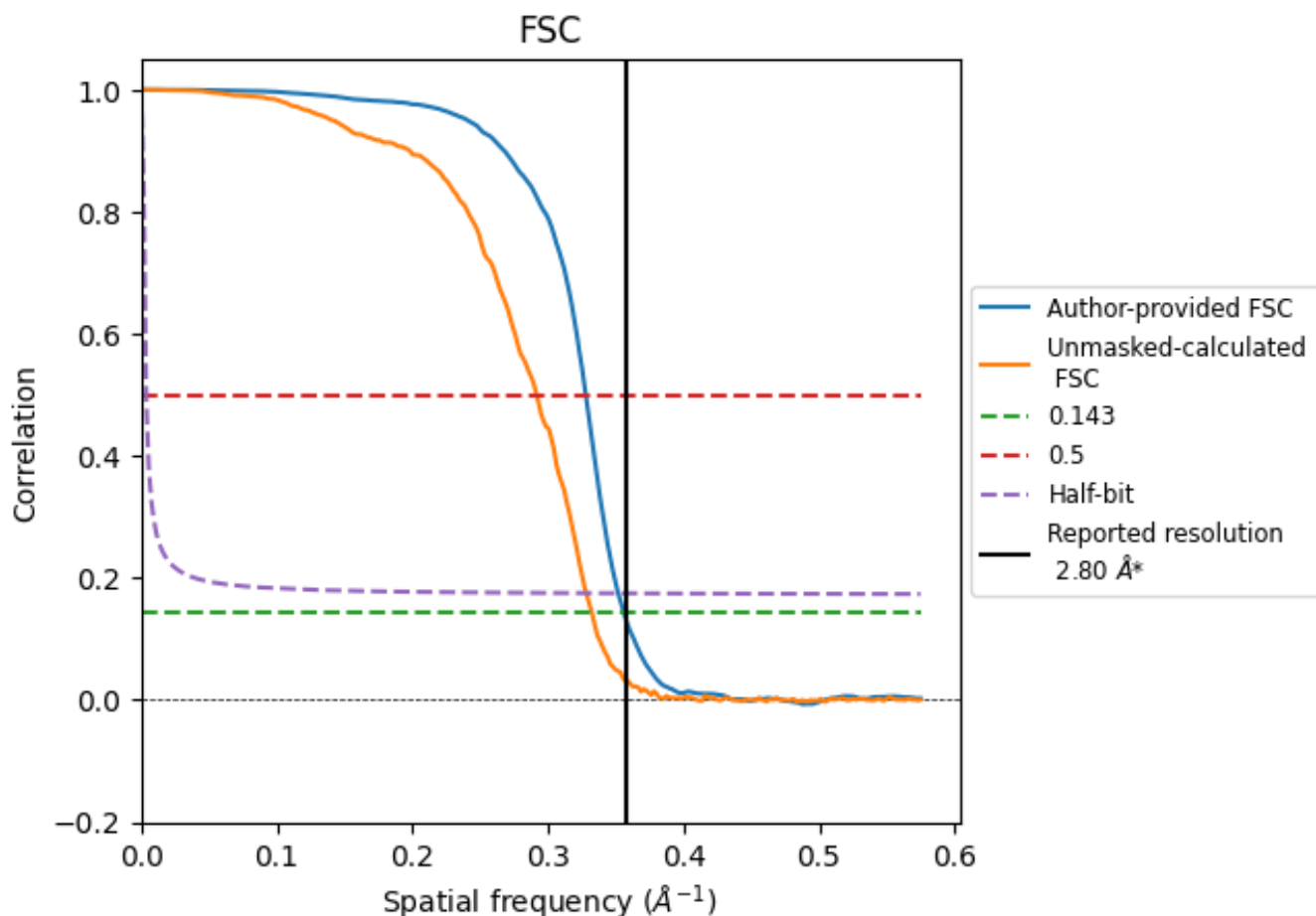


*Reported resolution corresponds to spatial frequency of 0.357 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.357 Å⁻¹

8.2 Resolution estimates [i](#)

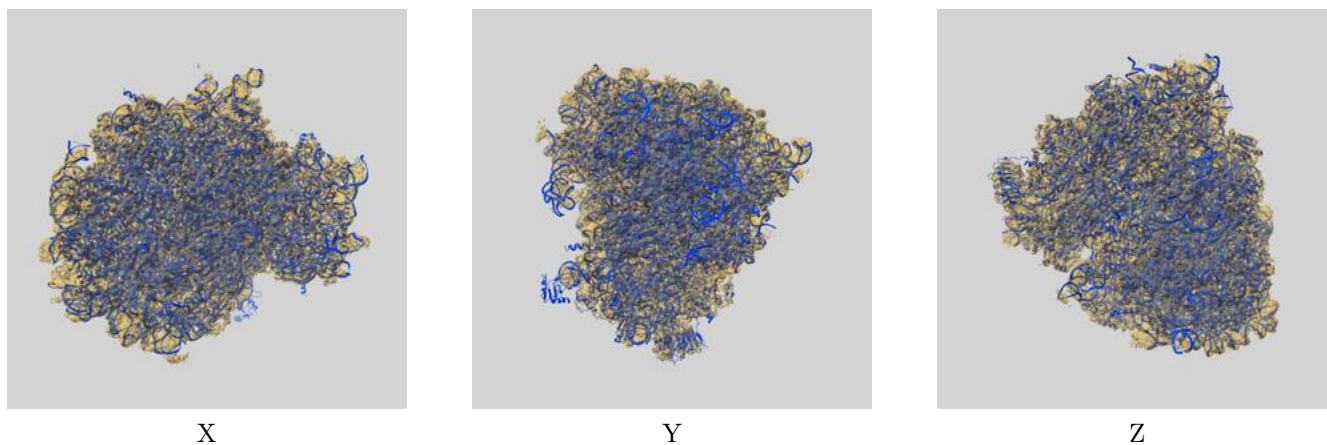
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.80	-	-
Author-provided FSC curve	2.81	3.05	2.84
Unmasked-calculated*	3.01	3.43	3.05

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

9 Map-model fit [i](#)

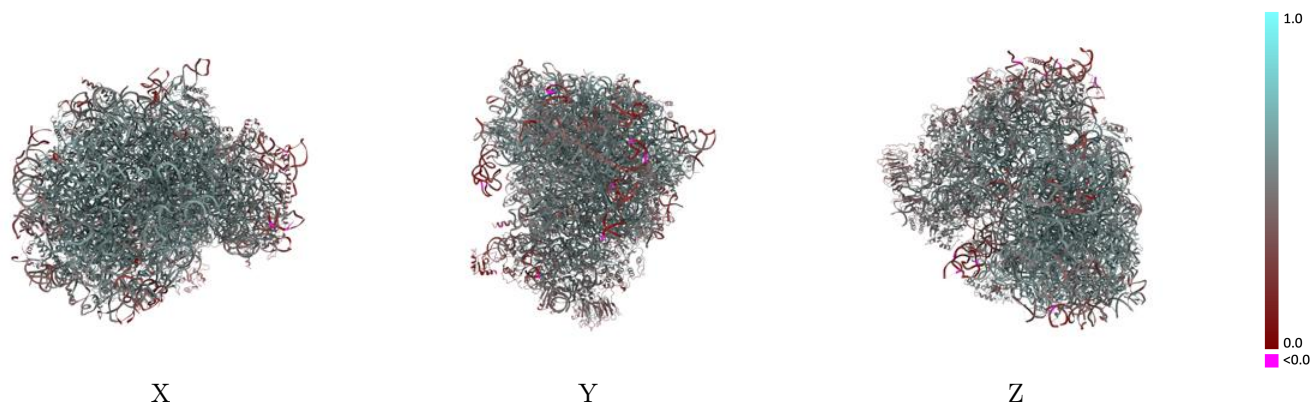
This section contains information regarding the fit between EMDB map EMD-44461 and PDB model 9BDL. Per-residue inclusion information can be found in section 3 on page 22.

9.1 Map-model overlay [i](#)



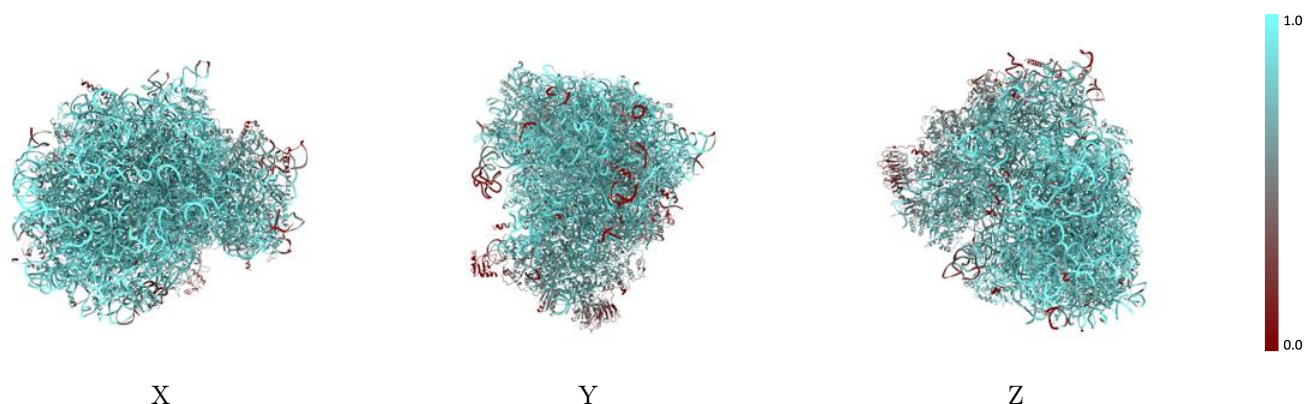
The images above show the 3D surface view of the map at the recommended contour level 3.5 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



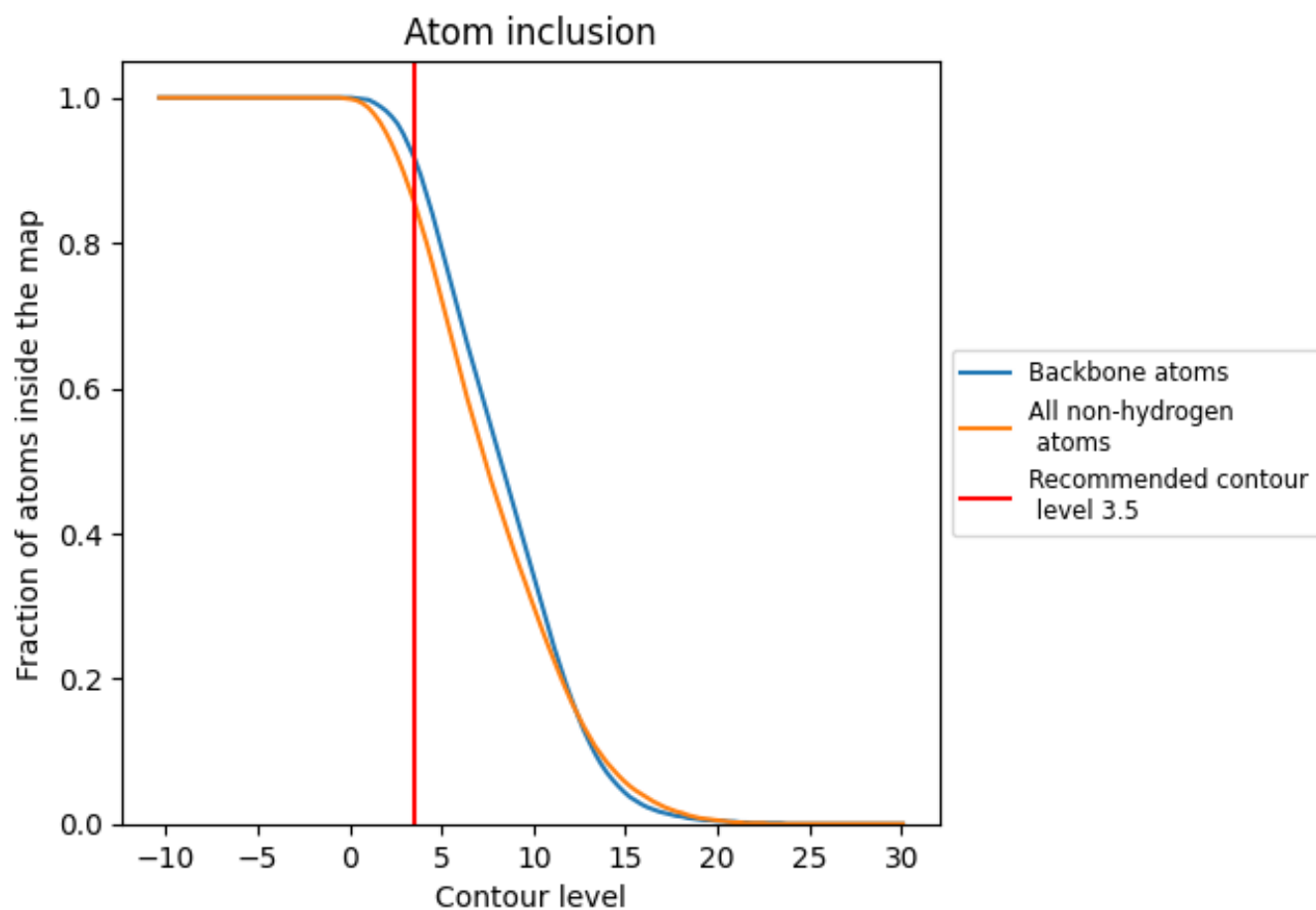
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (3.5).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 92% of all backbone atoms, 86% 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 (3.5) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8570	 0.5150
A18S	 0.9000	 0.5000
A28S	 0.9350	 0.5310
A58S	 0.9690	 0.5520
A5S	 0.9940	 0.5690
AL02	 0.9090	 0.5840
AL03	 0.8750	 0.5600
AL04	 0.8770	 0.5600
AL05	 0.8380	 0.5150
AL06	 0.8030	 0.5110
AL07	 0.8750	 0.5630
AL08	 0.7290	 0.5010
AL09	 0.7990	 0.5210
AL10	 0.8590	 0.5510
AL11	 0.8050	 0.5020
AL13	 0.8140	 0.5340
AL14	 0.8370	 0.5230
AL15	 0.9260	 0.5830
AL16	 0.8810	 0.5590
AL17	 0.8950	 0.5640
AL18	 0.8830	 0.5710
AL19	 0.8060	 0.5220
AL20	 0.8740	 0.5610
AL21	 0.8650	 0.5470
AL22	 0.7570	 0.4690
AL23	 0.8740	 0.5680
AL24	 0.8660	 0.5560
AL25	 0.8490	 0.5420
AL26	 0.8480	 0.5390
AL27	 0.8380	 0.5250
AL28	 0.9150	 0.5790
AL29	 0.7030	 0.4840
AL30	 0.8390	 0.5210
AL31	 0.8570	 0.5390
AL32	 0.8880	 0.5770















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Chain	Atom inclusion	Q-score
AL33	 0.9090	 0.5800
AL34	 0.8520	 0.5470
AL35	 0.8250	 0.5300
AL36	 0.8330	 0.5180
AL37	 0.9310	 0.5850
AL38	 0.7380	 0.4760
AL39	 0.8950	 0.5590
AL40	 0.8630	 0.5430
AL41	 0.7840	 0.5560
AL42	 0.8560	 0.5530
AL43	 0.8610	 0.5630
ALNW	 0.8830	 0.5590
ANG	 0.7230	 0.5010
ARAC	 0.3740	 0.3800
AS00	 0.6480	 0.4750
AS01	 0.7560	 0.5040
AS02	 0.7160	 0.4990
AS03	 0.5650	 0.4320
AS04	 0.7160	 0.4850
AS05	 0.7310	 0.4800
AS06	 0.5870	 0.4100
AS07	 0.5040	 0.4280
AS08	 0.7350	 0.4880
AS09	 0.7220	 0.4800
AS10	 0.4830	 0.3640
AS11	 0.7740	 0.5340
AS12	 0.0030	 0.1970
AS13	 0.7880	 0.5130
AS14	 0.7970	 0.5130
AS15	 0.6480	 0.4370
AS16	 0.6870	 0.4690
AS17	 0.5550	 0.4440
AS18	 0.7220	 0.4560
AS19	 0.6870	 0.4410
AS20	 0.5060	 0.4220
AS21	 0.6460	 0.4750
AS22	 0.8070	 0.5270
AS23	 0.8100	 0.5340
AS24	 0.6920	 0.4410
AS25	 0.6740	 0.4570
AS26	 0.8100	 0.5240
AS27	 0.6850	 0.4580

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Chain	Atom inclusion	Q-score
AS28	 0.7130	 0.4760
AS29	 0.6210	 0.4320
AS30	 0.6380	 0.4380
ETRN	 0.7450	 0.3420
MRNA	 0.6710	 0.4510
PTRN	 0.8890	 0.4990