



wwPDB EM Validation Summary Report ⓘ

Jul 1, 2025 – 05:49 pm BST

PDB ID : 9GXX / pdb_00009gxx
EMDB ID : EMD-51679
Title : 70S ribosome with doublet-decoding tRNA^{Ser3} bound to A-site GCA codon
Authors : Krishnaswamy, S.; Larsson, D.S.D.; Selmer, M.
Deposited on : 2024-10-01
Resolution : 2.61 Å (reported)
Based on initial models : 8CGJ, 8CF1, 8CGK, 7K00

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.dev118
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0rc1
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.44

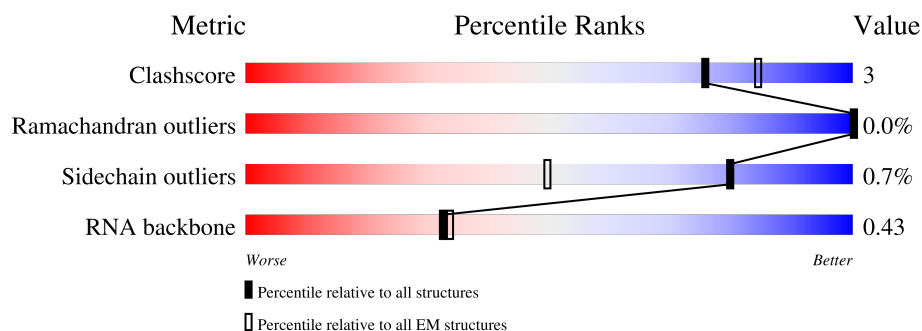
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.61 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.



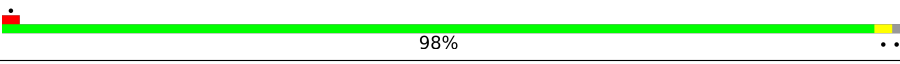
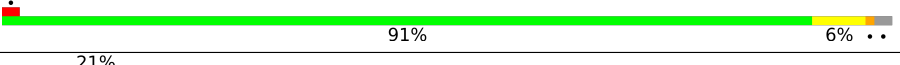

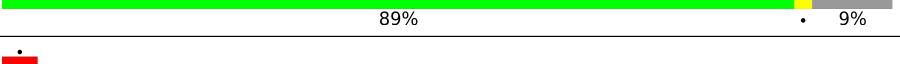
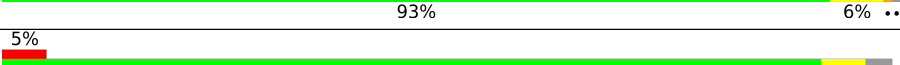
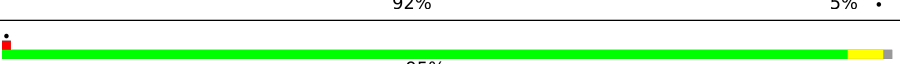
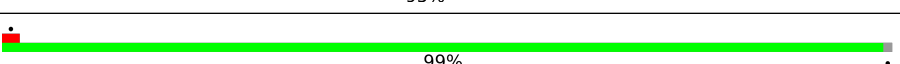
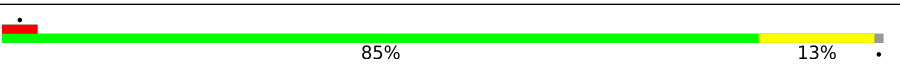


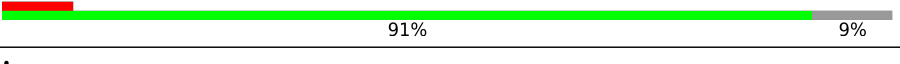
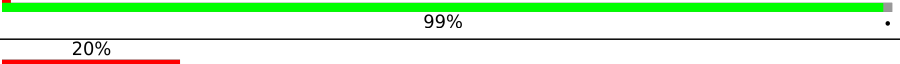

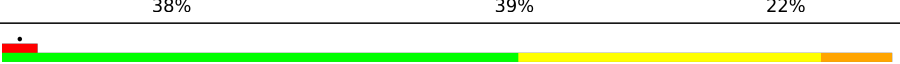
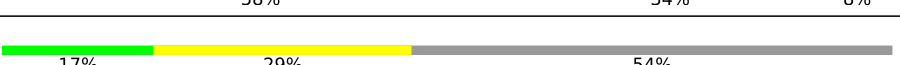
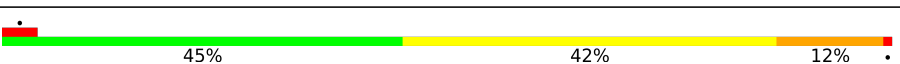


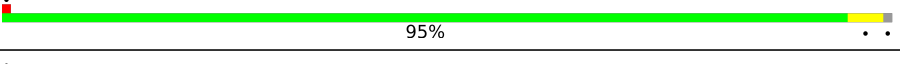
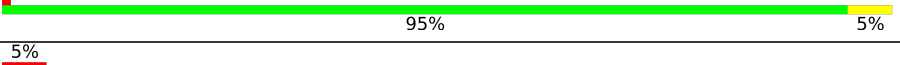
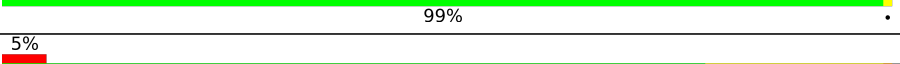

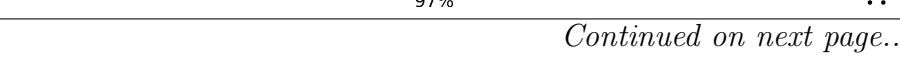


Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415
RNA backbone	6643	2191

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 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	1542	
2	B	241	
3	C	233	
4	D	206	
5	E	167	
6	F	135	
7	G	179	


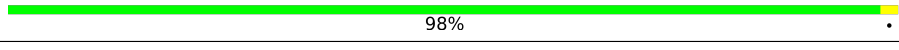
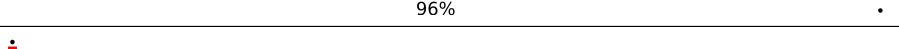
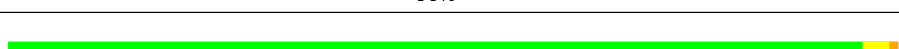
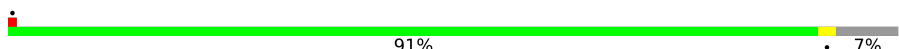

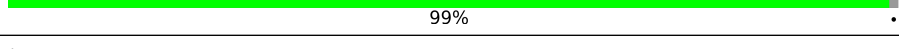
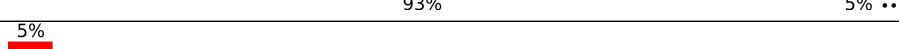
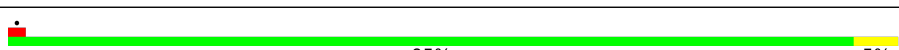

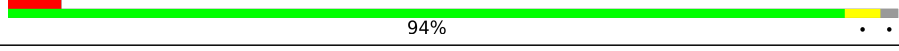
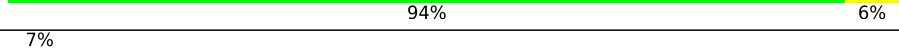
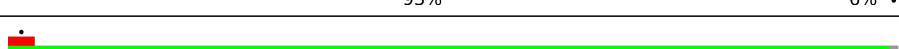
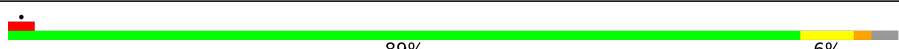
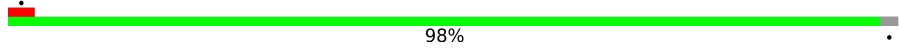
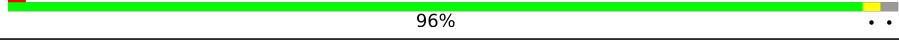
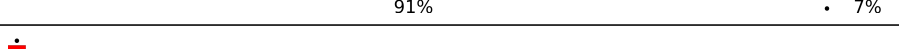







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Mol	Chain	Length	Quality of chain
8	H	130	
9	I	130	
10	J	103	
11	K	129	
12	L	124	
13	M	118	
14	N	101	
15	O	89	
16	P	82	
17	Q	84	
18	R	75	
19	S	92	
20	T	87	
21	U	71	
22	V	76	
22	Z	76	
23	X	24	
24	Y	93	
25	a	2904	
26	b	120	
27	c	273	
28	d	209	
29	e	201	
30	f	179	
31	g	177	

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Mol	Chain	Length	Quality of chain
32	h	149	
33	i	142	
34	j	123	
35	k	144	
36	l	136	
37	m	127	
38	n	117	
39	o	115	
40	p	118	
41	q	103	
42	r	110	
43	s	100	
44	t	104	
45	u	94	
46	v	85	
47	w	78	
48	x	63	
49	y	59	
50	z	57	
51	0	55	
52	1	46	
53	2	65	
54	3	38	
55	4	70	

2 Entry composition

There are 60 unique types of molecules in this entry. The entry contains 144033 atoms, of which 0 are hydrogens and 0 are deuteriums.

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

- Molecule 1 is a RNA chain called 16S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	1520	Total	C	N	O	P	0	0
			32634	14562	5991	10561	1520		

- Molecule 2 is a protein called Small ribosomal subunit protein uS2.

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	206	Total	C	N	O	S	0	0
			1624	1028	305	288	3		

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	156	Total	C	N	O	S	0	0
			1152	717	217	212	6		

- Molecule 6 is a protein called Small ribosomal subunit protein bS6, fully modified isoform.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	103	Total	C	N	O	S	0	0
			839	530	151	151	7		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	153	Total	C	N	O	S	0	0
			1203	750	231	218	4		

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

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	98	Total	C	N	O	S	0	0
			786	493	150	142	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	117	Total	C	N	O	S	0	0
			877	540	173	161	3		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	119	IAS	ASN	conflict	UNP P0A7R9

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
13	M	115	Total	C	N	O	S	0	0
			891	552	179	157	3		

- Molecule 14 is a protein called Small ribosomal subunit protein uS14.

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

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

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

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

Mol	Chain	Residues	Atoms					AltConf	Trace
16	P	81	Total	C	N	O	S	0	0
			643	403	127	112	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
17	Q	79	Total	C	N	O	S	0	0
			641	406	120	112	3		

- Molecule 18 is a protein called Small ribosomal subunit protein bS18.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	R	66	Total	C	N	O	S	0	0
			544	345	102	96	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
19	S	84	Total	C	N	O	S	0	0
			668	427	127	112	2		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
20	T	86	Total	C	N	O	S	0	0
			670	414	138	115	3		

- Molecule 21 is a protein called Small ribosomal subunit protein bS21.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	U	70	Total	C	N	O	S	0	0
			589	366	125	97	1		

- Molecule 22 is a RNA chain called tRNA(fMet).

Mol	Chain	Residues	Atoms					AltConf	Trace
22	V	76	Total	C	N	O	P	0	0
			1623	723	294	530	76		
22	Z	76	Total	C	N	O	P	0	0
			1623	723	294	530	76		

- Molecule 23 is a RNA chain called mRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	X	11	Total	C	N	O	P	0	0
			240	108	50	71	11		

- Molecule 24 is a RNA chain called tRNA(Ser3).

Mol	Chain	Residues	Atoms					AltConf	Trace	
24	Y	93	Total	C	N	O	P	S	0	0
			2000	893	365	648	92	2		

- Molecule 25 is a RNA chain called 23S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	a	2753	Total	C	N	O	P	2	0
			59172	26403	10905	19109	2755		

- Molecule 26 is a RNA chain called 5S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	b	119	Total	C	N	O	P	0	0
			2549	1135	466	829	119		

- Molecule 27 is a protein called Large ribosomal subunit protein uL2.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	c	271	Total	C	N	O	S	0	0
			2082	1288	423	364	7		

- Molecule 28 is a protein called Large ribosomal subunit protein uL3.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	d	209	Total	C	N	O	S	0	0
			1566	980	288	294	4		

- Molecule 29 is a protein called Large ribosomal subunit protein uL4.

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

- Molecule 30 is a protein called Large ribosomal subunit protein uL5.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	f	177	Total	C	N	O	S	0	0
			1410	899	249	256	6		

- Molecule 31 is a protein called Large ribosomal subunit protein uL6.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	g	173	Total	C	N	O	S	0	0
			1295	814	237	242	2		

- Molecule 32 is a protein called Large ribosomal subunit protein bL9.

Mol	Chain	Residues	Atoms					AltConf	Trace
32	h	41	Total	C	N	O	S	0	0
			303	194	54	54	1		

- Molecule 33 is a protein called Large ribosomal subunit protein uL13.

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

- Molecule 34 is a protein called Large ribosomal subunit protein uL14.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	j	123	Total	C	N	O	S	0	0
			946	593	181	166	6		

- Molecule 35 is a protein called Large ribosomal subunit protein uL15.

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

- Molecule 36 is a protein called Large ribosomal subunit protein uL16.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	l	136	Total	C	N	O	S	0	0
			1075	686	205	177	7		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
l	82	MS6	MET	conflict	UNP P0ADY7

- Molecule 37 is a protein called Large ribosomal subunit protein bL17.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	m	118	Total	C	N	O	S	0	0
			945	585	194	161	5		

- Molecule 38 is a protein called Large ribosomal subunit protein uL18.

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 Large ribosomal subunit protein bL19.

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 Large ribosomal subunit protein bL20.

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 Large ribosomal subunit protein bL21.

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 Large ribosomal subunit protein uL22.

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 Large ribosomal subunit protein uL23.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	s	93	Total	C	N	O	S	0	0
			738	466	139	131	2		

- Molecule 44 is a protein called Large ribosomal subunit protein uL24.

Mol	Chain	Residues	Atoms				AltConf	Trace
44	t	102	Total	C	N	O	0	0
			779	492	146	141		

- Molecule 45 is a protein called Large ribosomal subunit protein bL25.

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 Large ribosomal subunit protein bL27.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	v	84	Total	C	N	O	S	0	0
			628	388	126	113	1		

- Molecule 47 is a protein called Large ribosomal subunit protein bL28.

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 Large ribosomal subunit protein uL29.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	x	61	Total	C	N	O	S	0	0
			495	305	97	92	1		

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

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

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

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

- Molecule 51 is a protein called Large ribosomal subunit protein bL33.

Mol	Chain	Residues	Atoms				AltConf	Trace
51	0	51	Total	C	N	O	0	0
			417	269	76	72		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
52	1	46	Total	C	N	O	S	0	0
			377	228	90	57	2		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
53	2	64	Total	C	N	O	S	0	0
			504	323	105	74	2		

- Molecule 54 is a protein called Large ribosomal subunit protein bL36A.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	3	38	Total	C	N	O	S	0	0
			302	185	65	48	4		

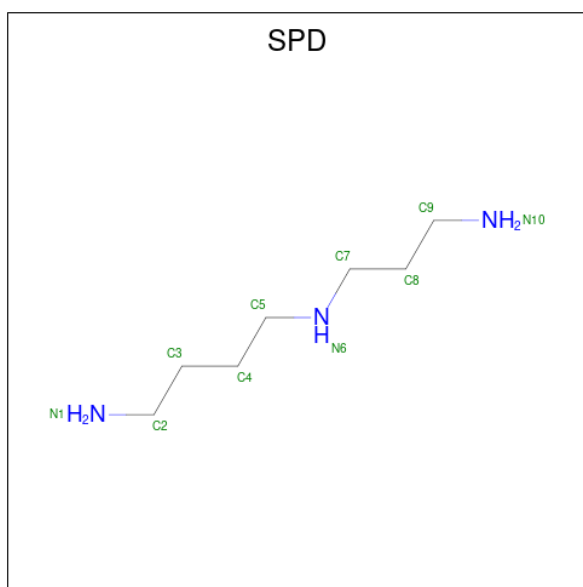
- Molecule 55 is a protein called Large ribosomal subunit protein bL31A.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	4	60	Total	C	N	O	S	0	0
			480	299	90	85	6		

- Molecule 56 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

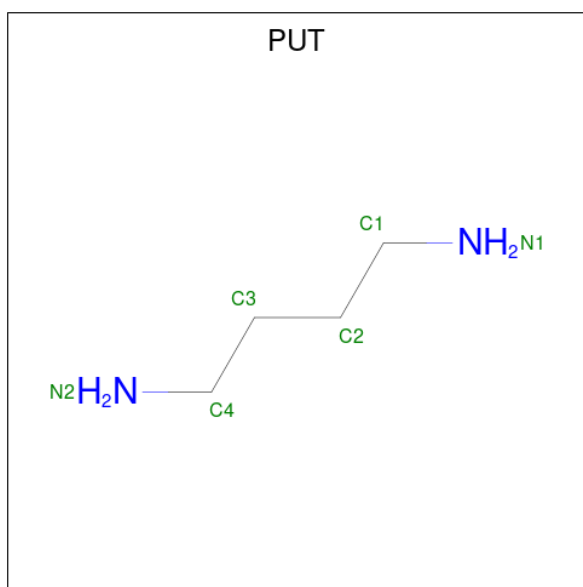
Mol	Chain	Residues	Atoms		AltConf
56	A	87	Total	Mg	0
			87	87	
56	Y	4	Total	Mg	0
			4	4	
56	a	187	Total	Mg	0
			187	187	
56	b	5	Total	Mg	0
			5	5	
56	c	1	Total	Mg	0
			1	1	
56	d	2	Total	Mg	0
			2	2	
56	m	1	Total	Mg	0
			1	1	
56	z	1	Total	Mg	0
			1	1	

- Molecule 57 is SPERMIDINE (CCD ID: SPD) (formula: C₇H₁₉N₃).



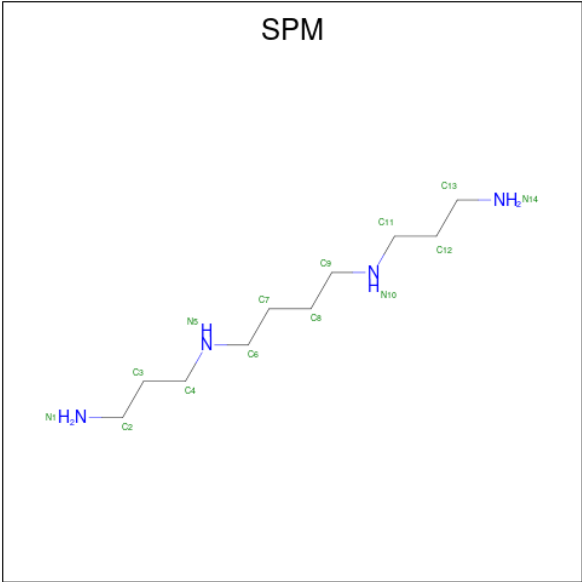
Mol	Chain	Residues	Atoms			AltConf
57	A	1	Total	C	N	0
			10	7	3	
57	a	1	Total	C	N	0
			10	7	3	
57	a	1	Total	C	N	0
			10	7	3	
57	a	1	Total	C	N	0
			10	7	3	
57	a	1	Total	C	N	0
			10	7	3	
57	a	1	Total	C	N	0
			10	7	3	

- Molecule 58 is 1,4-DIAMINOBTANE (CCD ID: PUT) (formula: C₄H₁₂N₂).



Mol	Chain	Residues	Atoms			AltConf
58	A	1	Total	C	N	0
			6	4	2	
58	a	1	Total	C	N	0
			6	4	2	
58	a	1	Total	C	N	0
			6	4	2	
58	a	1	Total	C	N	0
			6	4	2	
58	a	1	Total	C	N	0
			6	4	2	
58	a	1	Total	C	N	0
			6	4	2	

- Molecule 59 is SPERMINE (CCD ID: SPM) (formula: C₁₀H₂₆N₄).



Mol	Chain	Residues	Atoms			AltConf
59	a	1	Total	C	N	0
			14	10	4	

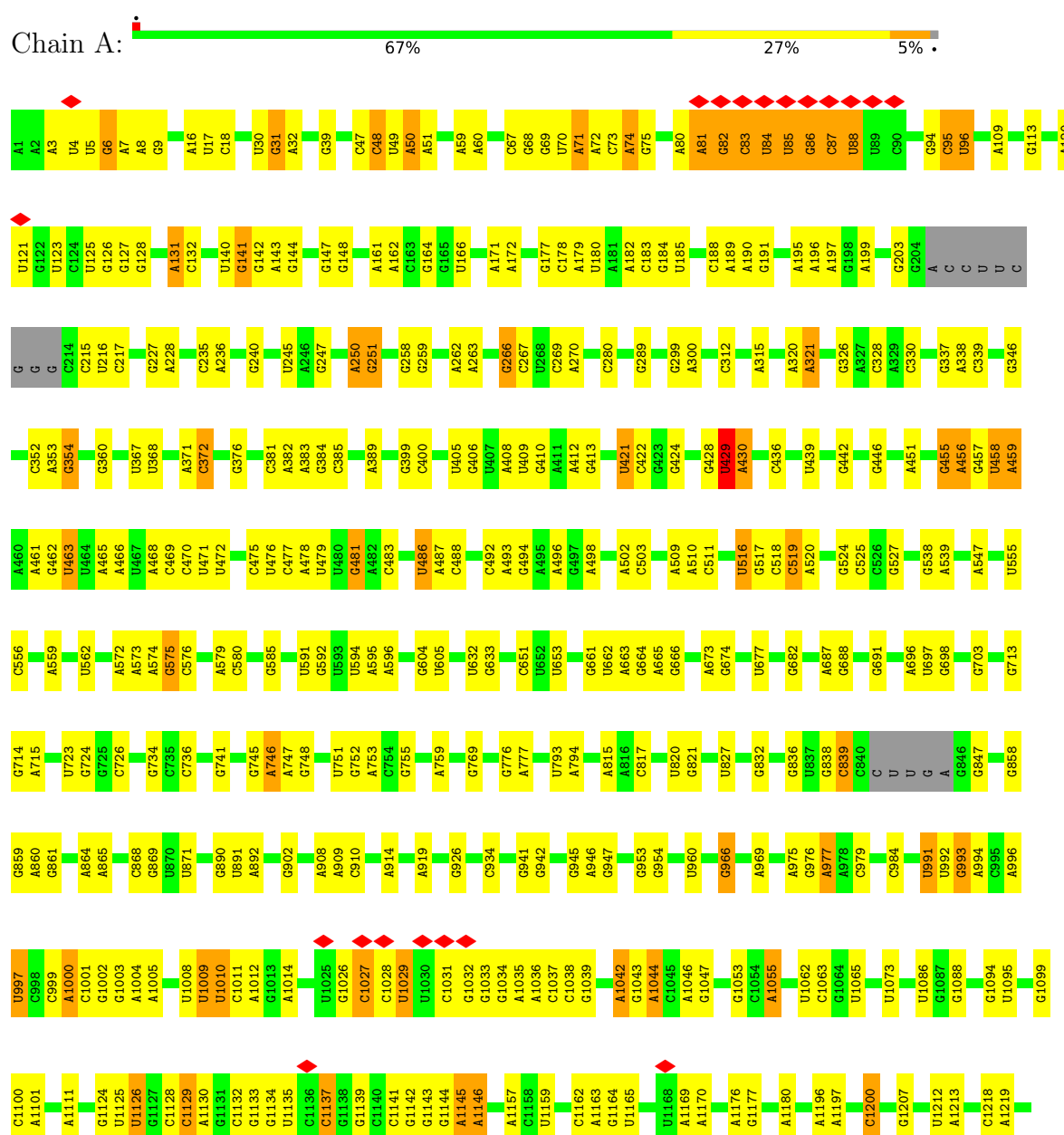
- Molecule 60 is ZINC ION (CCD ID: ZN) (formula: Zn).

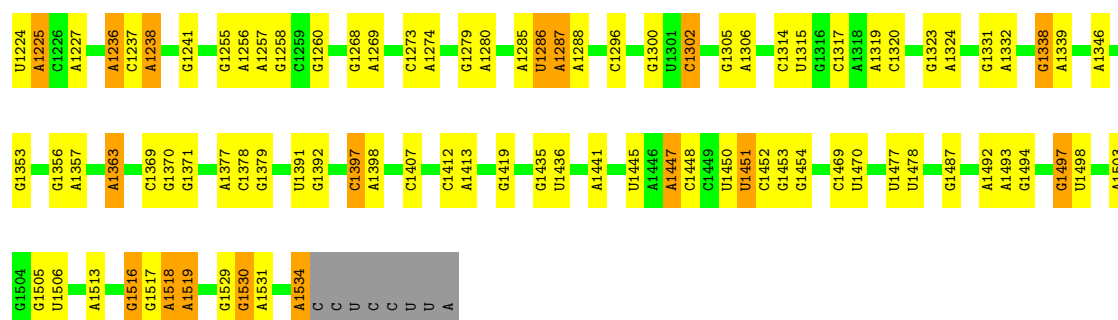
Mol	Chain	Residues	Atoms		AltConf
60	3	1	Total	Zn	0
			1	1	
60	4	1	Total	Zn	0
			1	1	

3 Residue-property plots

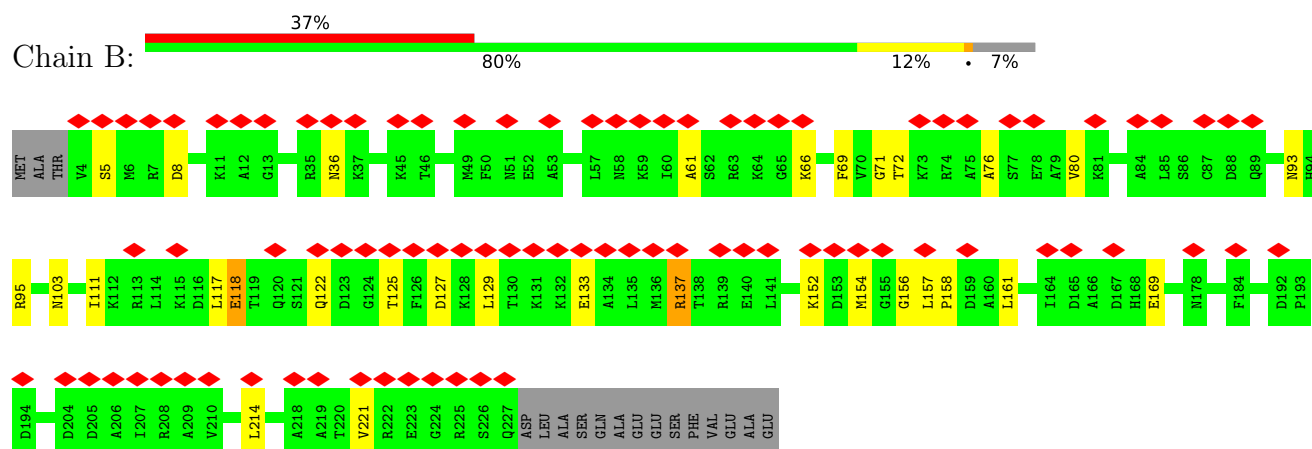
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

• Molecule 1: 16S rRNA

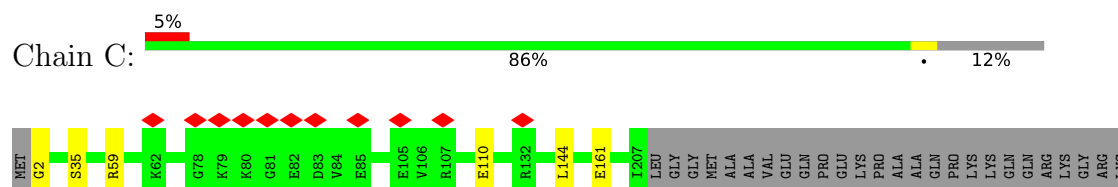




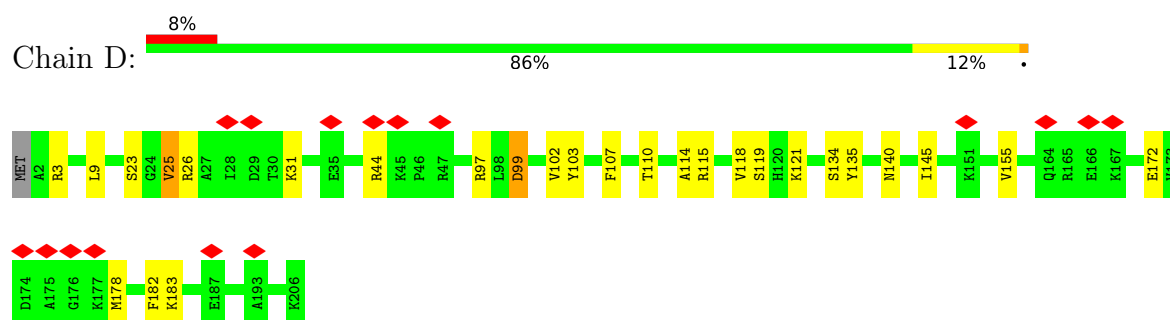
• Molecule 2: Small ribosomal subunit protein uS2



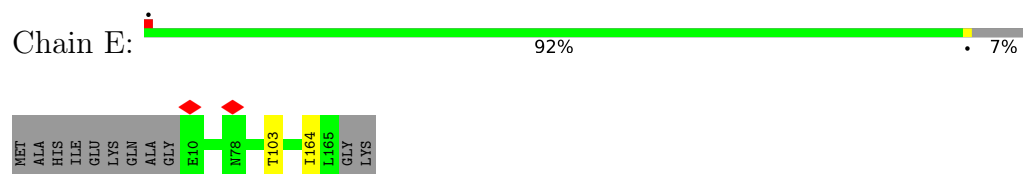
• Molecule 3: Small ribosomal subunit protein uS3



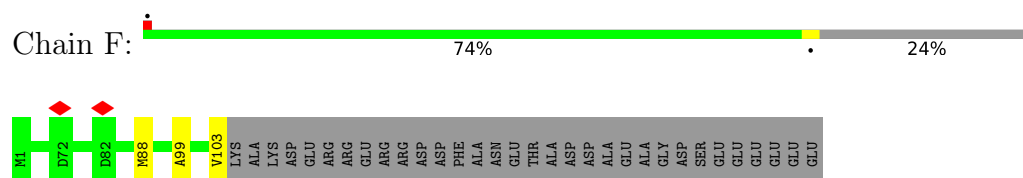
• Molecule 4: Small ribosomal subunit protein uS4



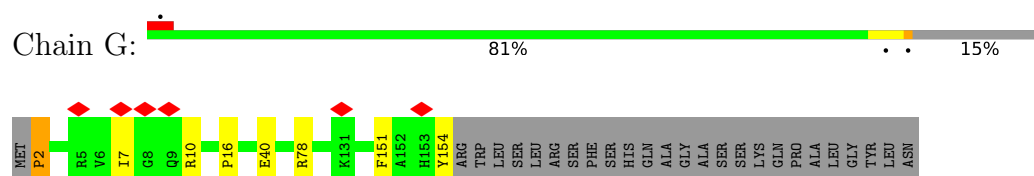
• Molecule 5: Small ribosomal subunit protein uS5



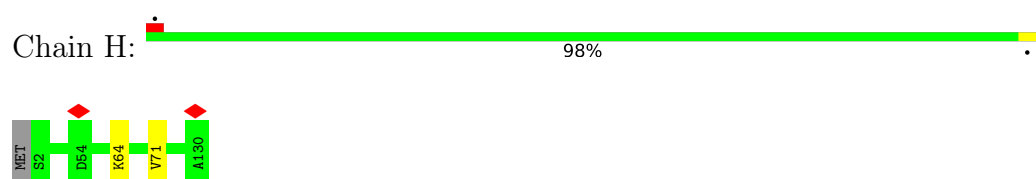
- Molecule 6: Small ribosomal subunit protein bS6, fully modified isoform



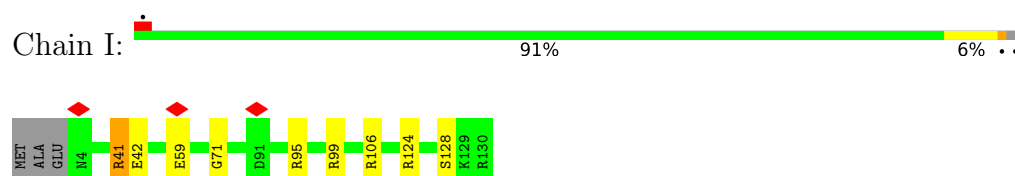
- Molecule 7: Small ribosomal subunit protein uS7



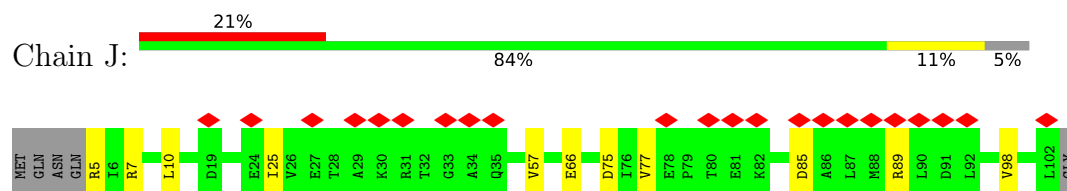
- Molecule 8: Small ribosomal subunit protein uS8



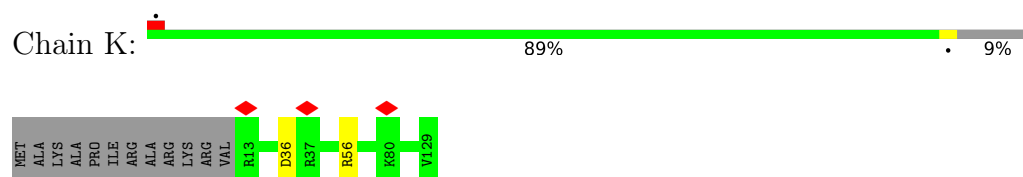
- Molecule 9: Small ribosomal subunit protein uS9



- Molecule 10: Small ribosomal subunit protein uS10

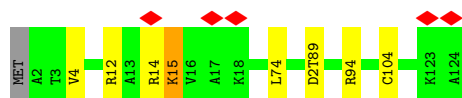


- Molecule 11: Small ribosomal subunit protein uS11

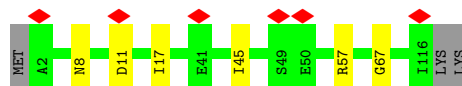
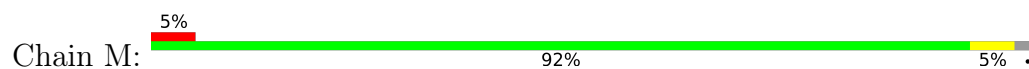


- Molecule 12: Small ribosomal subunit protein uS12

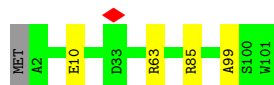




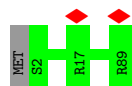
- Molecule 13: Small ribosomal subunit protein uS13



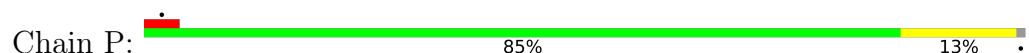
- Molecule 14: Small ribosomal subunit protein uS14



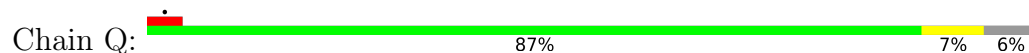
- Molecule 15: Small ribosomal subunit protein uS15



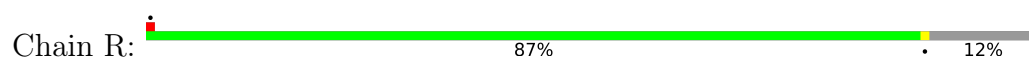
- Molecule 16: Small ribosomal subunit protein bS16



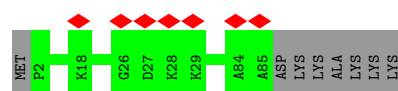
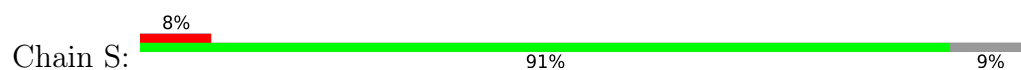
- Molecule 17: Small ribosomal subunit protein uS17



- Molecule 18: Small ribosomal subunit protein bS18



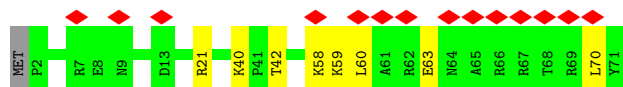
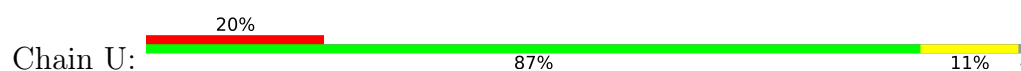
- Molecule 19: Small ribosomal subunit protein uS19



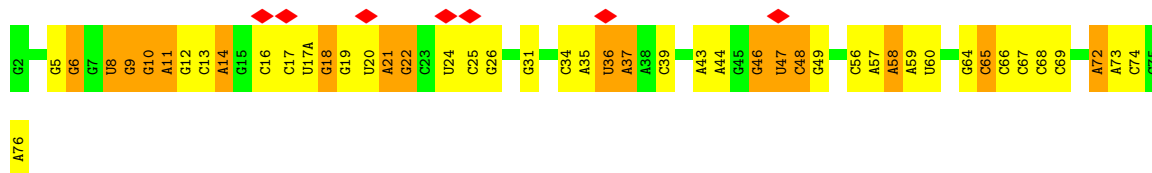
- Molecule 20: Small ribosomal subunit protein bS20



- Molecule 21: Small ribosomal subunit protein bS21



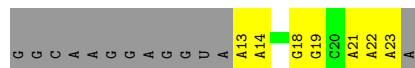
- Molecule 22: tRNA(fMet)



- Molecule 22: tRNA(fMet)

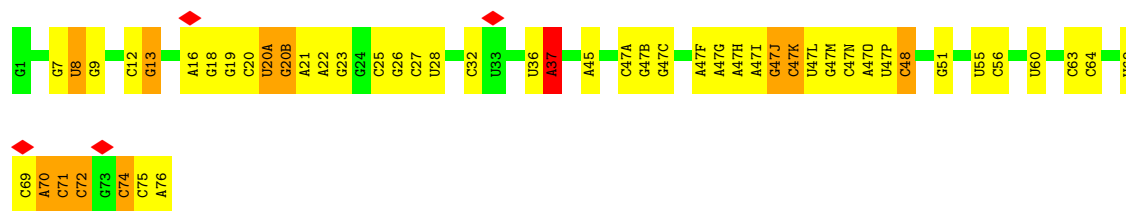


- Molecule 23: mRNA

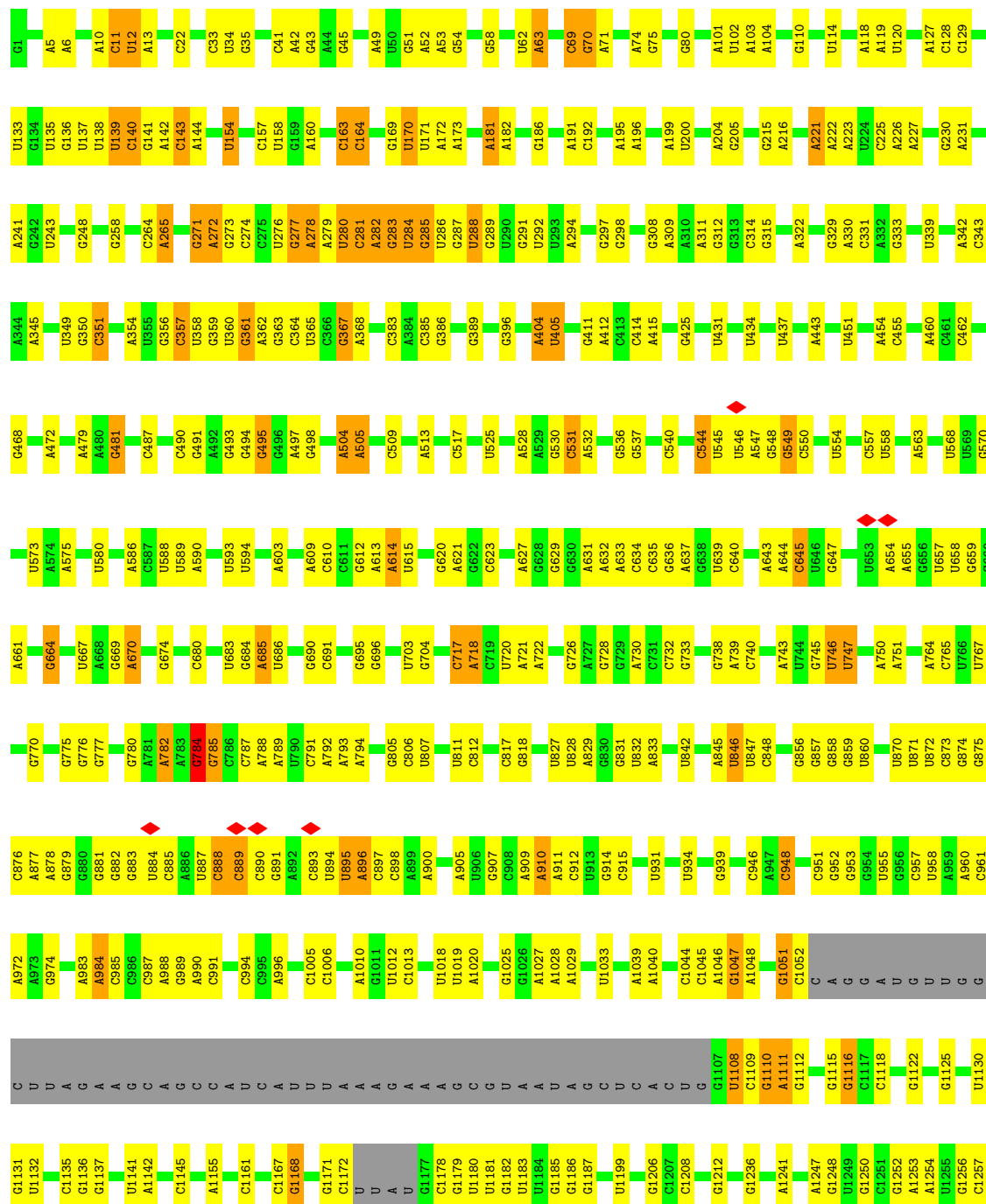


- Molecule 24: tRNA(Ser3)





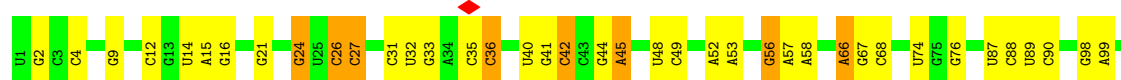
● Molecule 25: 23S rRNA



A2657	U2548	U2441	A2322	G2204	G	G2067	U1963	A1847	C1732	A1593	A1495	A1383	A1268
A2660	U2552	G2445	G2323	G2209	U	U2068	G1964	A1848	U1736	A1594	A1496	C1386	A1271
G2663	C2556	G2445	U2324	U2210	G	G1849	G1965	G1850	G	A1597	C1497	A1387	A1272
C2666	A2448	A2448	G2326	A2211	A	C2072	A1966	G1857	G1738	U1602	A1505	U1391	U1273
A2670	U2449	U2449	A2327	A2212	C	A2077	G1968	G1857	A1739	U1606	U1506	U1405	A1274
A2682	A2451	A2451	A2328	U2220	C	C2078	A1970	A1858	G1740	C1607	A1509	U1406	A1275
C2683	C2456	C2456	A2333	G2221	C	U2079	U1971	U1865	A1746	A1608	G1510	G1407	A1286
U2689	U2457	U2457	A2335	C2222	A	A2080	G1972	A1866	U1747	C1611	U1513	A1287	A1287
U2690	C2481	C2481	A2336	A2225	U	U2085	G1973	G1867	G1750	G1612	U1514	U1411	G1288
C2715	G2488	G2488	C2339	A2230	C	C2091	G1980	G1868	G1756	G1613	A1515	U1414	C1289
C2716	U2491	U2491	A2340	G2230	U	U2092	C1985	G1869	A1757	A1618	U1514	C1414	C1300
C2717	U2492	U2492	C2341	G2231	G	G2093	C1986	A1870	U1758	G1619	G1524	U1415	A1301
U2724	U2493	U2493	U2342	G2232	C	A2094	U1991	A1871	A1759	C1417	G1524	U1416	A1301
A2725	G2494	G2494	U2343	G2233	A	U2095	G1992	A1872	C1760	A1626	A1528	G1418	A1304
A2726	C2495	C2495	A2344	U2243	G	C2096	U1993	G1873	C1764	A1634	G1529	A1419	C1305
G2732	C2496	C2496	U2345	U2244	G	U2097	G1994	G1874	U1765	A1637	U1534	A1420	C1306
A2733	U2497	U2497	C2346	G2251	A	U2098	U1995	A1875	U1769	A1637	A1535	A1427	U1312
G2742	G2502	G2502	A2347	G2255	C	G	C1996	A1877	G1770	A1637	C1536	C1428	C1315
U2743	U2503	U2503	U2348	C2258	C	A	G2002	C1878	A1773	U1647	A1434	A1434	C1320
G2744	U2504	U2504	C2349	U2259	A	C	C2006	U1880	G1776	U1648	C1435	C1435	C1321
C2745	G2505	G2505	A2350	G2260	C	U	A2014	C1879	C1788	G1649	G1436	A1437	A1322
U2746	U2511	U2511	U2351	A2268	U	U	A2015	U1880	U1796	A1651	G1542	G1542	C1323
G2747	A2513	A2513	G2394	G2281	C	G	G2032	U1881	C1800	G1653	G1543	A1439	G1324
A2748	C2514	C2514	C2395	G2282	A	U	A2031	G1881	A1801	U1667	U1549	A1449	A1327
U2756	C2515	C2515	G2396	C2283	C	G	G2032	C1920	A1802	U1667	G1450	G1450	A1328
A2757	U2401	U2401	U2402	A2287	C	U	C2033	U1916	A1803	G1674	U1554	G1451	U1329
C2758	U2402	U2402	C2403	G2288	U	G	G2034	A1917	U1709	U1683	G1555	G1452	C1330
U2759	C2406	C2406	A2406	G2289	C	U	G2035	G1930	G1715	C1694	C1556	A1453	G1331
A2778	G2409	G2409	G2423	A2291	U	G	C2036	U1931	G1718	U1570	U1458	U1458	G1332
U2779	U2423	U2423	A2424	U2292	A	G	G2037	A1932	G1719	G1479	U1485	U1485	C1333
G2780	A2425	A2425	A2426	U2293	C	U	A2038	G1933	U1720	G1479	C1480	C1480	G1334
C2781	U2426	U2426	G2426	U2294	U	G	U2039	A1934	G1721	U1578	U1481	U1481	G1341
U2788	G2535	G2535	G2429	U2295	C	U	U2043	U1935	U1722	A1579	G1483	G1483	A1342
C2789	U2536	U2536	U2430	G2296	U	G	A2052	U1955	G1723	A1580	G1483	G1483	G1343
U2790	G2544	G2544	U2431	A2297	A	U	C2055	C1958	G1724	G1581	U1484	U1484	U1344
C2791	U2545	U2545	U2431	U2298	G	U	G2056	U1958	U1725	C1582	U1485	U1485	C1345
A2792	C2652	C2652	A2435	U2299	U	U	A2059	U1958	C1726	A1583	U1486	U1486	U1352
				U2202	A	G	A2060	U1958	G1835	U1584	U1486	U1486	U1352
				U2203	A	U	G2061	U1958	U1841	C1585	C1489	C1489	U1372
					G	U	A2062	U1958	G1846	G1587	C1493	C1493	A1373
					U	U	C2066	U1958		A1591	U1379	U1379	U1379
										C1592			



- Molecule 26: 5S rRNA



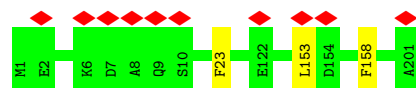
- Molecule 27: Large ribosomal subunit protein uL2



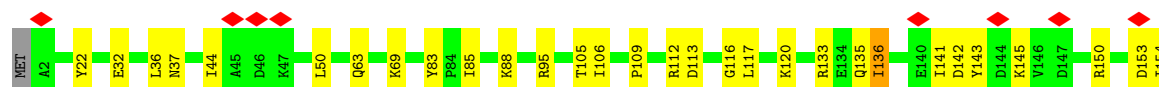
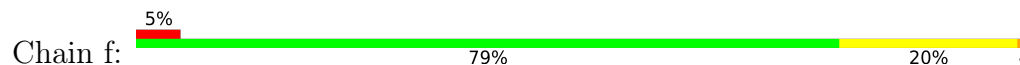
- Molecule 28: Large ribosomal subunit protein uL3



- Molecule 29: Large ribosomal subunit protein uL4



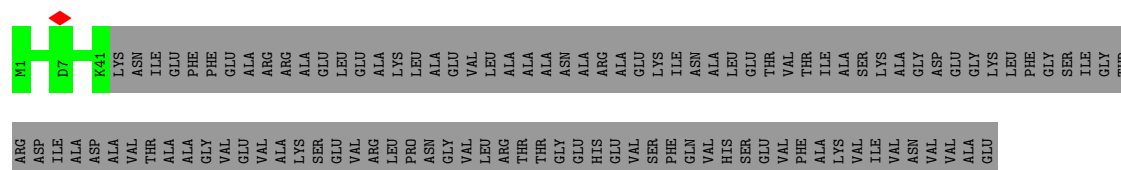
- Molecule 30: Large ribosomal subunit protein uL5



- Molecule 31: Large ribosomal subunit protein uL6



- Molecule 32: Large ribosomal subunit protein bL9



- Molecule 33: Large ribosomal subunit protein uL13



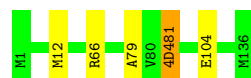
- Molecule 34: Large ribosomal subunit protein uL14



- Molecule 35: Large ribosomal subunit protein uL15



- Molecule 36: Large ribosomal subunit protein uL16



- Molecule 37: Large ribosomal subunit protein bL17





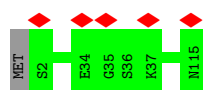
- Molecule 38: Large ribosomal subunit protein uL18

Chain n: 85% 14% ..



- Molecule 39: Large ribosomal subunit protein bL19

Chain o: 99% .



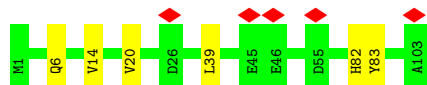
- Molecule 40: Large ribosomal subunit protein bL20

Chain p: 93% 5% ..



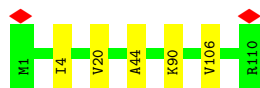
- Molecule 41: Large ribosomal subunit protein bL21

Chain q: 5% 94% 6%



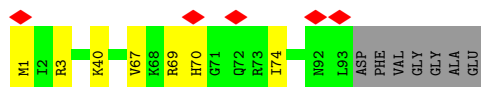
- Molecule 42: Large ribosomal subunit protein uL22

Chain r: 95% 5%

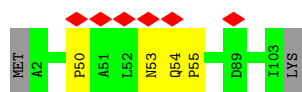
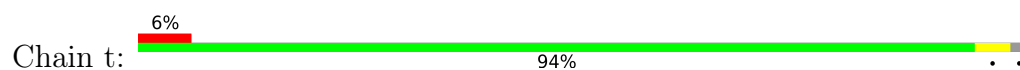


- Molecule 43: Large ribosomal subunit protein uL23

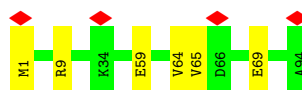
Chain s: 5% 86% 7% 7%



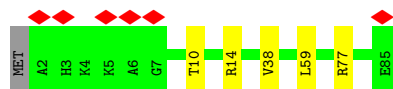
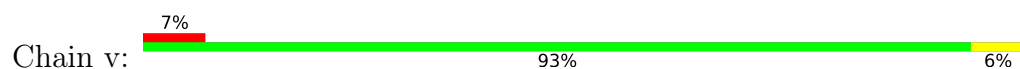
- Molecule 44: Large ribosomal subunit protein uL24



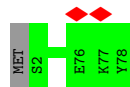
- Molecule 45: Large ribosomal subunit protein bL25



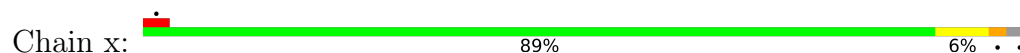
- Molecule 46: Large ribosomal subunit protein bL27



- Molecule 47: Large ribosomal subunit protein bL28



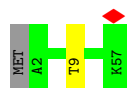
- Molecule 48: Large ribosomal subunit protein uL29




- Molecule 49: Large ribosomal subunit protein uL30

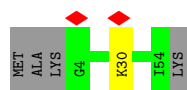


- Molecule 50: Large ribosomal subunit protein bL32



- Molecule 51: Large ribosomal subunit protein bL33

Chain 0:  91% 7%



- Molecule 52: Large ribosomal subunit protein bL34

Chain 1:  96%



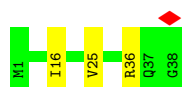
- Molecule 53: Large ribosomal subunit protein bL35

Chain 2:  94% 5%



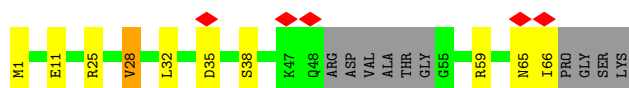
- Molecule 54: Large ribosomal subunit protein bL36A

Chain 3:  92% 8%



- Molecule 55: Large ribosomal subunit protein bL31A

Chain 4:  7% 71% 13% 14%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	196921	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS GLACIOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/\text{\AA}^2$)	28.29	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	1500	Depositor
Magnification	190000	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.586	Depositor
Minimum map value	-0.231	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.021	Depositor
Recommended contour level	0.085	Depositor
Map size (\AA)	447.78, 447.78, 447.78	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	0.8745703, 0.8745703, 0.8745703	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: G7M, MA6, RSP, MG, MS6, IAS, 2MG, UR3, 3TD, T6A, OMC, D2T, SPM, 5MU, OMU, 2MA, SPD, MEQ, 4SU, PSU, ZN, H2U, PUT, 4D4, 1MG, 6MZ, OMG, 4OC, 5MC

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	0.55	1/36261 (0.0%)	0.85	7/56559 (0.0%)
2	B	0.54	0/1784	1.03	2/2403 (0.1%)
3	C	0.52	0/1651	0.92	0/2225
4	D	0.51	0/1665	1.00	0/2227
5	E	0.58	0/1165	1.00	2/1568 (0.1%)
6	F	0.52	0/858	0.91	0/1160
7	G	0.52	0/1219	0.96	0/1635
8	H	0.54	0/989	0.97	0/1326
9	I	0.54	0/1034	0.96	1/1375 (0.1%)
10	J	0.53	0/796	0.96	0/1077
11	K	0.60	0/884	0.97	1/1191 (0.1%)
12	L	0.56	0/960	0.97	0/1286
13	M	0.53	0/900	0.98	0/1204
14	N	0.52	0/817	0.97	0/1088
15	O	0.52	0/722	0.94	0/964
16	P	0.55	0/653	0.90	1/877 (0.1%)
17	Q	0.54	0/650	0.95	0/871
18	R	0.54	0/553	0.93	0/742
19	S	0.55	0/685	0.93	0/922
20	T	0.53	0/676	1.02	0/895
21	U	0.62	1/597 (0.2%)	1.03	1/792 (0.1%)
22	V	0.58	0/1813	0.85	0/2825
22	Z	0.59	0/1813	0.83	0/2825
23	X	0.62	0/270	0.81	0/419
24	Y	0.58	2/2088 (0.1%)	0.79	0/3255
25	a	0.60	2/65697 (0.0%)	0.94	41/102482 (0.0%)
26	b	0.56	0/2850	0.92	1/4444 (0.0%)
27	c	0.64	1/2121 (0.0%)	1.02	0/2852
28	d	0.58	0/1576	0.95	0/2119
29	e	0.57	0/1571	0.97	0/2113
30	f	0.46	0/1434	0.80	0/1926

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
31	g	0.52	0/1315	0.92	0/1783
32	h	0.56	0/306	0.98	0/413
33	i	0.57	0/1152	0.97	0/1551
34	j	0.59	0/955	1.01	0/1279
35	k	0.61	0/1062	0.98	0/1413
36	l	0.58	0/1073	0.96	0/1433
37	m	0.62	1/958 (0.1%)	0.97	0/1281
38	n	0.52	0/902	0.83	0/1209
39	o	0.59	0/929	0.94	0/1242
40	p	0.59	0/960	1.05	3/1278 (0.2%)
41	q	0.56	0/829	0.93	0/1107
42	r	0.59	0/864	0.97	0/1156
43	s	0.52	0/744	0.95	0/994
44	t	0.52	0/787	0.91	0/1051
45	u	0.53	0/766	0.92	0/1025
46	v	0.63	0/636	0.96	0/841
47	w	0.62	0/635	1.01	0/848
48	x	0.44	0/496	0.97	0/660
49	y	0.55	0/453	0.93	0/605
50	z	0.64	0/450	1.05	0/599
51	0	0.56	0/424	0.93	0/565
52	1	0.65	0/380	1.16	2/498 (0.4%)
53	2	0.67	0/513	1.05	0/676
54	3	0.59	0/303	0.94	0/397
55	4	0.53	0/488	0.92	0/649
All	All	0.58	8/155132 (0.0%)	0.92	62/232200 (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
2	B	0	1
4	D	0	2
7	G	0	1
9	I	0	4
10	J	0	1
12	L	0	1
14	N	0	1
27	c	0	1
40	p	0	2

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Mol	Chain	#Chirality outliers	#Planarity outliers
46	v	0	1
48	x	0	3
52	1	0	1
53	2	0	1
54	3	0	1
55	4	0	1
All	All	0	22

The worst 5 of 8 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
21	U	21	ARG	NE-CZ	6.77	1.40	1.33
27	c	53	HIS	CG-CD2	-6.12	1.29	1.35
37	m	3	HIS	CG-CD2	-5.89	1.29	1.35
25	a	2069	G7M	O3'-P	5.54	1.61	1.56
25	a	2552	OMU	O3'-P	5.53	1.61	1.56

The worst 5 of 62 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
25	a	531	C	O3'-P-O5'	11.59	121.38	104.00
21	U	21	ARG	CD-NE-CZ	8.30	136.02	124.40
25	a	504	A	O3'-P-O5'	-7.60	92.60	104.00
25	a	271	G	C2'-C3'-O3'	7.49	120.74	109.50
25	a	557	C	O3'-P-O5'	-6.96	93.55	104.00

There are no chirality outliers.

5 of 22 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	B	137	ARG	Sidechain
4	D	44	ARG	Sidechain
4	D	97	ARG	Sidechain
7	G	10	ARG	Sidechain
9	I	41	ARG	Sidechain

5.2 Too-close contacts

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

the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	32634	0	16443	233	0
2	B	1753	0	1780	20	0
3	C	1624	0	1696	4	0
4	D	1643	0	1707	20	0
5	E	1152	0	1196	1	0
6	F	839	0	833	2	0
7	G	1203	0	1254	6	0
8	H	979	0	1031	1	0
9	I	1022	0	1070	4	0
10	J	786	0	828	6	0
11	K	877	0	884	1	0
12	L	957	0	1017	4	0
13	M	891	0	952	8	0
14	N	805	0	844	2	0
15	O	714	0	734	0	0
16	P	643	0	661	6	0
17	Q	641	0	682	5	0
18	R	544	0	565	0	0
19	S	668	0	693	0	0
20	T	670	0	719	0	0
21	U	589	0	629	3	0
22	V	1623	0	825	21	0
22	Z	1623	0	825	9	0
23	X	240	0	121	4	0
24	Y	2000	0	1020	39	0
25	a	59172	0	29791	237	0
26	b	2549	0	1291	16	0
27	c	2082	0	2154	6	0
28	d	1566	0	1618	5	0
29	e	1552	0	1619	2	0
30	f	1410	0	1444	22	0
31	g	1295	0	1332	0	0
32	h	303	0	327	0	0
33	i	1129	0	1162	2	0
34	j	946	0	1023	3	0
35	k	1053	0	1129	6	0
36	l	1075	0	1145	4	0
37	m	945	0	989	4	0
38	n	892	0	923	12	0
39	o	917	0	962	0	0
40	p	947	0	1019	4	0
41	q	816	0	839	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
42	r	857	0	922	4	0
43	s	738	0	807	6	0
44	t	779	0	831	5	0
45	u	753	0	780	3	0
46	v	628	0	642	2	0
47	w	625	0	652	0	0
48	x	495	0	526	1	0
49	y	449	0	488	0	0
50	z	444	0	458	1	0
51	0	417	0	451	1	0
52	1	377	0	418	1	0
53	2	504	0	572	2	0
54	3	302	0	340	1	0
55	4	480	0	478	10	0
56	A	87	0	0	0	0
56	Y	4	0	0	0	0
56	a	187	0	0	0	0
56	b	5	0	0	0	0
56	c	1	0	0	0	0
56	d	2	0	0	0	0
56	m	1	0	0	0	0
56	z	1	0	0	0	0
57	A	10	0	19	0	0
57	a	60	0	114	3	0
58	A	6	0	12	0	0
58	a	36	0	72	0	0
59	a	14	0	26	2	0
60	3	1	0	0	0	0
60	4	1	0	0	0	0
All	All	144033	0	96384	680	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 3.

The worst 5 of 680 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)
1:A:84:U:H3'	1:A:86:G:H21	1.34	0.90
25:a:276:U:H2'	25:a:277:G:C8	2.19	0.77
25:a:284:U:H2'	25:a:285:G:H5''	1.67	0.76
30:f:158:THR:HG22	30:f:160:ALA:H	1.49	0.76

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
25:a:280:U:H2'	25:a:281:C:C6	2.20	0.76

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	B	222/241 (92%)	208 (94%)	14 (6%)	0	100	100
3	C	204/233 (88%)	199 (98%)	5 (2%)	0	100	100
4	D	203/206 (98%)	194 (96%)	9 (4%)	0	100	100
5	E	154/167 (92%)	149 (97%)	5 (3%)	0	100	100
6	F	101/135 (75%)	97 (96%)	4 (4%)	0	100	100
7	G	151/179 (84%)	147 (97%)	4 (3%)	0	100	100
8	H	127/130 (98%)	119 (94%)	8 (6%)	0	100	100
9	I	125/130 (96%)	119 (95%)	6 (5%)	0	100	100
10	J	96/103 (93%)	94 (98%)	1 (1%)	1 (1%)	13	26
11	K	113/129 (88%)	110 (97%)	3 (3%)	0	100	100
12	L	120/124 (97%)	118 (98%)	2 (2%)	0	100	100
13	M	113/118 (96%)	111 (98%)	2 (2%)	0	100	100
14	N	98/101 (97%)	98 (100%)	0	0	100	100
15	O	86/89 (97%)	85 (99%)	1 (1%)	0	100	100
16	P	79/82 (96%)	78 (99%)	1 (1%)	0	100	100
17	Q	77/84 (92%)	76 (99%)	1 (1%)	0	100	100
18	R	64/75 (85%)	63 (98%)	1 (2%)	0	100	100
19	S	82/92 (89%)	81 (99%)	1 (1%)	0	100	100
20	T	84/87 (97%)	84 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
21	U	68/71 (96%)	66 (97%)	2 (3%)	0	100	100
27	c	269/273 (98%)	262 (97%)	7 (3%)	0	100	100
28	d	206/209 (99%)	200 (97%)	6 (3%)	0	100	100
29	e	199/201 (99%)	194 (98%)	5 (2%)	0	100	100
30	f	175/179 (98%)	171 (98%)	4 (2%)	0	100	100
31	g	171/177 (97%)	165 (96%)	6 (4%)	0	100	100
32	h	39/149 (26%)	36 (92%)	3 (8%)	0	100	100
33	i	140/142 (99%)	140 (100%)	0	0	100	100
34	j	121/123 (98%)	118 (98%)	3 (2%)	0	100	100
35	k	142/144 (99%)	139 (98%)	3 (2%)	0	100	100
36	l	132/136 (97%)	128 (97%)	4 (3%)	0	100	100
37	m	116/127 (91%)	113 (97%)	3 (3%)	0	100	100
38	n	114/117 (97%)	112 (98%)	2 (2%)	0	100	100
39	o	112/115 (97%)	108 (96%)	4 (4%)	0	100	100
40	p	115/118 (98%)	114 (99%)	1 (1%)	0	100	100
41	q	101/103 (98%)	99 (98%)	2 (2%)	0	100	100
42	r	108/110 (98%)	107 (99%)	1 (1%)	0	100	100
43	s	91/100 (91%)	88 (97%)	3 (3%)	0	100	100
44	t	100/104 (96%)	96 (96%)	4 (4%)	0	100	100
45	u	92/94 (98%)	90 (98%)	2 (2%)	0	100	100
46	v	82/85 (96%)	78 (95%)	4 (5%)	0	100	100
47	w	75/78 (96%)	74 (99%)	1 (1%)	0	100	100
48	x	59/63 (94%)	59 (100%)	0	0	100	100
49	y	56/59 (95%)	55 (98%)	1 (2%)	0	100	100
50	z	54/57 (95%)	53 (98%)	1 (2%)	0	100	100
51	0	49/55 (89%)	49 (100%)	0	0	100	100
52	1	44/46 (96%)	44 (100%)	0	0	100	100
53	2	62/65 (95%)	60 (97%)	2 (3%)	0	100	100
54	3	36/38 (95%)	36 (100%)	0	0	100	100
55	4	56/70 (80%)	51 (91%)	5 (9%)	0	100	100
All	All	5483/5913 (93%)	5335 (97%)	147 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
10	J	57	VAL

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	B	186/199 (94%)	183 (98%)	3 (2%)	58	78
3	C	170/190 (90%)	170 (100%)	0	100	100
4	D	172/173 (99%)	170 (99%)	2 (1%)	67	84
5	E	119/126 (94%)	119 (100%)	0	100	100
6	F	90/116 (78%)	90 (100%)	0	100	100
7	G	126/147 (86%)	125 (99%)	1 (1%)	79	91
8	H	104/105 (99%)	104 (100%)	0	100	100
9	I	105/107 (98%)	104 (99%)	1 (1%)	73	87
10	J	86/90 (96%)	86 (100%)	0	100	100
11	K	89/98 (91%)	89 (100%)	0	100	100
12	L	102/103 (99%)	100 (98%)	2 (2%)	50	73
13	M	93/96 (97%)	93 (100%)	0	100	100
14	N	83/84 (99%)	83 (100%)	0	100	100
15	O	76/77 (99%)	76 (100%)	0	100	100
16	P	65/65 (100%)	64 (98%)	1 (2%)	60	80
17	Q	73/78 (94%)	73 (100%)	0	100	100
18	R	57/65 (88%)	56 (98%)	1 (2%)	54	76
19	S	72/79 (91%)	72 (100%)	0	100	100
20	T	65/66 (98%)	65 (100%)	0	100	100
21	U	60/61 (98%)	58 (97%)	2 (3%)	33	57
27	c	216/218 (99%)	215 (100%)	1 (0%)	86	95
28	d	163/163 (100%)	161 (99%)	2 (1%)	67	84

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
29	e	165/165 (100%)	165 (100%)	0	100	100
30	f	148/150 (99%)	143 (97%)	5 (3%)	32	56
31	g	134/138 (97%)	132 (98%)	2 (2%)	60	80
32	h	32/114 (28%)	32 (100%)	0	100	100
33	i	116/116 (100%)	114 (98%)	2 (2%)	56	77
34	j	104/104 (100%)	104 (100%)	0	100	100
35	k	103/103 (100%)	103 (100%)	0	100	100
36	l	107/107 (100%)	107 (100%)	0	100	100
37	m	98/103 (95%)	98 (100%)	0	100	100
38	n	86/87 (99%)	82 (95%)	4 (5%)	22	43
39	o	99/100 (99%)	99 (100%)	0	100	100
40	p	89/90 (99%)	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	80/84 (95%)	80 (100%)	0	100	100
44	t	83/85 (98%)	83 (100%)	0	100	100
45	u	78/78 (100%)	77 (99%)	1 (1%)	65	83
46	v	61/63 (97%)	60 (98%)	1 (2%)	58	78
47	w	67/68 (98%)	67 (100%)	0	100	100
48	x	54/55 (98%)	53 (98%)	1 (2%)	52	75
49	y	48/49 (98%)	48 (100%)	0	100	100
50	z	47/48 (98%)	47 (100%)	0	100	100
51	0	46/49 (94%)	46 (100%)	0	100	100
52	1	38/38 (100%)	38 (100%)	0	100	100
53	2	51/52 (98%)	51 (100%)	0	100	100
54	3	34/34 (100%)	34 (100%)	0	100	100
55	4	55/62 (89%)	54 (98%)	1 (2%)	54	76
All	All	4572/4825 (95%)	4539 (99%)	33 (1%)	80	92

5 of 33 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
38	n	87	ILE

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Mol	Chain	Res	Type
45	u	65	VAL
55	4	28	VAL
21	U	63	GLU
21	U	60	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 61 such sidechains are listed below:

Mol	Chain	Res	Type
15	O	35	GLN
48	x	15	ASN
20	T	61	GLN
44	t	53	ASN
55	4	20	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	1517/1542 (98%)	203 (13%)	27 (1%)
22	V	75/76 (98%)	35 (46%)	3 (4%)
22	Z	75/76 (98%)	27 (36%)	4 (5%)
23	X	10/24 (41%)	3 (30%)	1 (10%)
24	Y	92/93 (98%)	16 (17%)	1 (1%)
25	a	2747/2904 (94%)	689 (25%)	0
26	b	118/120 (98%)	32 (27%)	0
All	All	4634/4835 (95%)	1005 (21%)	36 (0%)

5 of 1005 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	3	A
1	A	4	U
1	A	5	U
1	A	6	G
1	A	7	A

5 of 36 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
22	V	17(A)	U
22	Z	48	C

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Mol	Chain	Res	Type
22	V	18	G
22	Z	7	G
1	A	776	G

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

45 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$
24	PSU	Y	55	24	18,21,22	0.88	1 (5%)	22,30,33	0.70	0
28	MEQ	d	150	28	8,9,10	0.50	0	5,10,12	0.60	0
25	2MA	a	2503	56,25	19,25,26	1.01	2 (10%)	21,37,40	1.82	4 (19%)
25	5MU	a	1939	25	19,22,23	0.48	0	28,32,35	0.63	0
25	PSU	a	2604	25	18,21,22	1.00	1 (5%)	22,30,33	1.45	4 (18%)
36	4D4	l	81	36	9,11,12	0.49	0	8,13,15	0.79	1 (12%)
1	PSU	A	516	1,56	18,21,22	0.86	1 (5%)	22,30,33	0.95	1 (4%)
25	OMG	a	2251	22,25	18,26,27	0.96	2 (11%)	19,38,41	0.77	1 (5%)
25	H2U	a	2449	25	18,21,22	0.79	1 (5%)	21,30,33	0.87	1 (4%)
25	PSU	a	746	56,25	18,21,22	1.13	2 (11%)	22,30,33	0.89	0
25	PSU	a	2580	25	18,21,22	0.95	1 (5%)	22,30,33	0.83	1 (4%)
1	2MG	A	966	1	18,26,27	1.07	2 (11%)	16,38,41	0.78	0
1	UR3	A	1498	1	19,22,23	0.40	0	26,32,35	0.68	0
11	IAS	K	119	11	6,7,8	0.95	0	6,8,10	0.95	0
1	MA6	A	1519	1	18,26,27	0.78	1 (5%)	19,38,41	0.79	1 (5%)
25	PSU	a	955	25	18,21,22	0.84	1 (5%)	22,30,33	0.85	0
25	G7M	a	2069	25	20,26,27	1.44	3 (15%)	17,39,42	0.76	0
1	MA6	A	1518	1	18,26,27	0.83	1 (5%)	19,38,41	0.47	0
12	D2T	L	89	12	7,9,10	1.01	0	6,11,13	1.92	4 (66%)
1	5MC	A	1407	1	18,22,23	0.42	0	26,32,35	0.76	1 (3%)
24	T6A	Y	37	24	27,34,35	0.80	0	29,49,52	0.79	1 (3%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	2MG	A	1207	1	18,26,27	1.01	1 (5%)	16,38,41	0.76	0
24	5MU	Y	54	24	19,22,23	0.37	0	28,32,35	0.33	0
25	2MG	a	2445	25	18,26,27	1.24	3 (16%)	16,38,41	0.68	0
1	4OC	A	1402	1,56	20,23,24	0.37	0	26,32,35	0.57	0
1	5MC	A	967	1	18,22,23	0.29	0	26,32,35	0.51	0
25	1MG	a	745	25	18,26,27	1.07	2 (11%)	19,39,42	0.95	1 (5%)
1	2MG	A	1516	1	18,26,27	1.08	3 (16%)	16,38,41	0.79	0
24	4SU	Y	8	24	18,21,22	0.23	0	26,30,33	0.29	0
25	PSU	a	2457	25	18,21,22	1.01	1 (5%)	22,30,33	0.69	0
25	6MZ	a	1618	25	18,25,26	0.95	0	16,36,39	0.69	0
24	H2U	Y	20(A)	24	18,21,22	0.62	0	21,30,33	0.82	1 (4%)
25	5MU	a	747	25	19,22,23	0.39	0	28,32,35	0.47	0
25	3TD	a	1915	25	18,22,23	0.92	1 (5%)	22,32,35	0.65	0
25	PSU	a	1917	25	18,21,22	0.86	1 (5%)	22,30,33	0.85	0
24	RSP	Y	32	24	17,21,22	0.26	0	22,30,33	0.33	0
1	G7M	A	527	1	20,26,27	1.12	2 (10%)	17,39,42	0.50	0
25	PSU	a	1911	25	18,21,22	0.86	1 (5%)	22,30,33	0.71	0
25	5MC	a	1962	25	18,22,23	0.41	0	26,32,35	0.45	0
25	2MG	a	1835	25	18,26,27	1.00	1 (5%)	16,38,41	0.87	0
25	6MZ	a	2030	25	18,25,26	0.72	0	16,36,39	0.77	0
25	PSU	a	2504	25	18,21,22	0.98	1 (5%)	22,30,33	0.87	1 (4%)
25	OMU	a	2552	25	19,22,23	0.49	0	26,31,34	0.68	1 (3%)
25	PSU	a	2605	25	18,21,22	1.03	1 (5%)	22,30,33	0.80	0
25	OMC	a	2498	56,25	19,22,23	0.62	0	26,31,34	0.52	0

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
24	PSU	Y	55	24	-	0/7/25/26	0/2/2/2
28	MEQ	d	150	28	-	2/8/9/11	-
25	2MA	a	2503	56,25	-	2/3/25/26	0/3/3/3
25	5MU	a	1939	25	-	0/7/25/26	0/2/2/2
25	PSU	a	2604	25	-	0/7/25/26	0/2/2/2
36	4D4	l	81	36	-	0/11/12/14	-
1	PSU	A	516	1,56	-	0/7/25/26	0/2/2/2
25	OMG	a	2251	22,25	-	1/5/27/28	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
25	H2U	a	2449	25	-	3/7/38/39	0/2/2/2
25	PSU	a	746	56,25	-	1/7/25/26	0/2/2/2
25	PSU	a	2580	25	-	0/7/25/26	0/2/2/2
1	2MG	A	966	1	-	0/5/27/28	0/3/3/3
1	UR3	A	1498	1	-	0/7/25/26	0/2/2/2
11	IAS	K	119	11	-	1/7/7/8	-
1	MA6	A	1519	1	-	0/7/29/30	0/3/3/3
25	PSU	a	955	25	-	0/7/25/26	0/2/2/2
25	G7M	a	2069	25	-	1/3/25/26	0/3/3/3
1	MA6	A	1518	1	-	0/7/29/30	0/3/3/3
12	D2T	L	89	12	-	4/7/12/14	-
1	5MC	A	1407	1	-	0/7/25/26	0/2/2/2
24	T6A	Y	37	24	-	6/19/41/42	0/3/3/3
1	2MG	A	1207	1	-	0/5/27/28	0/3/3/3
24	5MU	Y	54	24	-	0/7/25/26	0/2/2/2
25	2MG	a	2445	25	-	1/5/27/28	0/3/3/3
1	4OC	A	1402	1,56	-	1/9/29/30	0/2/2/2
1	5MC	A	967	1	-	0/7/25/26	0/2/2/2
25	1MG	a	745	25	-	0/3/25/26	0/3/3/3
1	2MG	A	1516	1	-	0/5/27/28	0/3/3/3
24	4SU	Y	8	24	-	0/7/25/26	0/2/2/2
25	PSU	a	2457	25	-	0/7/25/26	0/2/2/2
25	6MZ	a	1618	25	-	2/5/27/28	0/3/3/3
24	H2U	Y	20(A)	24	-	3/7/38/39	0/2/2/2
25	5MU	a	747	25	-	1/7/25/26	0/2/2/2
25	3TD	a	1915	25	-	2/7/25/26	0/2/2/2
25	PSU	a	1917	25	-	0/7/25/26	0/2/2/2
24	RSP	Y	32	24	-	2/7/25/26	0/2/2/2
1	G7M	A	527	1	-	1/3/25/26	0/3/3/3
25	PSU	a	1911	25	-	0/7/25/26	0/2/2/2
25	5MC	a	1962	25	-	0/7/25/26	0/2/2/2
25	2MG	a	1835	25	-	0/5/27/28	0/3/3/3
25	6MZ	a	2030	25	-	2/5/27/28	0/3/3/3
25	PSU	a	2504	25	-	2/7/25/26	0/2/2/2
25	OMU	a	2552	25	-	0/9/27/28	0/2/2/2
25	PSU	a	2605	25	-	0/7/25/26	0/2/2/2
25	OMC	a	2498	56,25	-	0/9/27/28	0/2/2/2

The worst 5 of 37 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
25	a	1915	3TD	C6-C5	3.62	1.39	1.35
25	a	2069	G7M	C8-N9	3.52	1.39	1.33
1	A	527	G7M	C8-N9	3.50	1.39	1.33
25	a	2605	PSU	C6-C5	3.42	1.39	1.35
25	a	2504	PSU	C6-C5	3.38	1.39	1.35

The worst 5 of 24 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
25	a	2503	2MA	C5-C6-N1	-5.47	117.42	121.01
25	a	2503	2MA	C5-C6-N6	4.55	127.26	120.35
25	a	2604	PSU	C6-C5-C4	3.62	120.73	118.20
25	a	2604	PSU	C5-C6-N1	-3.23	117.27	122.11
25	a	2503	2MA	CM2-C2-N1	2.70	121.36	117.15

There are no chirality outliers.

5 of 38 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	1402	4OC	C1'-C2'-O2'-CM2
12	L	89	D2T	CA-CB-CG-OD1
12	L	89	D2T	CA-CB-CG-OD2
24	Y	32	RSP	C3'-C4'-C5'-O5'
24	Y	32	RSP	O4'-C4'-C5'-O5'

There are no ring outliers.

10 monomers are involved in 11 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
36	l	81	4D4	1	0
25	a	2251	OMG	1	0
1	A	1519	MA6	2	0
1	A	1518	MA6	2	0
24	Y	37	T6A	1	0
1	A	1516	2MG	1	0
24	Y	20(A)	H2U	1	0
25	a	747	5MU	1	0
25	a	1915	3TD	1	0
25	a	2030	6MZ	2	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 305 ligands modelled in this entry, 290 are monoatomic - leaving 15 for Mogul analysis.

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	$\# Z > 2$
58	PUT	A	1686	-	5,5,5	0.15	0	4,4,4	0.12	0
57	SPD	a	6192	-	9,9,9	0.17	0	8,8,8	0.32	0
59	SPM	a	6193	-	13,13,13	0.17	0	12,12,12	0.31	0
57	SPD	A	1685	-	9,9,9	0.20	0	8,8,8	0.21	0
58	PUT	a	6194	-	5,5,5	0.19	0	4,4,4	0.32	0
57	SPD	a	6189	-	9,9,9	0.20	0	8,8,8	0.35	0
58	PUT	a	6197	-	5,5,5	0.18	0	4,4,4	0.30	0
58	PUT	a	6196	-	5,5,5	0.16	0	4,4,4	0.41	0
57	SPD	a	6188	-	9,9,9	0.17	0	8,8,8	0.40	0
57	SPD	a	6187	-	9,9,9	0.25	0	8,8,8	0.27	0
57	SPD	a	6191	-	9,9,9	0.20	0	8,8,8	0.22	0
58	PUT	a	6198	-	5,5,5	0.20	0	4,4,4	0.35	0
58	PUT	a	6200	-	5,5,5	0.13	0	4,4,4	0.23	0
57	SPD	a	6190	-	9,9,9	0.19	0	8,8,8	0.20	0
58	PUT	a	6199	-	5,5,5	0.19	0	4,4,4	0.25	0

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
58	PUT	A	1686	-	-	0/3/3/3	-
57	SPD	a	6192	-	-	0/7/7/7	-
59	SPM	a	6193	-	-	3/11/11/11	-
57	SPD	A	1685	-	-	0/7/7/7	-
58	PUT	a	6194	-	-	0/3/3/3	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
57	SPD	a	6189	-	-	1/7/7/7	-
58	PUT	a	6197	-	-	0/3/3/3	-
58	PUT	a	6196	-	-	0/3/3/3	-
57	SPD	a	6188	-	-	0/7/7/7	-
57	SPD	a	6187	-	-	1/7/7/7	-
57	SPD	a	6191	-	-	0/7/7/7	-
58	PUT	a	6198	-	-	1/3/3/3	-
58	PUT	a	6200	-	-	1/3/3/3	-
57	SPD	a	6190	-	-	0/7/7/7	-
58	PUT	a	6199	-	-	0/3/3/3	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
59	a	6193	SPM	C7-C6-N5-C4
59	a	6193	SPM	N10-C11-C12-C13
58	a	6198	PUT	C1-C2-C3-C4
58	a	6200	PUT	C1-C2-C3-C4
57	a	6187	SPD	C2-C3-C4-C5

There are no ring outliers.

4 monomers are involved in 5 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
57	a	6192	SPD	1	0
59	a	6193	SPM	2	0
57	a	6191	SPD	1	0
57	a	6190	SPD	1	0

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

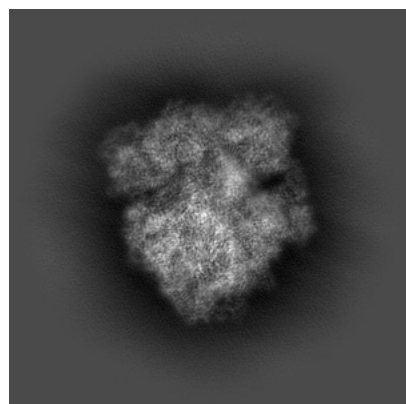
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-51679. These allow visual inspection of the internal detail of the map and identification of artifacts.

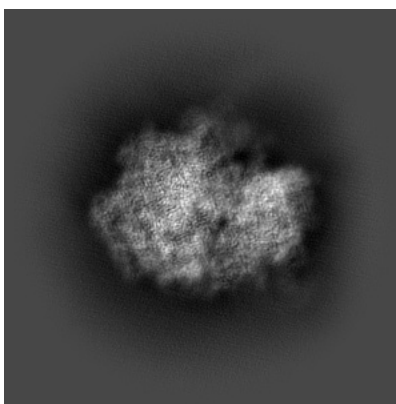
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

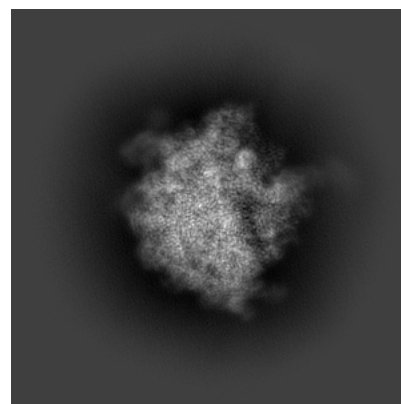
6.1.1 Primary map



X

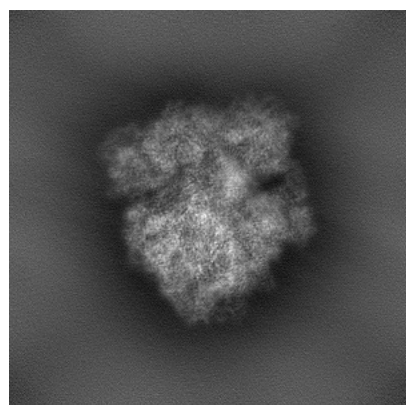


Y

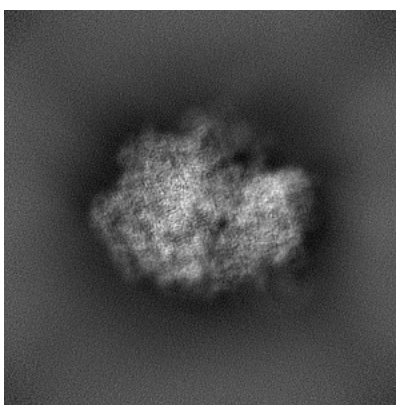


Z

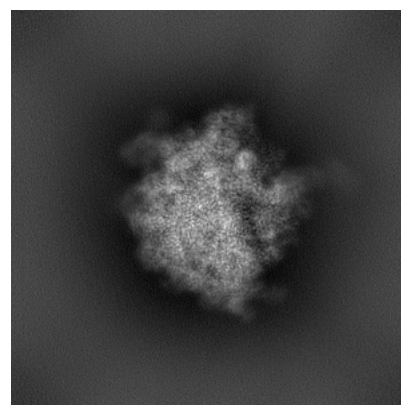
6.1.2 Raw map



X



Y

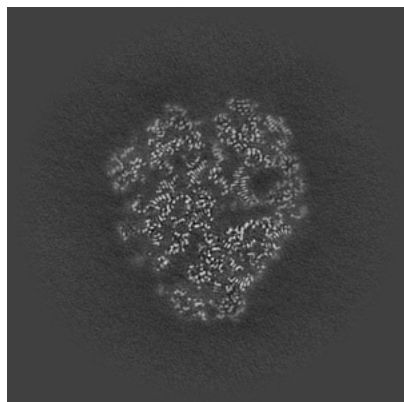


Z

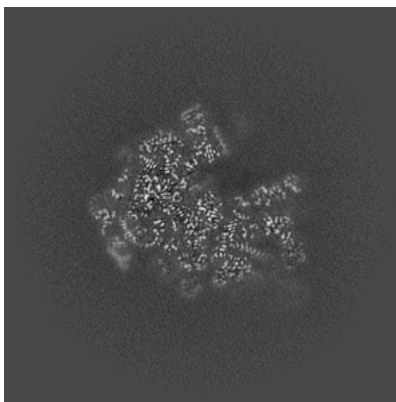
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

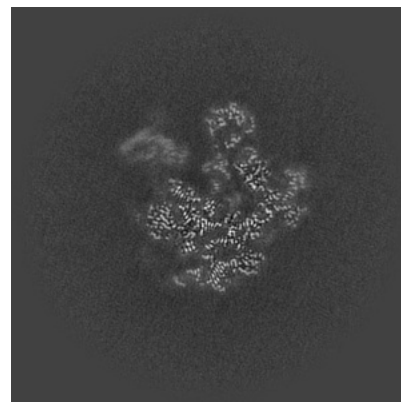
6.2.1 Primary map



X Index: 256

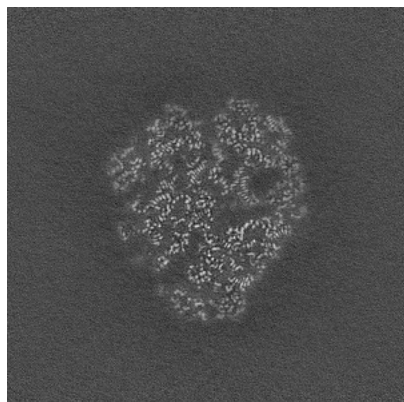


Y Index: 256

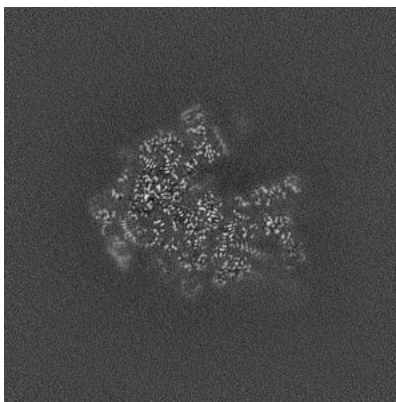


Z Index: 256

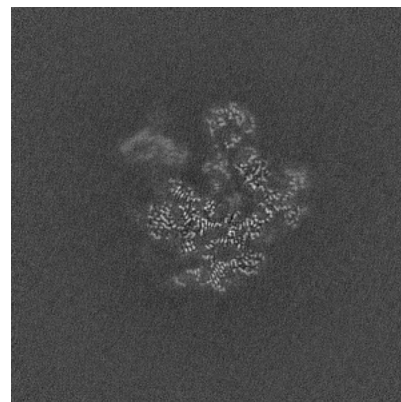
6.2.2 Raw map



X Index: 256



Y Index: 256

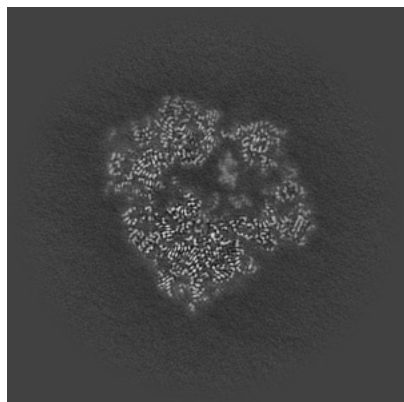


Z Index: 256

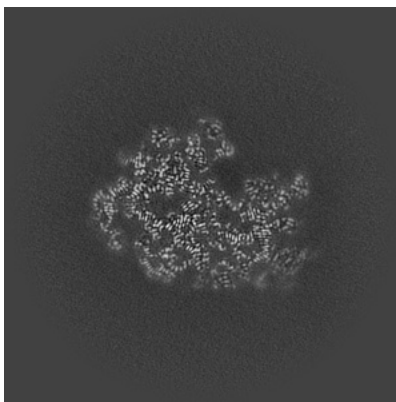
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

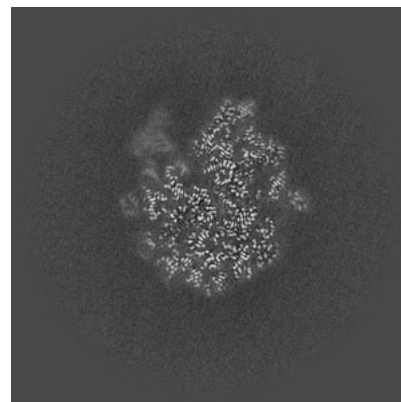
6.3.1 Primary map



X Index: 278

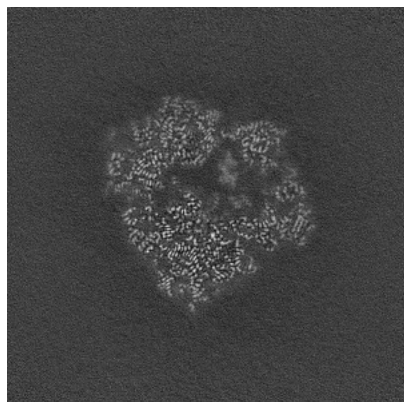


Y Index: 241

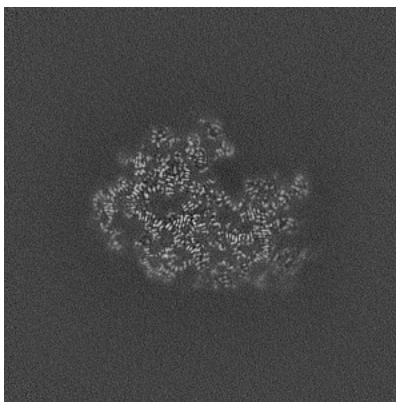


Z Index: 232

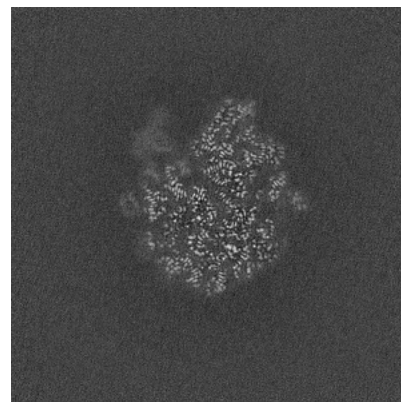
6.3.2 Raw map



X Index: 278



Y Index: 241

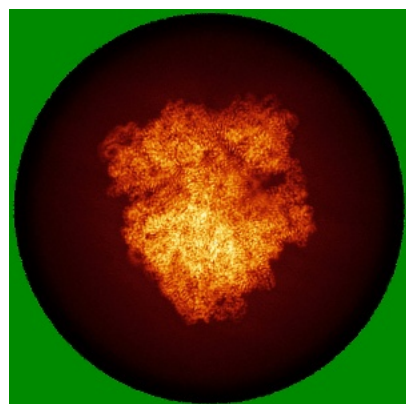


Z Index: 231

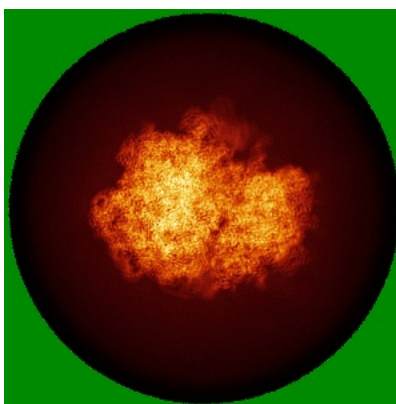
The images above show the largest variance slices of the map in three orthogonal directions.

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

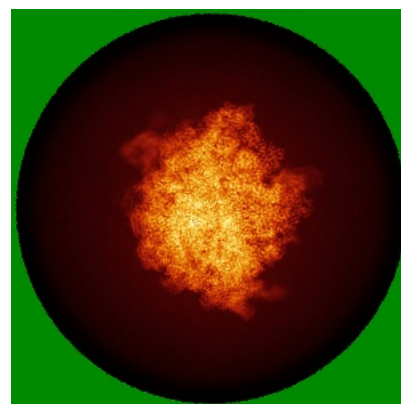
6.4.1 Primary map



X



Y

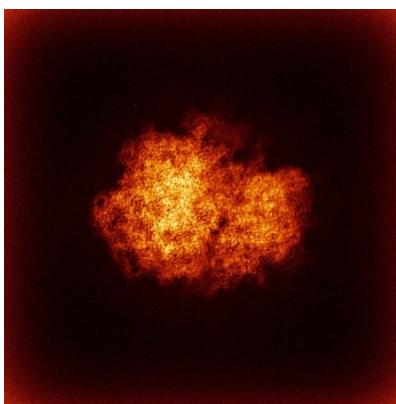


Z

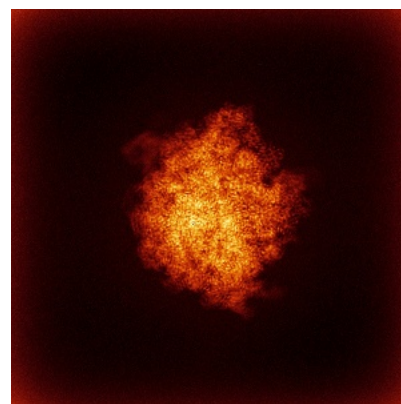
6.4.2 Raw map



X



Y

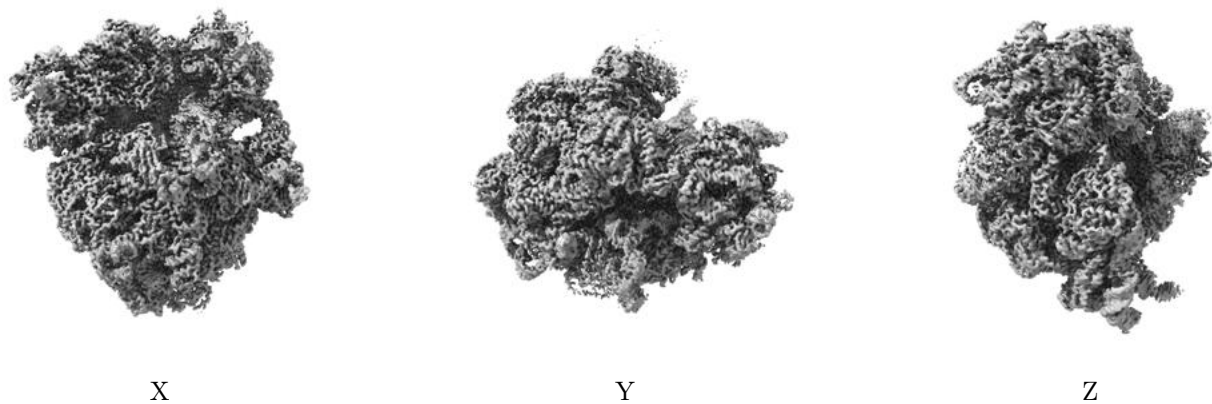


Z

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

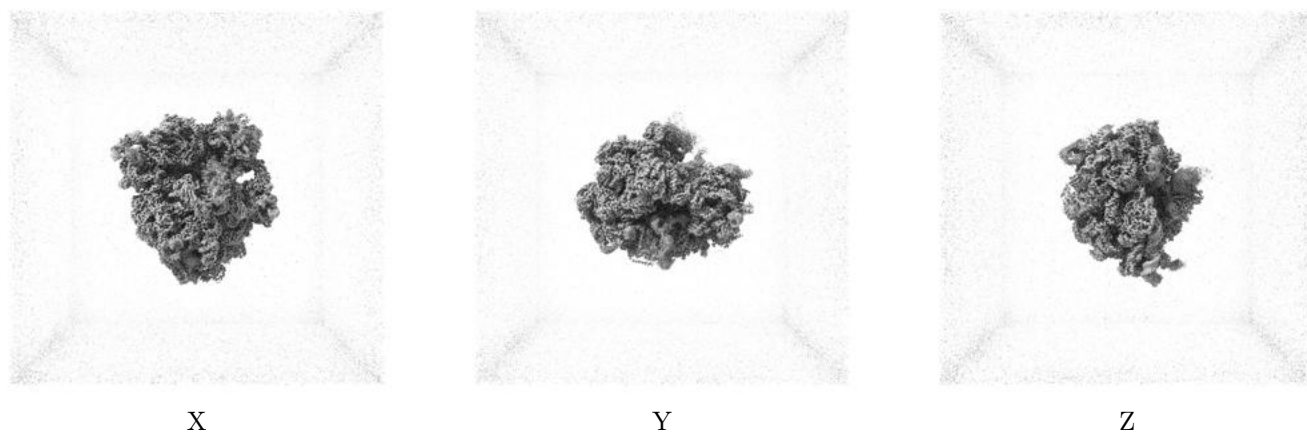
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.085. 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.

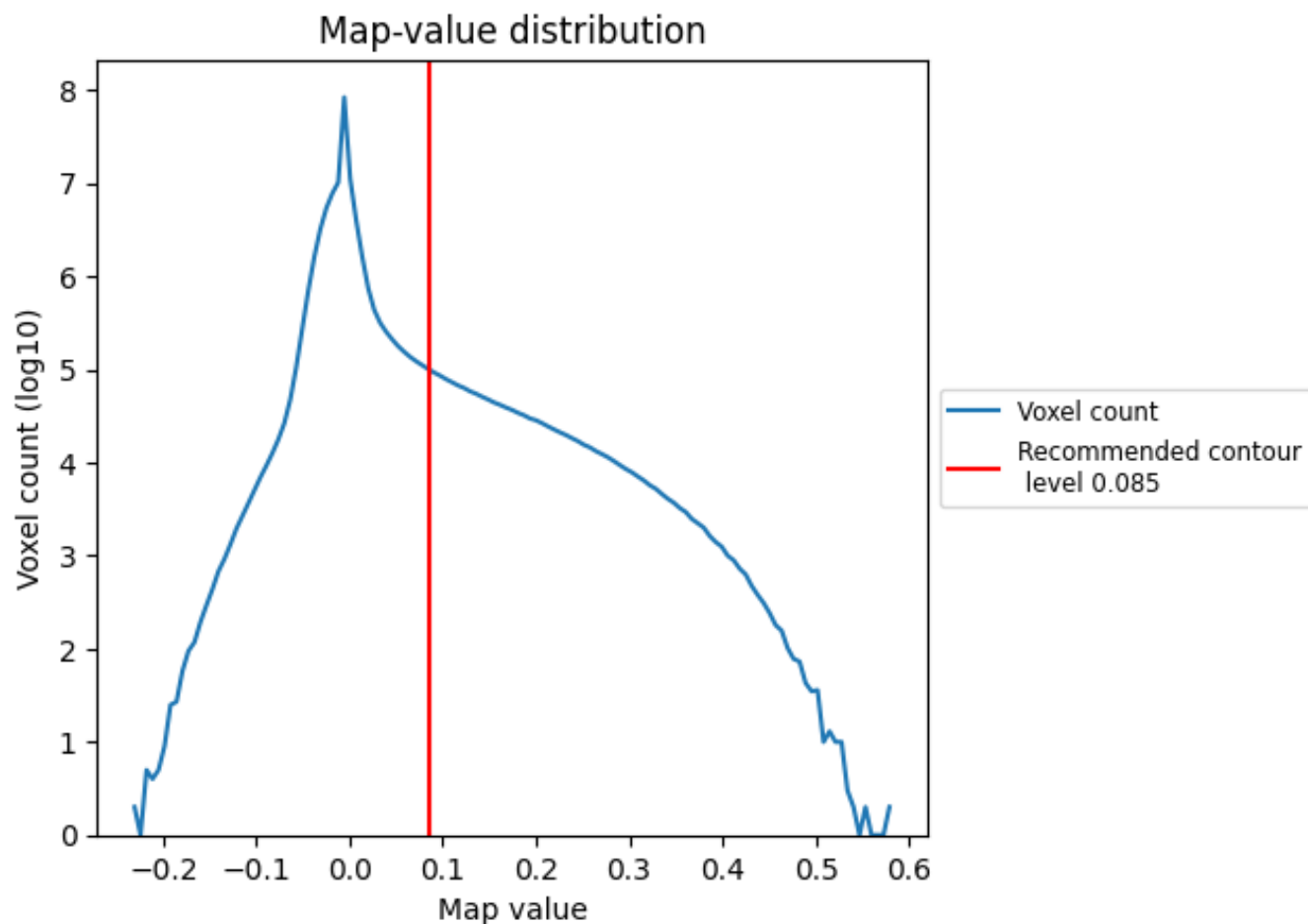
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

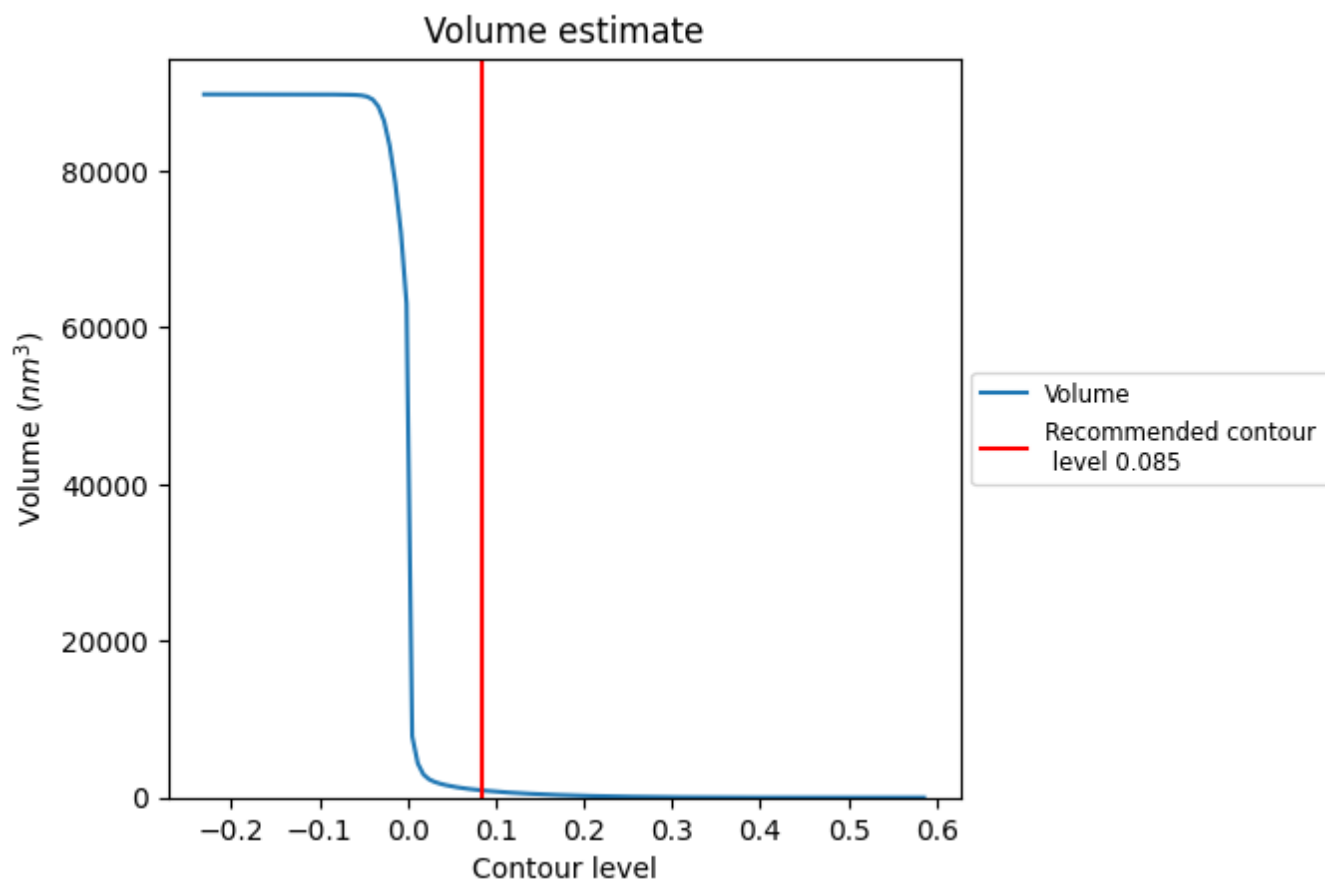
This section contains the results of statistical analysis of the map.

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.

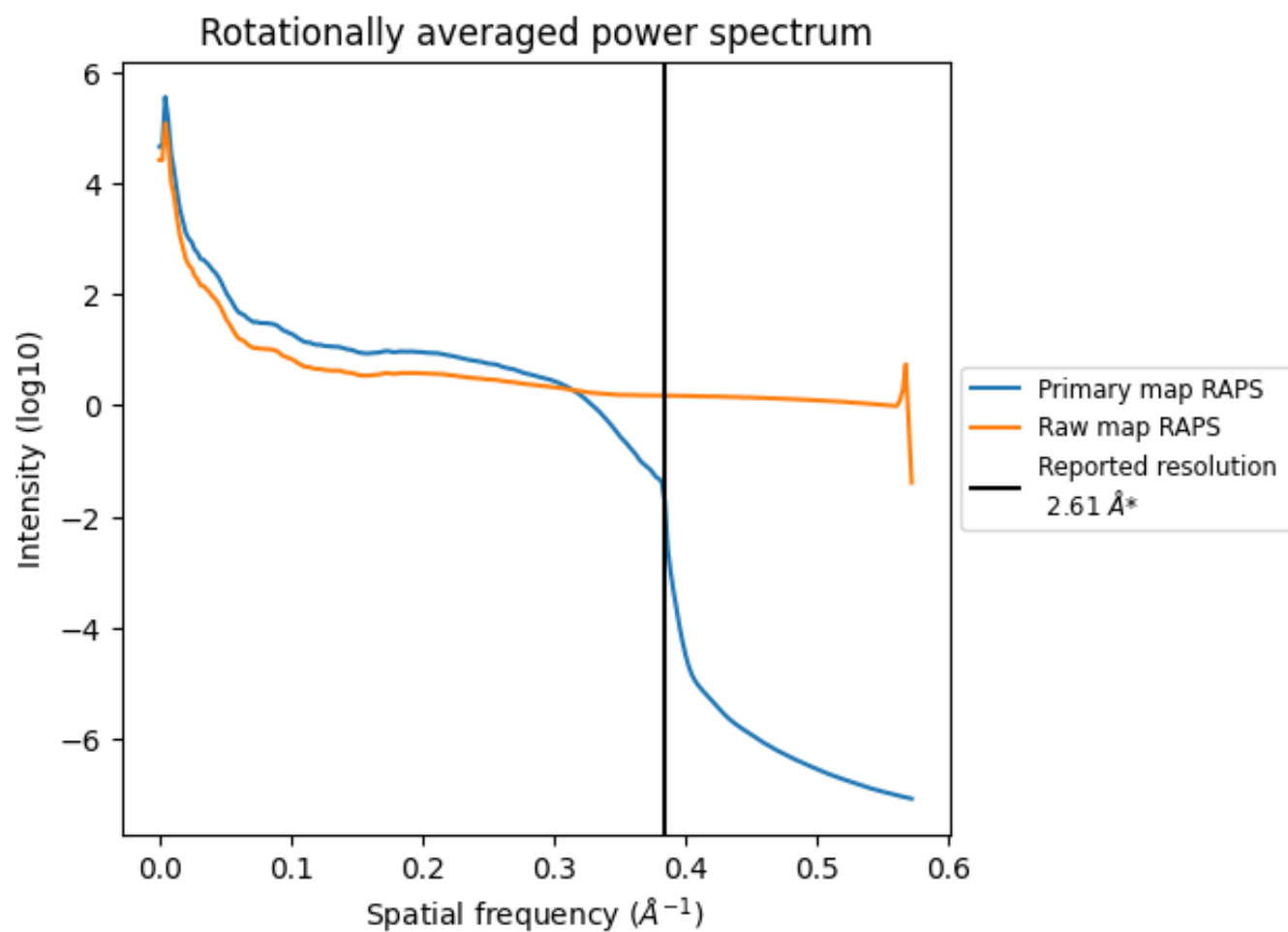
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 912 nm³; this corresponds to an approximate mass of 824 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 ⓘ

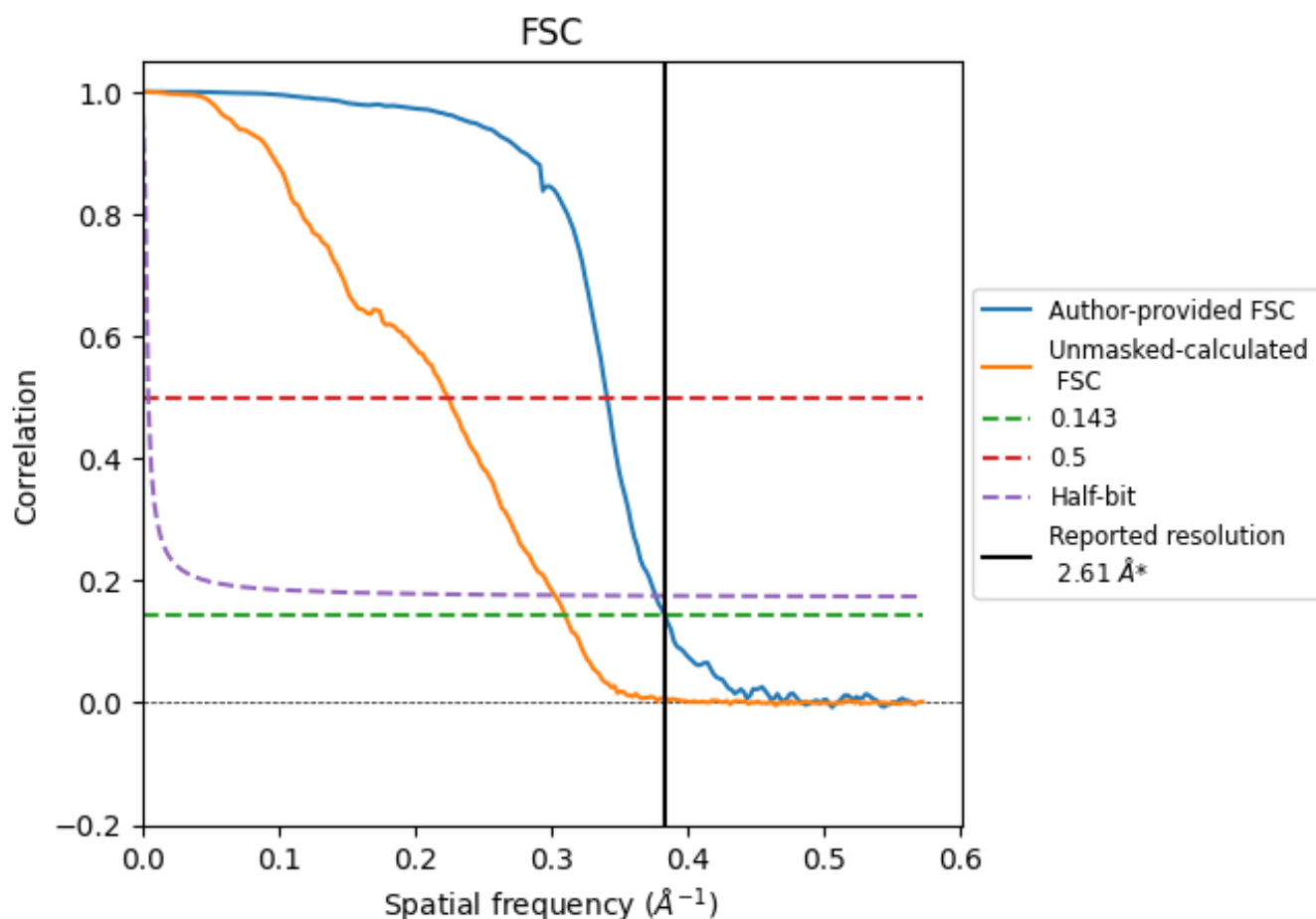


*Reported resolution corresponds to spatial frequency of 0.383 \AA^{-1}

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 0.383 Å⁻¹

8.2 Resolution estimates [i](#)

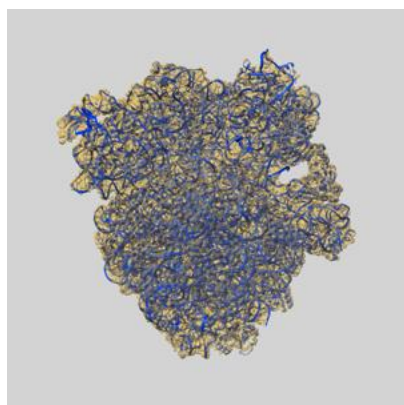
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.61	-	-
Author-provided FSC curve	2.61	2.94	2.66
Unmasked-calculated*	3.23	4.46	3.31

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.23 differs from the reported value 2.61 by more than 10 %

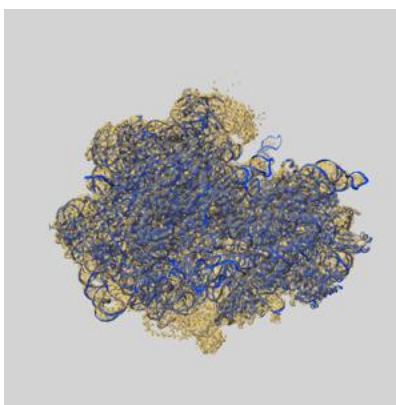
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-51679 and PDB model 9GXX. Per-residue inclusion information can be found in section [3](#) on page [17](#).

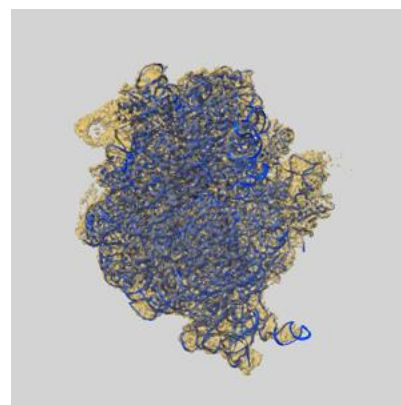
9.1 Map-model overlay [i](#)



X



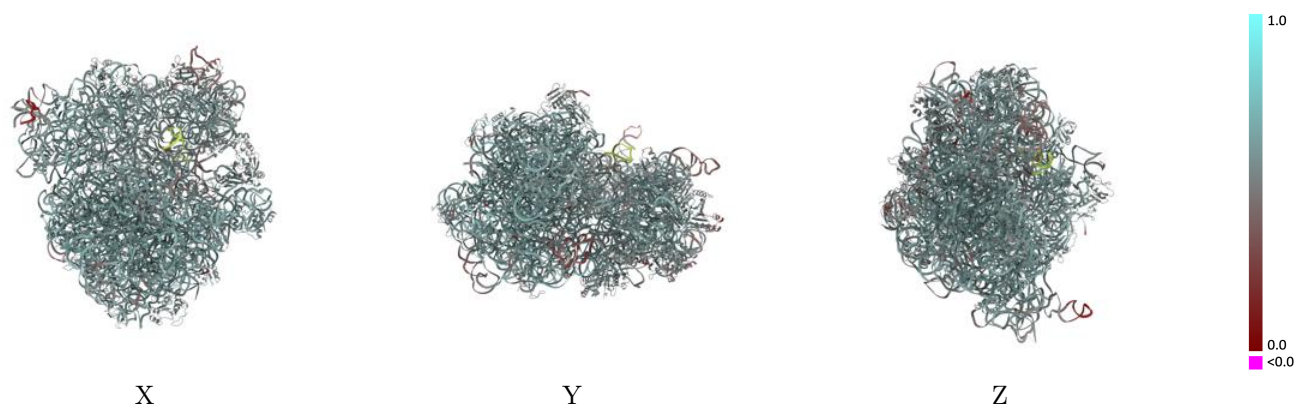
Y



Z

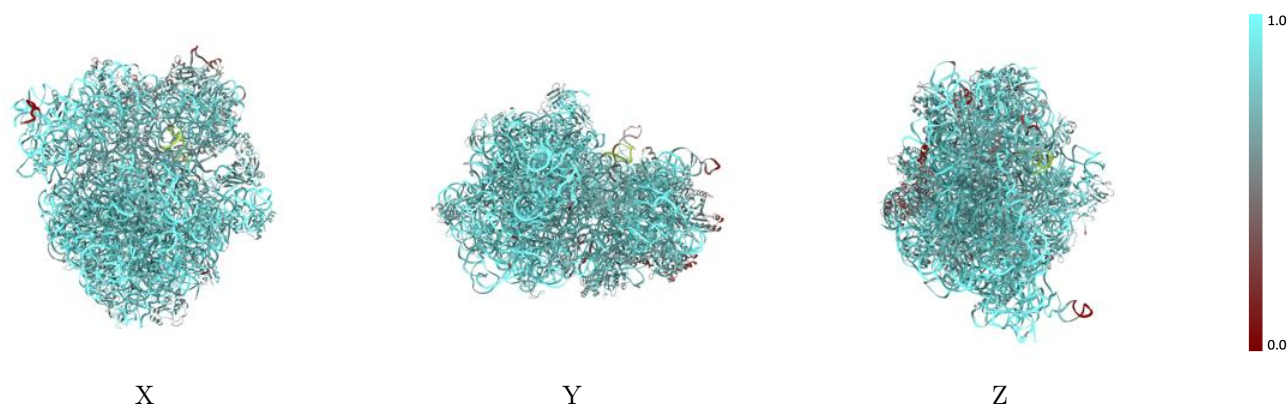
The images above show the 3D surface view of the map at the recommended contour level 0.085 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



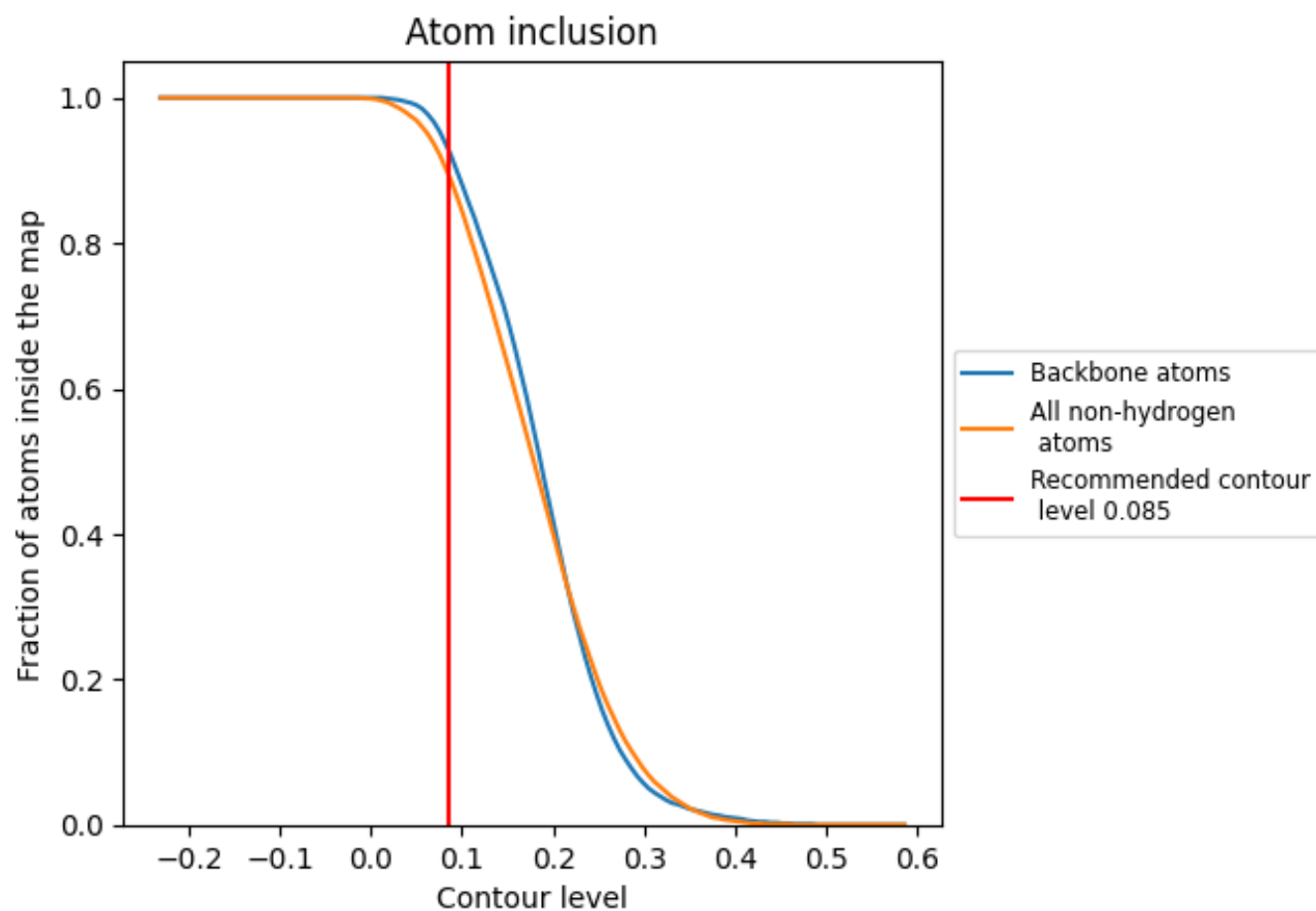
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.085).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 90% 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 (0.085) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8950	 0.5780
0	 0.8120	 0.5740
1	 0.9150	 0.6230
2	 0.9210	 0.6220
3	 0.8460	 0.5910
4	 0.6460	 0.5090
A	 0.9450	 0.5780
B	 0.4610	 0.4770
C	 0.7140	 0.5390
D	 0.7210	 0.5410
E	 0.7820	 0.5670
F	 0.7410	 0.5230
G	 0.7060	 0.5220
H	 0.7940	 0.5690
I	 0.7390	 0.5320
J	 0.6050	 0.4810
K	 0.7800	 0.5590
L	 0.7630	 0.5750
M	 0.7290	 0.5410
N	 0.7870	 0.5460
O	 0.7800	 0.5580
P	 0.8500	 0.5690
Q	 0.7580	 0.5390
R	 0.7920	 0.5450
S	 0.7300	 0.5340
T	 0.8260	 0.5600
U	 0.6050	 0.4680
V	 0.6830	 0.3790
X	 0.8960	 0.5800
Y	 0.7150	 0.4650
Z	 0.8290	 0.5610
a	 0.9670	 0.6020
b	 0.9560	 0.5840
c	 0.8950	 0.6090
d	 0.8600	 0.5990



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Chain	Atom inclusion	Q-score
e	 0.7660	 0.5630
f	 0.7280	 0.5210
g	 0.7120	 0.5090
h	 0.7000	 0.5180
i	 0.8670	 0.5850
j	 0.7740	 0.5930
k	 0.8530	 0.5870
l	 0.8420	 0.5990
m	 0.9040	 0.6080
n	 0.8130	 0.5590
o	 0.7970	 0.5880
p	 0.8950	 0.6020
q	 0.8040	 0.5770
r	 0.8390	 0.5910
s	 0.7730	 0.5500
t	 0.7240	 0.5260
u	 0.7850	 0.5630
v	 0.8320	 0.6000
w	 0.8570	 0.5880
x	 0.7180	 0.5270
y	 0.8050	 0.5820
z	 0.8350	 0.5870