



wwPDB EM Validation Summary Report ⓘ

Jul 13, 2026 – 12:29 pm BST

PDB ID : 9SRA / pdb_00009sra
EMDB ID : EMD-55135
Title : Cryo-EM structure of P. abyssi HibA:ribosome with an SD:antiSD duplex
Authors : Madru, C.M.; Bourgeois, G.B.; Mechulam, Y.M.; Schmitt, E.S.
Deposited on : 2025-09-24
Resolution : 2.20 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.dev133
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

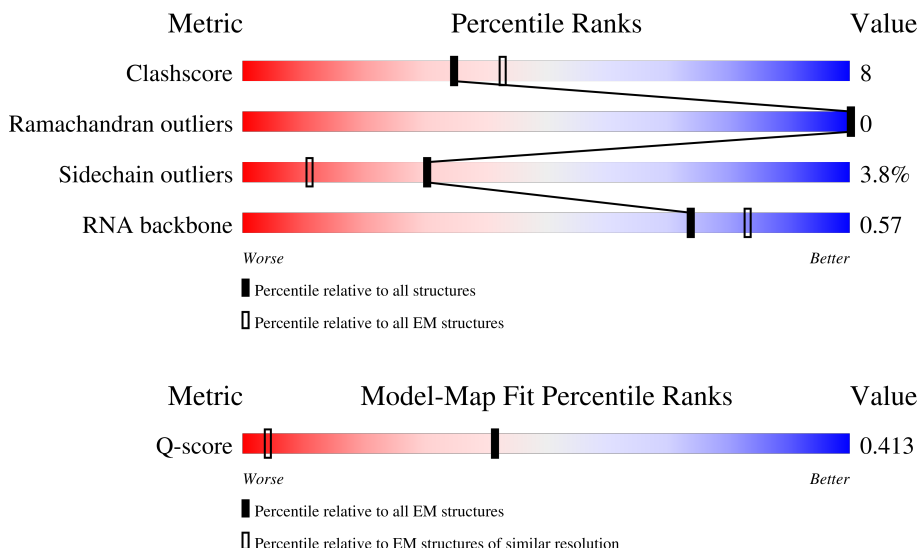
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.20 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	3184 (1.71 - 2.70)

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	1	3018	
2	2	1512	
3	3	128	

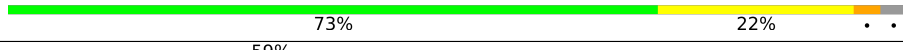
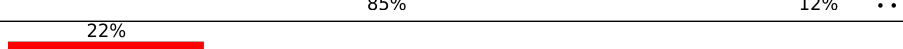
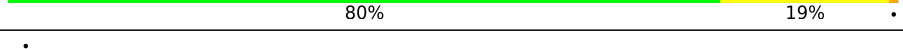
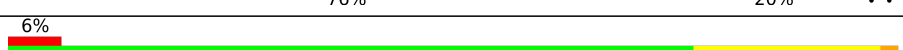
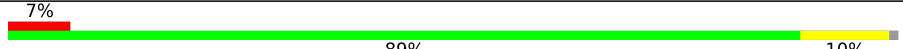
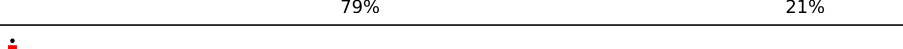
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Mol	Chain	Length	Quality of chain
4	H	392	 66% 76% 20% . .
5	AA	199	 70% 23% . 6%
6	AB	202	 83% 14% .
7	AC	63	 71% 19% 10%
8	AD	180	 85% 11% .
9	AE	243	 79% 20%
10	AF	236	 83% 14% .
11	AG	125	 72% 26% ..
12	AH	215	 76% 20% .
13	AI	130	 85% 14% ..
14	AJ	127	 87% 9% . .
15	AK	135	 84% 16%
16	AL	102	 74% 22% . .
17	AM	137	 82% 10% . 7%
18	AN	147	 84% 15% .
19	AP	56	 62% 34% . .
20	AQ	158	 85% 13% ..
21	AR	113	 78% 16% . 5%
22	AS	67	 70% 25% .
23	AU	150	 81% 17% ..
24	AV	99	 85% 12% .
24	B6	99	 61% 64% 28% . 5%
25	AW	65	 82% 12% 6%
26	AX	71	 75% 11% 6% 8%
27	AY	51	 63% 61% 35% .

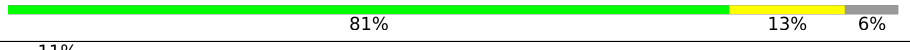
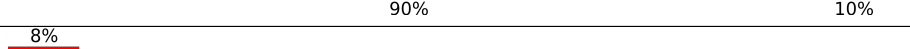
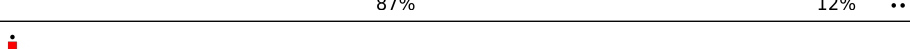
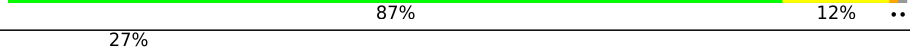
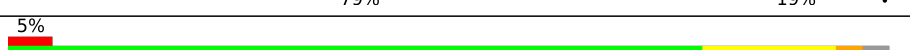
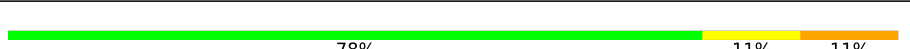
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Mol	Chain	Length	Quality of chain
28	AZ	210	
29	A0	37	
30	A3	123	
30	B4	123	
30	BG	123	
31	AT	132	
32	AO	148	
33	BB	239	
34	BY	155	
35	BO	203	
36	BC	361	
37	B5	82	
37	BK	82	
38	BL	147	
39	Bf	51	
40	BU	121	
41	Bb	130	
42	Be	62	
43	BE	188	
44	Ba	94	
45	BT	86	
46	BW	68	
47	Bi	83	
48	BI	142	
49	BR	97	

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Mol	Chain	Length	Quality of chain
50	BQ	151	 81% 19% .
51	BV	67	 81% 13% 6%
52	Bj	94	 11% 90% 10%
53	BD	255	 8% 80% 19% .
54	BF	184	 5% 79% 18% ..
55	BZ	99	 79% 17% ..
56	BP	120	 12% 85% 14% .
57	BM	194	 82% 16% ..
58	BS	155	 87% 12% ..
59	Bd	91	 91% 9%
60	BN	181	 9% 76% 22% ..
61	Bg	51	 76% 18% 6%
62	Bc	87	 6% 80% 20%
63	BJ	141	 87% 12% ..
64	Bl	77	 27% 79% 19% .
65	Bk	65	 5% 78% 15% . .
66	sd	9	 78% 11% 11%

2 Entry composition [i](#)

There are 68 unique types of molecules in this entry. The entry contains 170684 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 rRNA 23S.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	1	3018	Total	C	N	O	P	0	0
			65181	29055	12094	21014	3018		

- Molecule 2 is a RNA chain called rRNA 16S.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	2	1497	Total	C	N	O	P	0	0
			32311	14417	5959	10438	1497		

- Molecule 3 is a RNA chain called rRNA 5S.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	3	126	Total	C	N	O	P	0	0
			2703	1203	498	876	126		

- Molecule 4 is a protein called Dehydrogenase.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	H	386	Total	C	N	O	S	0	0
			3074	1962	535	567	10		

- Molecule 5 is a protein called 30S ribosomal protein S3Ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	AA	188	Total	C	N	O	S	0	0
			1531	993	268	266	4		

- Molecule 6 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	AB	196	Total	C	N	O	S	0	0
			1571	1017	269	281	4		

- Molecule 7 is a protein called Zn-ribbon RNA-binding protein involved in translation.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	AC	57	Total	C	N	O	S	0	0
			449	285	80	76	8		

- Molecule 8 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	AD	173	Total	C	N	O	S	0	0
			1452	913	280	255	4		

- Molecule 9 is a protein called 30S ribosomal protein S4e.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	AE	242	Total	C	N	O	S	0	0
			1983	1281	358	339	5		

- Molecule 10 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	AF	229	Total	C	N	O	S	0	0
			1808	1147	334	320	7		

- Molecule 11 is a protein called 30S ribosomal protein S6e.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	AG	124	Total	C	N	O	S	0	0
			977	621	178	176	2		

- Molecule 12 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	AH	214	Total	C	N	O	S	0	0
			1725	1095	323	300	7		

- Molecule 13 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	AI	129	Total	C	N	O	S	0	0
			1034	668	184	180	2		

- Molecule 14 is a protein called 30S ribosomal protein S8e.

Mol	Chain	Residues	Atoms				AltConf	Trace
14	AJ	125	Total	C	N	O		
			986	612	205	169	0	0

- Molecule 15 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms				AltConf	Trace
15	AK	135	Total	C	N	O	S	
			1073	673	207	189	4	0

- Molecule 16 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms				AltConf	Trace
16	AL	100	Total	C	N	O	S	
			809	502	157	147	3	0

- Molecule 17 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms				AltConf	Trace
17	AM	127	Total	C	N	O	S	
			955	591	190	172	2	0

- Molecule 18 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms				AltConf	Trace
18	AN	146	Total	C	N	O	S	
			1148	727	224	194	3	0

- Molecule 19 is a protein called 30S ribosomal protein S14 type Z.

Mol	Chain	Residues	Atoms				AltConf	Trace
19	AP	55	Total	C	N	O	S	
			455	288	95	67	5	0

- Molecule 20 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms				AltConf	Trace
20	AQ	157	Total	C	N	O	S	
			1305	833	249	219	4	0

- Molecule 21 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	AR	107	Total	C	N	O	S	0	0
			884	562	172	147	3		

- Molecule 22 is a protein called 30S ribosomal protein S17e.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	AS	64	Total	C	N	O	S	0	0
			541	343	104	93	1		

- Molecule 23 is a protein called 30S ribosomal protein S19e.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	AU	149	Total	C	N	O	S	0	0
			1223	790	221	212			

- Molecule 24 is a protein called 30S ribosomal protein S24e.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	AV	96	Total	C	N	O	S	0	0
			808	528	129	148	3		
24	B6	94	Total	C	N	O	S	0	0
			790	516	125	146	3		

- Molecule 25 is a protein called 30S ribosomal protein S27e.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	AW	61	Total	C	N	O	S	0	0
			470	294	91	80	5		

- Molecule 26 is a protein called 30S ribosomal protein S28e.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	AX	65	Total	C	N	O	S	0	0
			516	316	103	97			

- Molecule 27 is a protein called 30S ribosomal protein S27ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
27	AY	49	Total	C	N	O	S	0	0
			400	257	76	62	5		

- Molecule 28 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	AZ	196	Total	C	N	O	S	0	0
			1541	983	284	270	4		

- Molecule 29 is a protein called eS32.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	A0	36	Total	C	N	O	S	0	0
			343	218	84	39	2		

- Molecule 30 is a protein called 50S ribosomal protein L7Ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	A3	122	Total	C	N	O	S	0	0
			933	594	156	180	3		
30	BG	122	Total	C	N	O	S	0	0
			932	594	156	179	3		
30	B4	122	Total	C	N	O	S	0	0
			933	594	156	180	3		

- Molecule 31 is a protein called 30S ribosomal protein S19.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	AT	130	Total	C	N	O	S	0	0
			1057	675	201	174	7		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
32	AO	138	Total	C	N	O	S	0	0
			1116	700	221	190	5		

- Molecule 33 is a protein called Large ribosomal subunit protein uL2.

Mol	Chain	Residues	Atoms					AltConf	Trace
33	BB	238	Total	C	N	O	S	0	0
			1838	1163	354	317	4		

- Molecule 34 is a protein called Large ribosomal subunit protein uL30.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	BY	154	Total	C	N	O	S	0	0
			1235	783	234	212	6		

- Molecule 35 is a protein called Large ribosomal subunit protein uL18.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	BO	198	Total	C	N	O	S	0	0
			1607	1024	302	279	2		

- Molecule 36 is a protein called Large ribosomal subunit protein uL3.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	BC	360	Total	C	N	O	S	0	0
			2870	1843	522	494	11		

- Molecule 37 is a protein called Large ribosomal subunit protein eL14.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	B5	81	Total	C	N	O	S	0	0
			614	388	116	108	2		
37	BK	80	Total	C	N	O	S	0	0
			607	383	115	107	2		

- Molecule 38 is a protein called Large ribosomal subunit protein uL15.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	BL	147	Total	C	N	O	S	0	0
			1149	720	224	202	3		

- Molecule 39 is a protein called Large ribosomal subunit protein eL39.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	Bf	50	Total	C	N	O	S	0	0
			435	275	99	60	1		

- Molecule 40 is a protein called Large ribosomal subunit protein uL24.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	BU	121	Total	C	N	O	S	0	0
			1011	638	198	170	5		

- Molecule 41 is a protein called Large ribosomal subunit protein eL32.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	Bb	129	Total	C	N	O	S	0	0
			1095	702	221	171	1		

- Molecule 42 is a protein called Large ribosomal subunit protein eL37.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	Be	61	Total	C	N	O	S	0	0
			503	312	112	75	4		

- Molecule 43 is a protein called Large ribosomal subunit protein uL5.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	BE	185	Total	C	N	O	S	0	0
			1485	934	277	265	9		

- Molecule 44 is a protein called Large ribosomal subunit protein eL31.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	Ba	94	Total	C	N	O	S	0	0
			778	503	147	127	1		

- Molecule 45 is a protein called Large ribosomal subunit protein uL23.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	BT	86	Total	C	N	O	S	0	0
			696	451	118	126	1		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
46	BW	68	Total	C	N	O	S	0	0
			565	349	111	99	6		

- Molecule 47 is a protein called Large ribosomal subunit protein eL43.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	Bi	82	Total	C	N	O	S	0	0
			620	389	129	97	5		

- Molecule 48 is a protein called Large ribosomal subunit protein uL13.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	BI	142	Total	C	N	O	S	0	0
			1148	734	217	194	3		

- Molecule 49 is a protein called Large ribosomal subunit protein eL21.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	BR	96	Total	C	N	O	S	0	0
			797	507	164	125	1		

- Molecule 50 is a protein called Large ribosomal subunit protein eL19.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	BQ	150	Total	C	N	O	S	0	0
			1265	795	261	203	6		

- Molecule 51 is a protein called Large ribosomal subunit protein eL24.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	BV	63	Total	C	N	O	S	0	0
			532	339	103	84	6		

- Molecule 52 is a protein called Large ribosomal subunit protein eL42.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	Bj	94	Total	C	N	O	S	0	0
			789	502	163	119	5		

- Molecule 53 is a protein called Large ribosomal subunit protein uL4.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	BD	255	Total	C	N	O	S	0	0
			2026	1290	388	343	5		

- Molecule 54 is a protein called Large ribosomal subunit protein uL6.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	BF	183	Total	C	N	O	S	0	0
			1456	943	251	261	1		

- Molecule 55 is a protein called Large ribosomal subunit protein eL30.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	BZ	98	Total	C	N	O		0	0
			749	486	122	141			

- Molecule 56 is a protein called Large ribosomal subunit protein eL18.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	BP	120	Total	C	N	O	S	0	0
			971	610	188	170	3		

- Molecule 57 is a protein called Large ribosomal subunit protein eL15.

Mol	Chain	Residues	Atoms					AltConf	Trace
57	BM	193	Total	C	N	O	S	0	0
			1588	1014	318	251	5		

- Molecule 58 is a protein called Large ribosomal subunit protein uL22.

Mol	Chain	Residues	Atoms					AltConf	Trace
58	BS	154	Total	C	N	O	S	0	0
			1247	786	246	211	4		

- Molecule 59 is a protein called Large ribosomal subunit protein eL34.

Mol	Chain	Residues	Atoms					AltConf	Trace
59	Bd	91	Total	C	N	O	S	0	0
			759	477	163	109	10		

- Molecule 60 is a protein called Large ribosomal subunit protein uL16.

Mol	Chain	Residues	Atoms					AltConf	Trace
60	BN	178	Total	C	N	O	S	0	0
			1452	920	281	246	5		

- Molecule 61 is a protein called Large ribosomal subunit protein eL40.

Mol	Chain	Residues	Atoms					AltConf	Trace
61	Bg	48	Total	C	N	O	S	0	0
			387	243	80	60	4		

- Molecule 62 is a protein called Large ribosomal subunit protein eL33.

Mol	Chain	Residues	Atoms					AltConf	Trace
62	Bc	87	Total	C	N	O	S	0	0
			684	435	132	115	2		

- Molecule 63 is a protein called Large ribosomal subunit protein uL14.

Mol	Chain	Residues	Atoms					AltConf	Trace
63	BJ	140	Total	C	N	O	S	0	0
			1066	664	214	185	3		

- Molecule 64 is a protein called Large ribosomal subunit protein eL20.

Mol	Chain	Residues	Atoms					AltConf	Trace
64	Bl	77	Total	C	N	O	S	0	0
			659	425	120	113	1		

- Molecule 65 is a protein called C2H2-type domain-containing protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
65	Bk	63	Total	C	N	O	S	0	0
			528	339	108	78	3		

- Molecule 66 is a RNA chain called antiSD-RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
66	sd	9	Total	C	N	O	P	0	0
			199	88	39	63	9		

- Molecule 67 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
67	1	166	Total	Mg	0
			166	166	
67	2	69	Total	Mg	0
			69	69	
67	3	1	Total	Mg	0
			1	1	
67	AK	1	Total	Mg	0
			1	1	
67	AN	1	Total	Mg	0
			1	1	
67	BB	2	Total	Mg	0
			2	2	
67	Bb	1	Total	Mg	0
			1	1	
67	BD	1	Total	Mg	0
			1	1	
67	BM	1	Total	Mg	0
			1	1	

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Mol	Chain	Residues	Atoms		AltConf
67	BS	1	Total	Mg	0
			1	1	

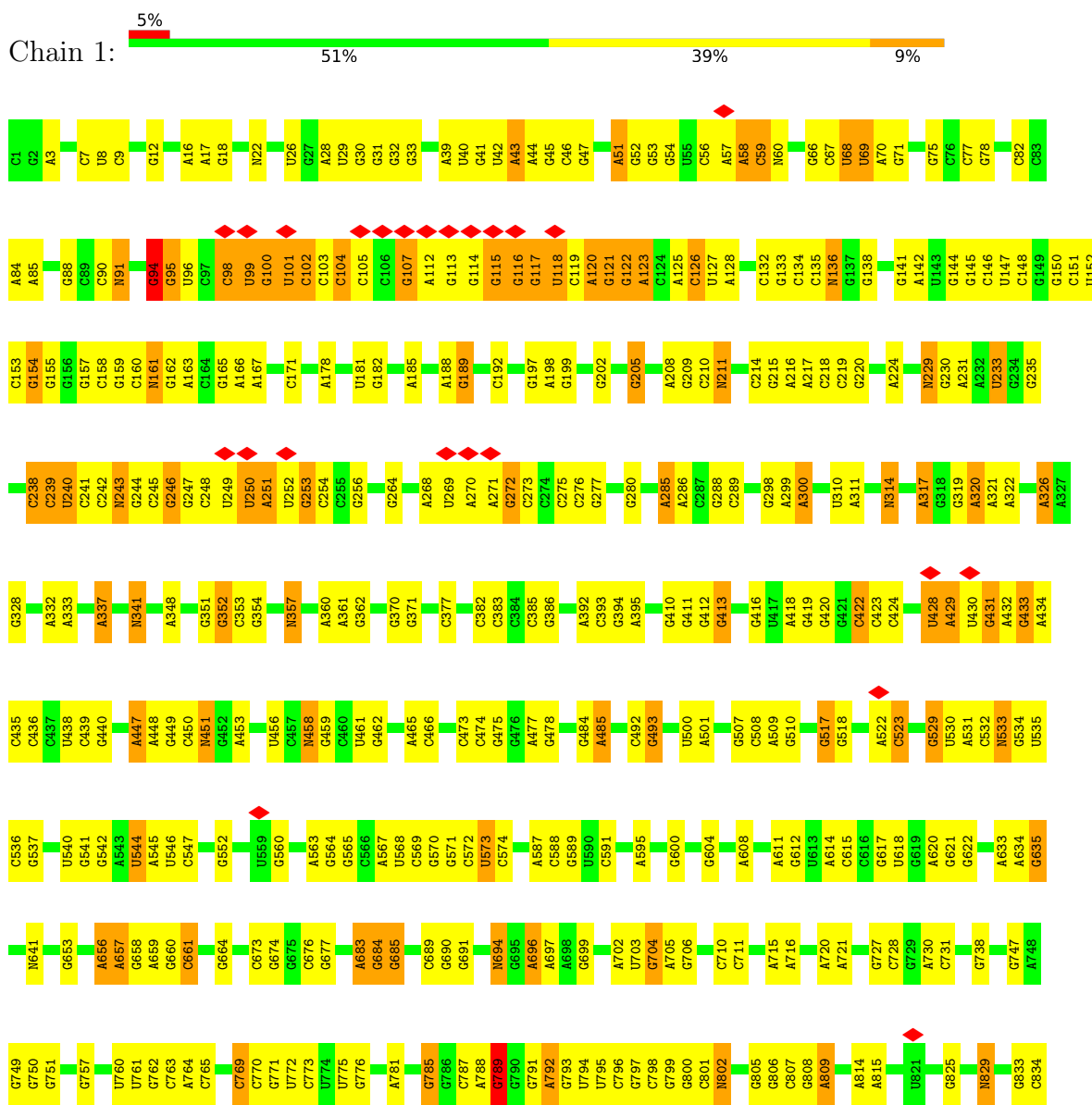
- Molecule 68 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
68	AC	2	Total	Zn	0
			2	2	
68	AF	1	Total	Zn	0
			1	1	
68	AP	1	Total	Zn	0
			1	1	
68	AR	1	Total	Zn	0
			1	1	
68	AW	1	Total	Zn	0
			1	1	
68	Be	1	Total	Zn	0
			1	1	
68	Bi	1	Total	Zn	0
			1	1	
68	BV	1	Total	Zn	0
			1	1	
68	Bj	1	Total	Zn	0
			1	1	
68	Bd	1	Total	Zn	0
			1	1	
68	Bg	1	Total	Zn	0
			1	1	
68	Bk	1	Total	Zn	0
			1	1	

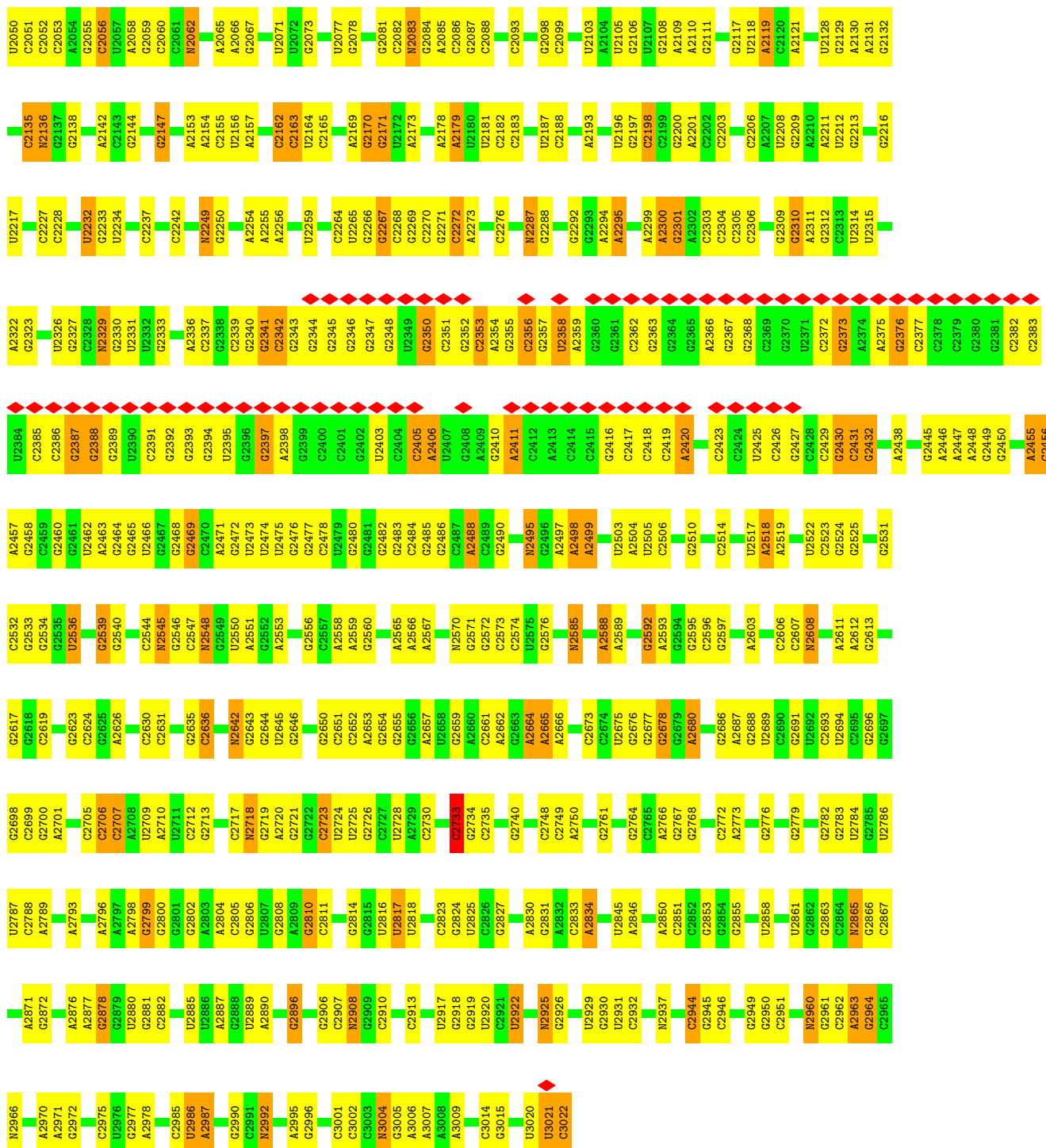
3 Residue-property plots

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: rRNA 23S

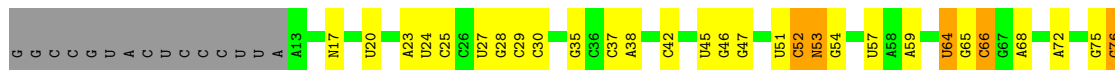


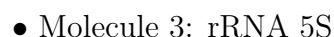
G1943	A1854	G1769	A1582	U1482	G1392	G1323	A1251	C1153	G1078	U993	A919	M835
G1950	A1855	A1780	U1583	A1482	G1393	A1324	G1252	C1154	G1079	U998	C920	G841
C1951	U1856	G1761	C1587	A1484	C1394	A1325	G1253	C1155	A1081		C921	N922
A1952	U1857	U1762	C1588	C1489	U1395	G1326	C1254	U1159	A1084	G1006	N923	A842
A1953	C1858	G1763	U1589	G1490	C1396	C1327	C1255	G1161	U1085	G1007	A924	A844
A1954	C1859	G1764	C1590	C1498	A1397	A1328	C1256	A1162	G1086	C1011	A925	C845
	U1862	N1765	A1591	G1499	C1401	G1329	A1259	A1163	G1087	G1012	A926	G847
	N1867	A1769	C1592	C1499	G1402	G1330	U1260	G1168		G1013	G929	A848
	U1770	N1594	C1593	G1508	G1406	C1331	U1261	C1093	G1094	U1014	G930	C849
			N1594	C1509	C1407	C1332	U1262	C1094	G1095	A1015	U931	A850
			G1598			U1333	G1263	G1095	G1096	C1019	A854	U851
			C1599		G1411	C1334	C1265	A1170		G1020	U935	C855
			U1600		C1415	G1335	G1266	C1171	A1101	G1026	G936	G856
			U1601				U1269	C1172	G1102	G1027	U937	C857
			G1517				U1272	C1173	G1103	G1031	A938	A858
			G1520		G1420	U1337		C1104	C1105		G939	
					C1421	U1338	G1272	G1105		A1032	C940	G864
					A1422	A1339	C1278	G1106		A1033	C941	C865
					C1423	A1340	U1279	U1107		A1034	C946	C866
					G1424	A1341	C1280	G1108		G1035	C947	U867
					A1426	A1342	C1281	G1109		G1036	C948	C868
						A1343	A1282	G1110		G1037	C949	G869
						A1344	A1283	G1111		G950	C950	A870
						A1345	A1284			A1038	G951	G871
						A1346	G1285	A1115		A1039	C952	C872
						A1347	U1288	G1116		A1040	U953	
						G1348	G1289	C1117		U1041	G954	U877
						U1349	U1290	C1118		C1042	A955	G878
						A1350	N1293	G1119		G1043	G956	A879
						A1351	G1294	U1120			G957	G880
						A1352	G1295	G1121		G1046	C958	G887
						A1353	U1299	A1122		N1048	C959	G888
						G1354	U1300	U1123		A1053	C960	U889
						C1355	A1301	G1124		U1054	U961	G890
						U1356	G1302	G1125		A1055	C962	C891
						U1357	A1303	G1129		U1056	A963	C892
						A1358	C1304	G1130		C1057	A964	A893
						C1359	A1305	G1131		C1058	C965	A894
						U1360	G1306	C1132		U1059	A966	A895
						C1361	C1307	A1133		G1060	C967	G896
						U1362	G1308	G1134			U973	C897
						C1363	C1309	G1135		C1063	G974	G898
						U1364	G1310	G1136		U1065	A976	G903
						C1365	G1311	G1137		C1066	G977	U906
						U1366	U1312	G1138		C1067	C978	U806
						G1367	U1313	C1139		U1068	C979	A908
						C1370	U1314	G1140		A1071	C980	C909
						U1371	U1315	A1141		U1072	G981	G910
						G1372	G1316	A1142		C1073	G982	
						U1373	A1317	A1143		U1074	U984	C913
						G1374	G1318	C1144		A1075	U990	G914
						G1375	C1319	C1145		C1076	G991	A917
						U1376	U1320	C1146		U1077	C992	G918
						G1377	U1321	C1147		C1078		
						U1378	U1322	C1148		C1079		
						G1379	U1323					
						U1380	U1324					
						G1381	U1325					
						U1382	U1326					
						G1383	U1327					
						U1384	U1328					
						G1385	U1329					
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						G1387	U1331					
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						G1670	U1424					
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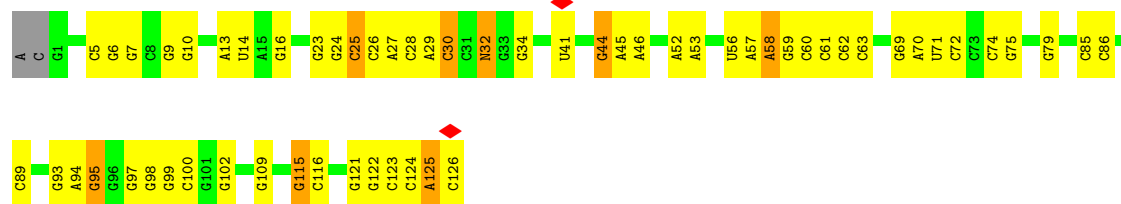
• Molecule 2: rRNA 16S

Chain 2: 58% 34% 7%



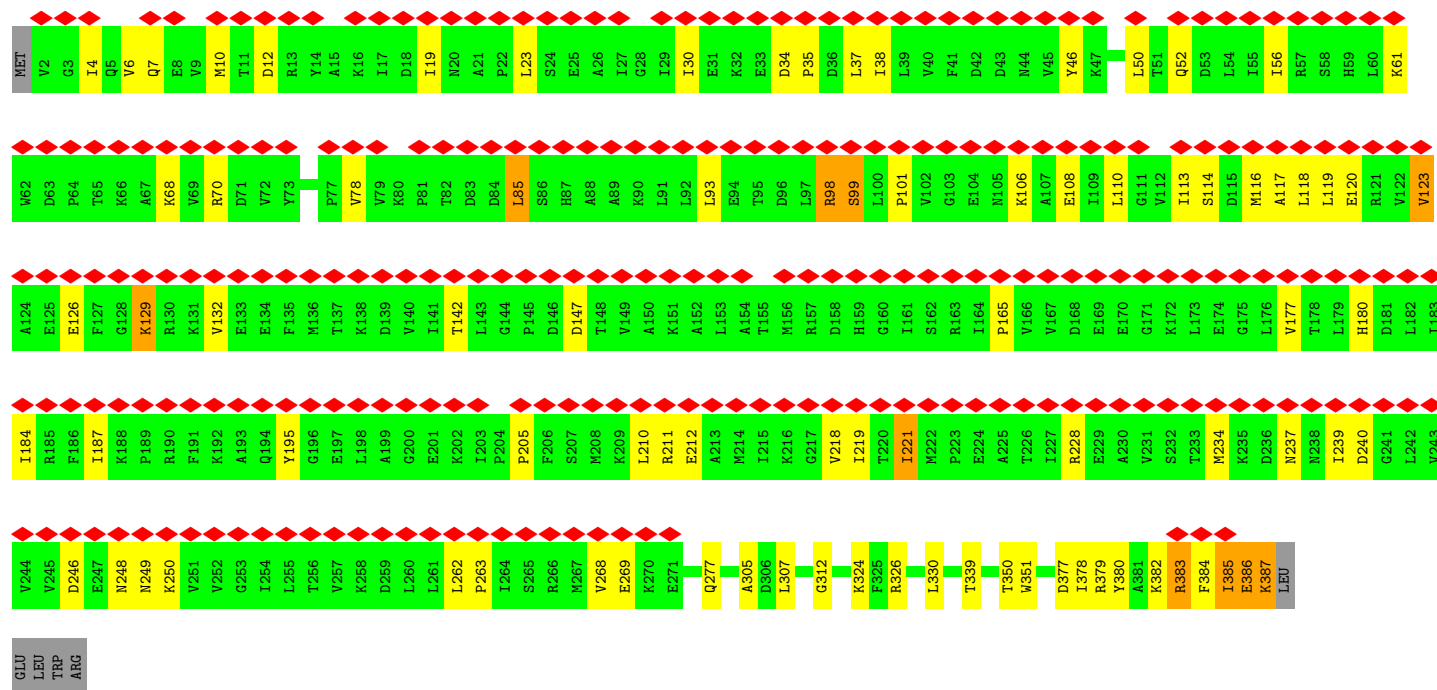


Chain 3: 



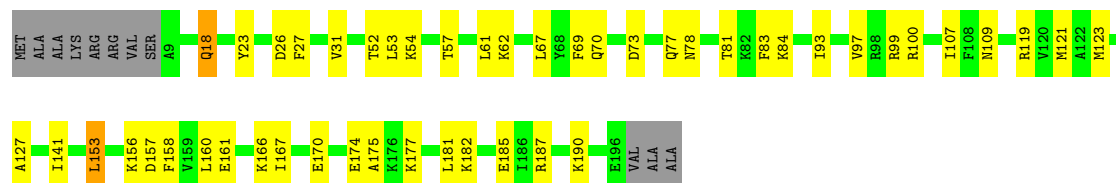
• Molecule 4: Dehydrogenase

Chain H: 



• Molecule 5: 30S ribosomal protein S3Ae

Chain AA: 

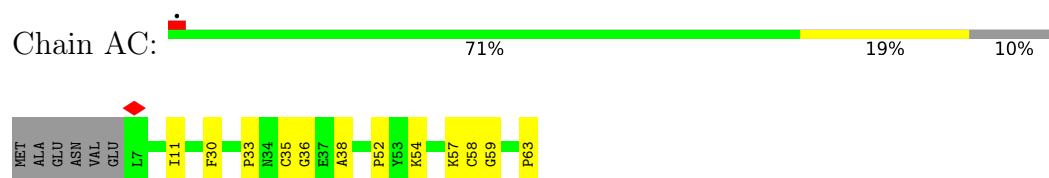


• Molecule 6: 30S ribosomal protein S2

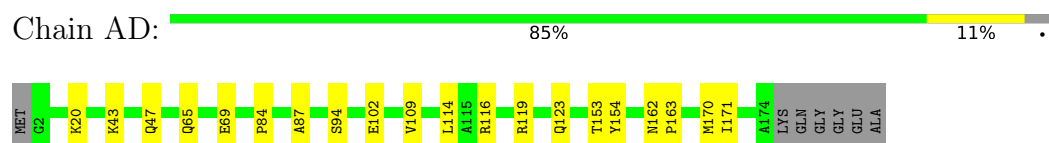
Chain AB: 



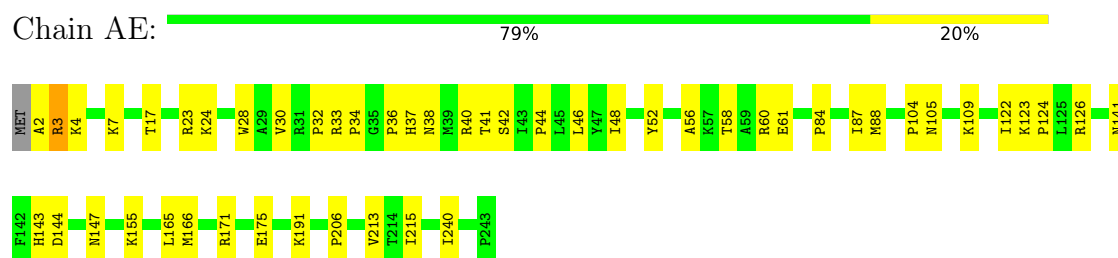
- Molecule 7: Zn-ribbon RNA-binding protein involved in translation



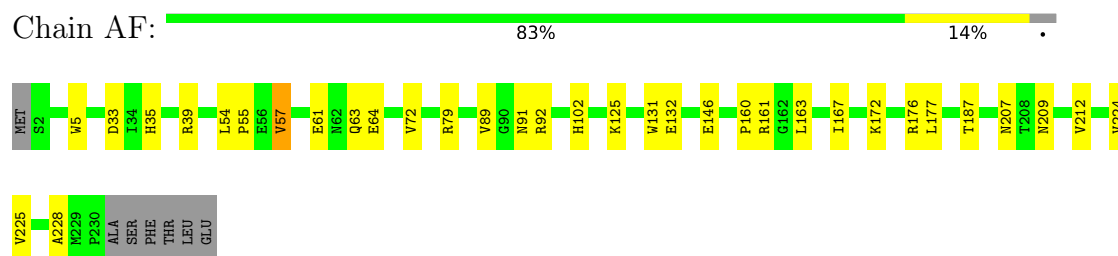
- Molecule 8: 30S ribosomal protein S4



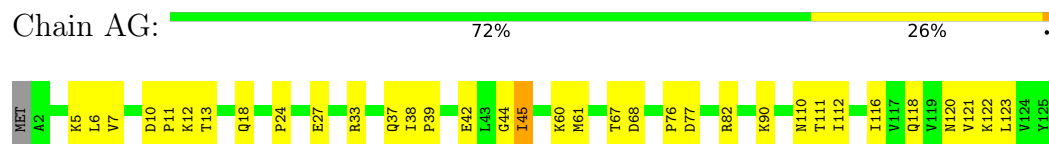
- Molecule 9: 30S ribosomal protein S4e



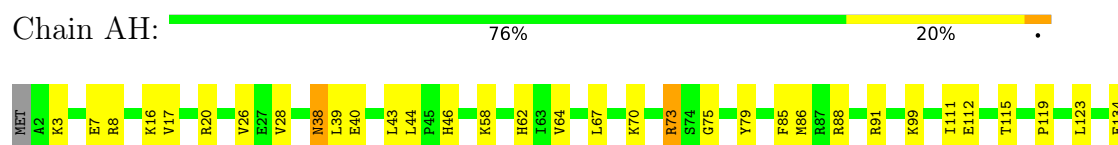
- Molecule 10: 30S ribosomal protein S5



- Molecule 11: 30S ribosomal protein S6e



- Molecule 12: 30S ribosomal protein S7





- Molecule 13: 30S ribosomal protein S8

Chain AI: 85% 14% ..



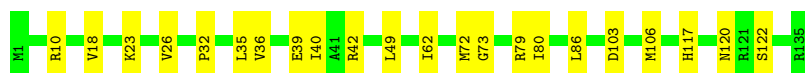
- Molecule 14: 30S ribosomal protein S8e

Chain AJ: 87% 9% ..



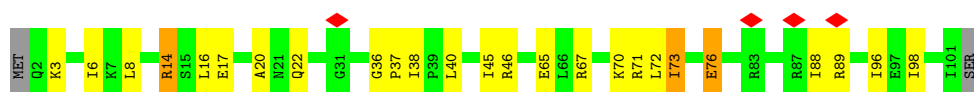
- Molecule 15: 30S ribosomal protein S9

Chain AK: 84% 16%



- Molecule 16: 30S ribosomal protein S10

Chain AL: 74% 22% ..



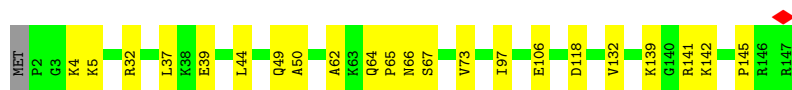
- Molecule 17: 30S ribosomal protein S11

Chain AM: 82% 10% 7%



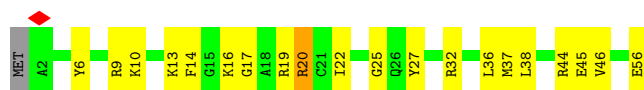
- Molecule 18: 30S ribosomal protein S12

Chain AN: 84% 15%



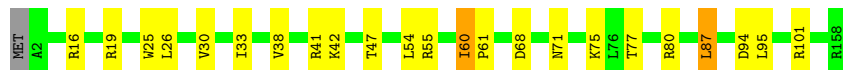
- Molecule 19: 30S ribosomal protein S14 type Z

Chain AP: 62% 34%



- Molecule 20: 30S ribosomal protein S15

Chain AQ: 85% 13% ..



- Molecule 21: 30S ribosomal protein S17

Chain AR: 78% 16% • 5%



- Molecule 22: 30S ribosomal protein S17e

Chain AS: 70% 25% •



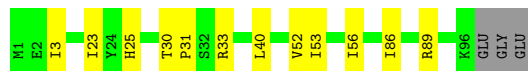
- Molecule 23: 30S ribosomal protein S19e

Chain AU: 81% 17% ..



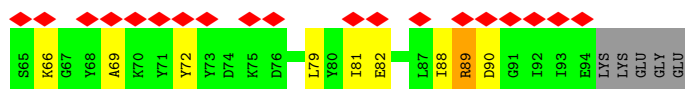
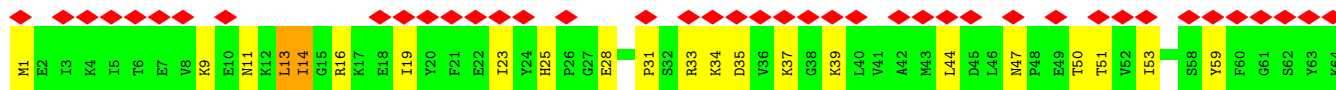
- Molecule 24: 30S ribosomal protein S24e

Chain AV: 85% 12% •



- Molecule 24: 30S ribosomal protein S24e

Chain B6: 61% 64% 28% • 5%



- Molecule 25: 30S ribosomal protein S27e

Chain AW: 



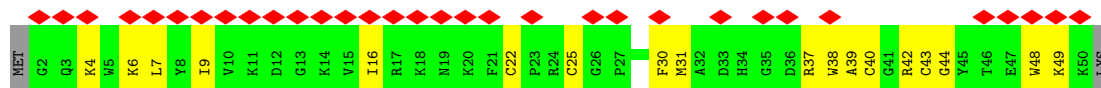
- Molecule 26: 30S ribosomal protein S28e

Chain AX: 



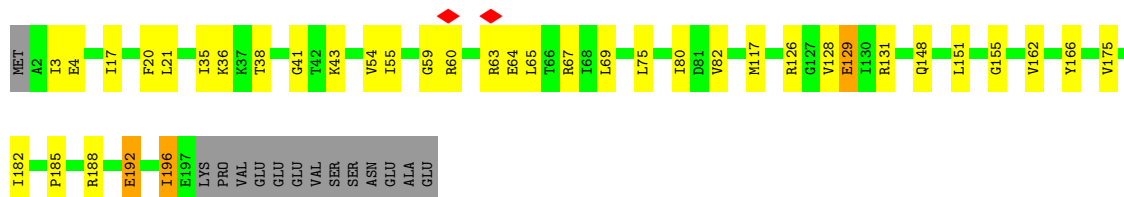
- Molecule 27: 30S ribosomal protein S27ae

Chain AY: 



- Molecule 28: 30S ribosomal protein S3

Chain AZ: 



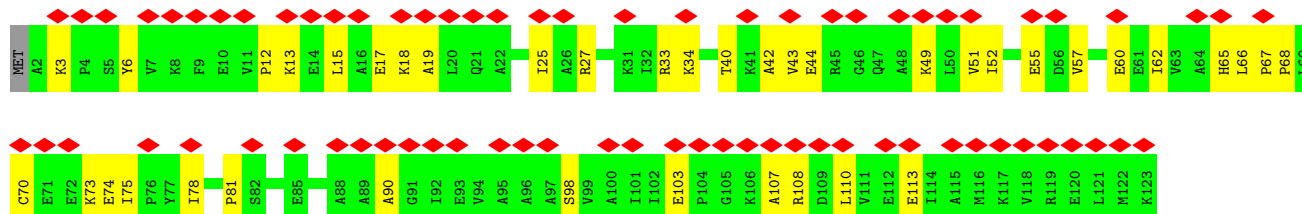
- Molecule 29: eS32

Chain A0: 

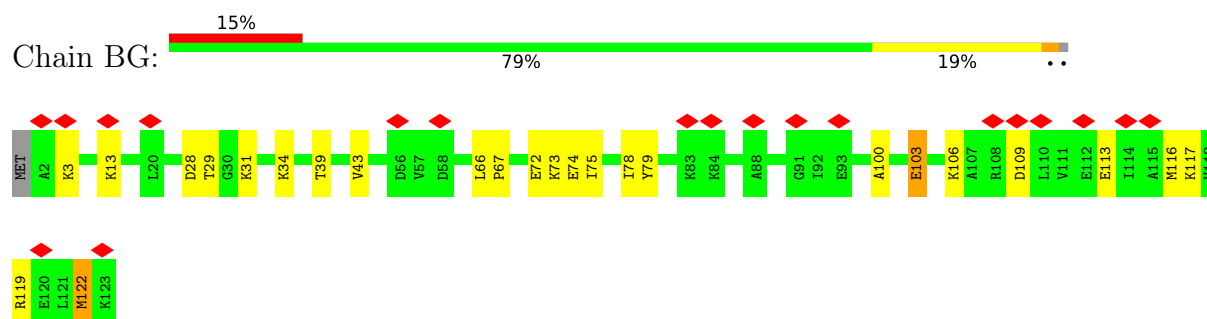


- Molecule 30: 50S ribosomal protein L7Ae

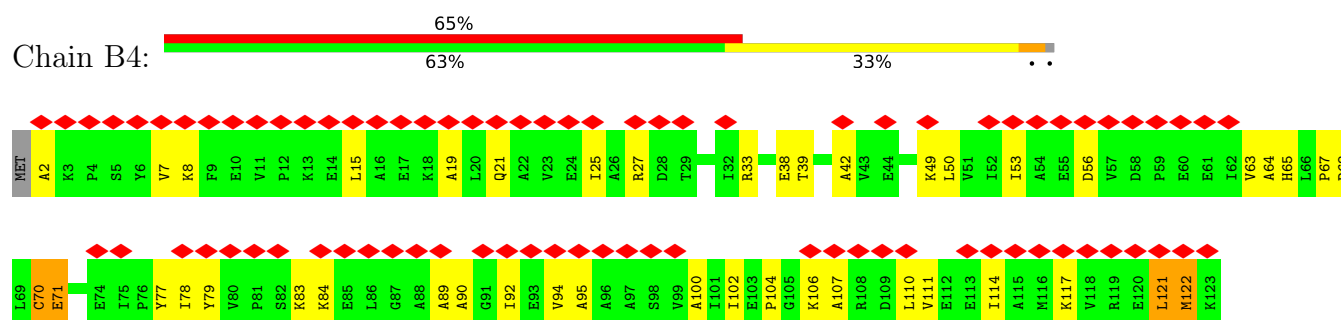
Chain A3: 



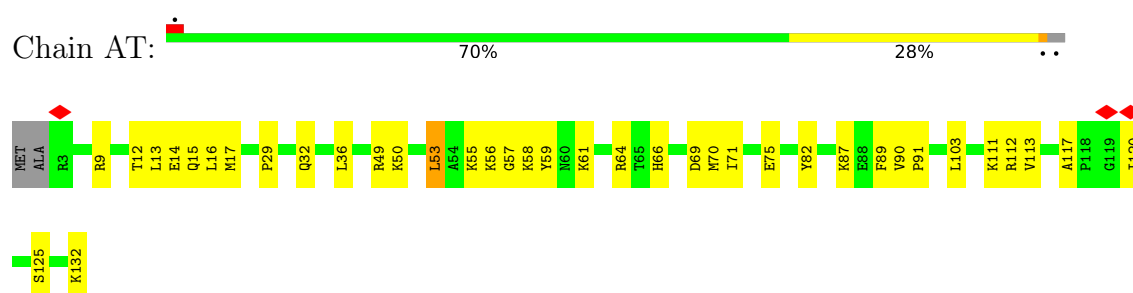
- Molecule 30: 50S ribosomal protein L7Ae



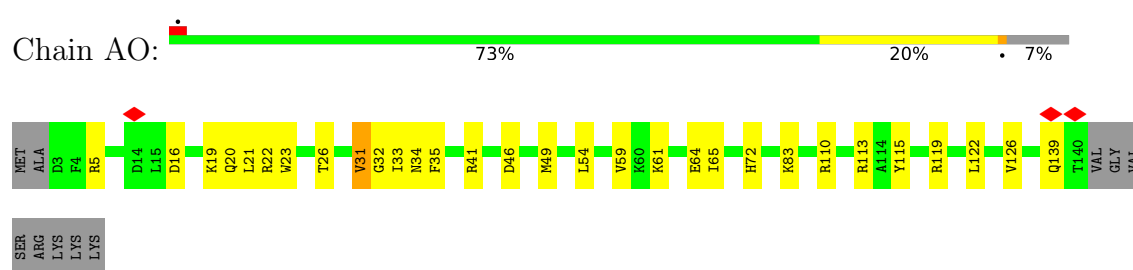
- Molecule 30: 50S ribosomal protein L7Ae



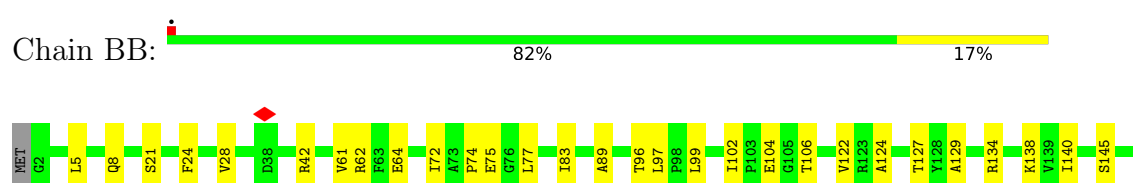
- Molecule 31: 30S ribosomal protein S19



- Molecule 32: 30S ribosomal protein S13

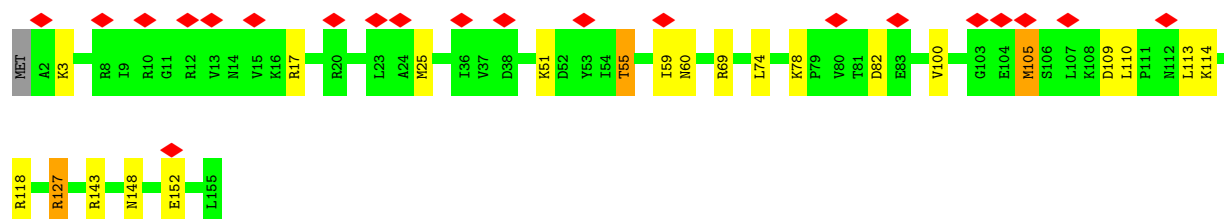
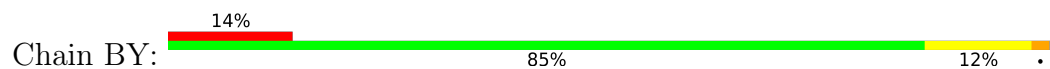


- Molecule 33: Large ribosomal subunit protein uL2

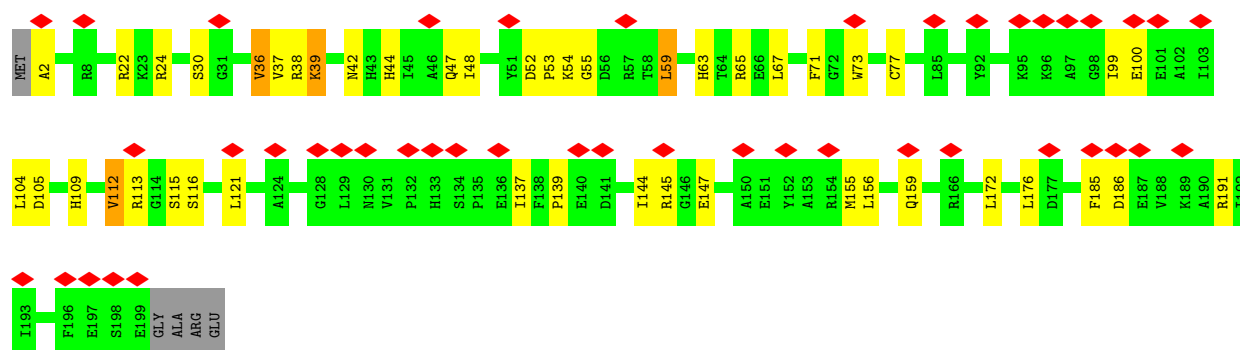
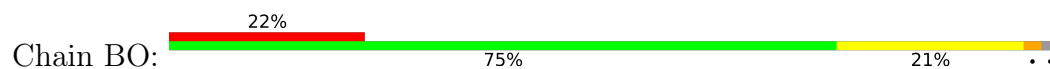




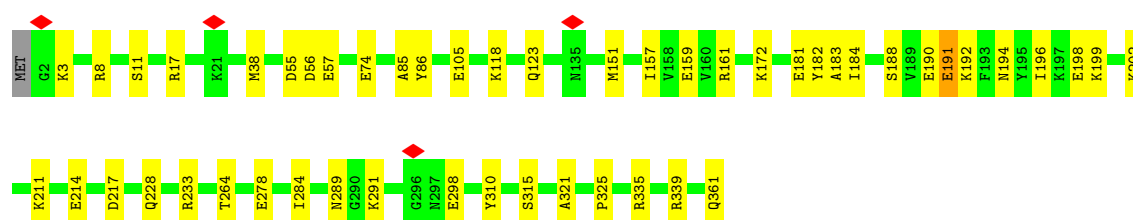
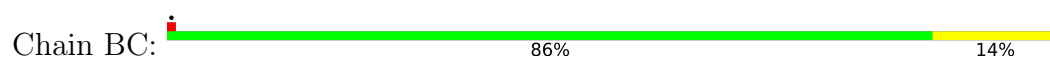
- Molecule 34: Large ribosomal subunit protein uL30



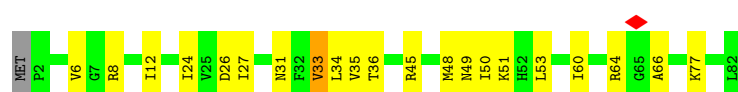
- Molecule 35: Large ribosomal subunit protein uL18



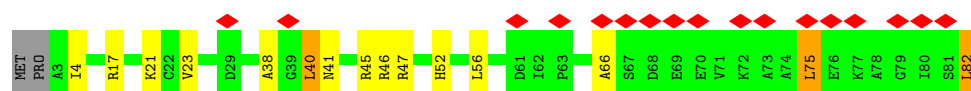
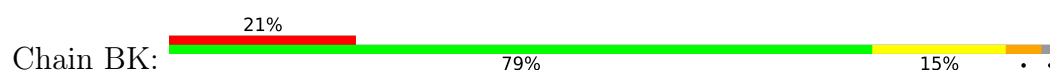
- Molecule 36: Large ribosomal subunit protein uL3



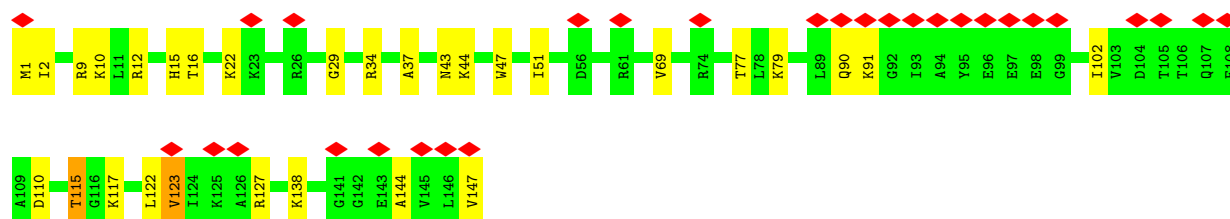
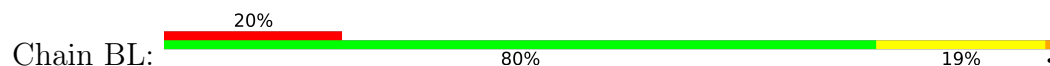
- Molecule 37: Large ribosomal subunit protein eL14



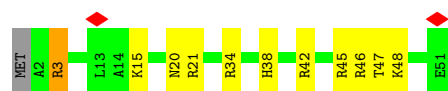
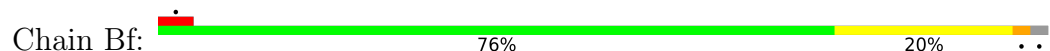
- Molecule 37: Large ribosomal subunit protein eL14



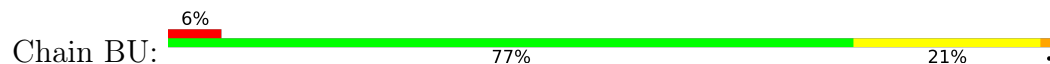
- Molecule 38: Large ribosomal subunit protein uL15



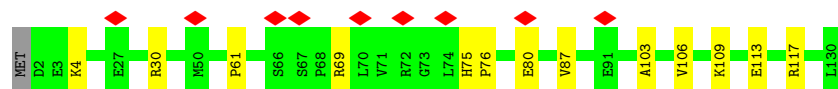
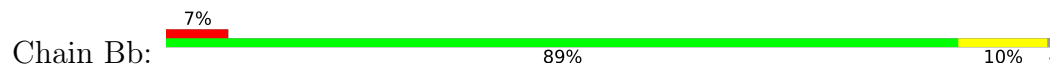
- Molecule 39: Large ribosomal subunit protein eL39



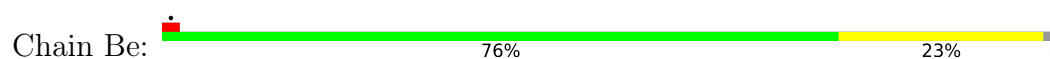
- Molecule 40: Large ribosomal subunit protein uL24



- Molecule 41: Large ribosomal subunit protein eL32

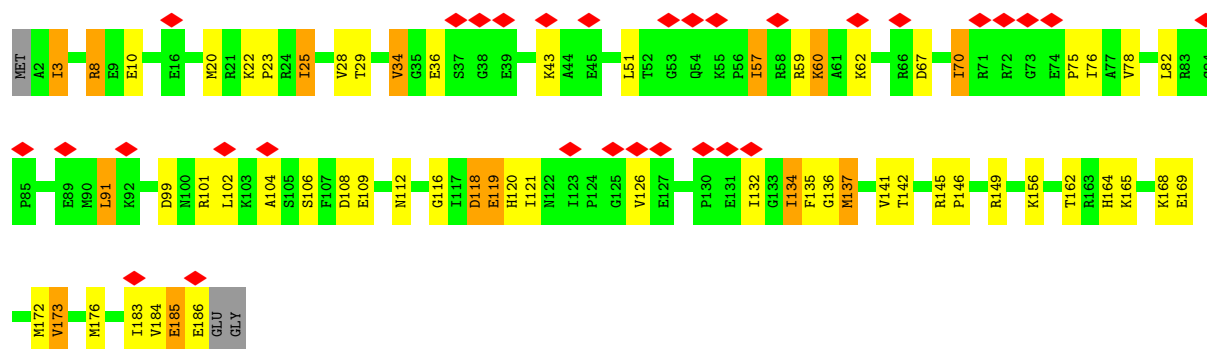


- Molecule 42: Large ribosomal subunit protein eL37

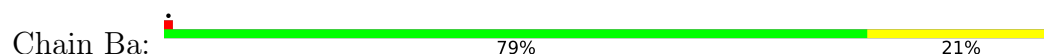


- Molecule 43: Large ribosomal subunit protein uL5

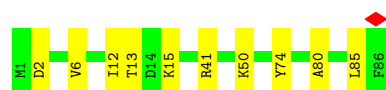




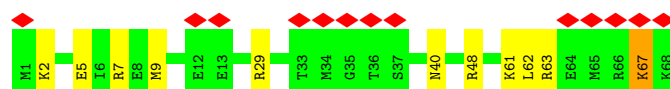
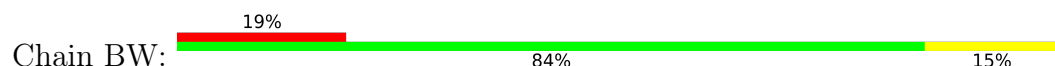
- Molecule 44: Large ribosomal subunit protein eL31



- Molecule 45: Large ribosomal subunit protein uL23



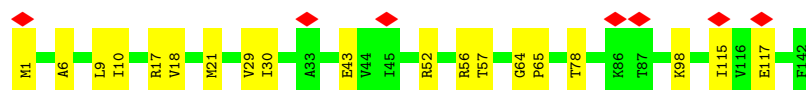
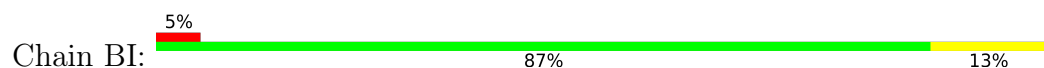
- Molecule 46: Large ribosomal subunit protein uL29



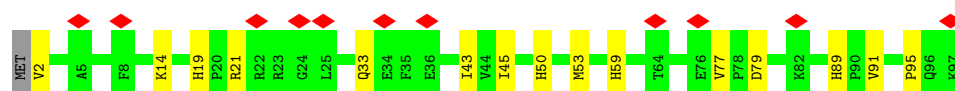
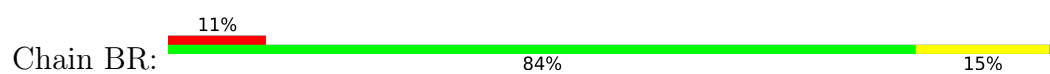
- Molecule 47: Large ribosomal subunit protein eL43



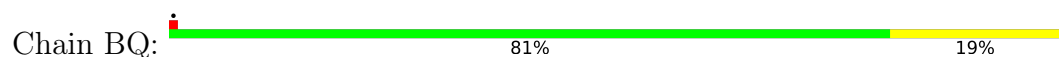
- Molecule 48: Large ribosomal subunit protein uL13



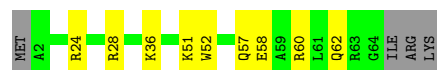
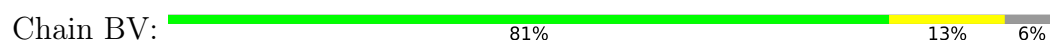
- Molecule 49: Large ribosomal subunit protein eL21



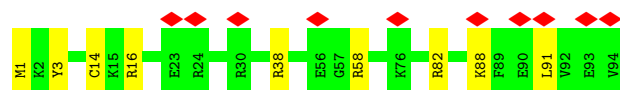
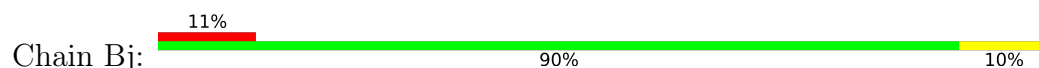
- Molecule 50: Large ribosomal subunit protein eL19



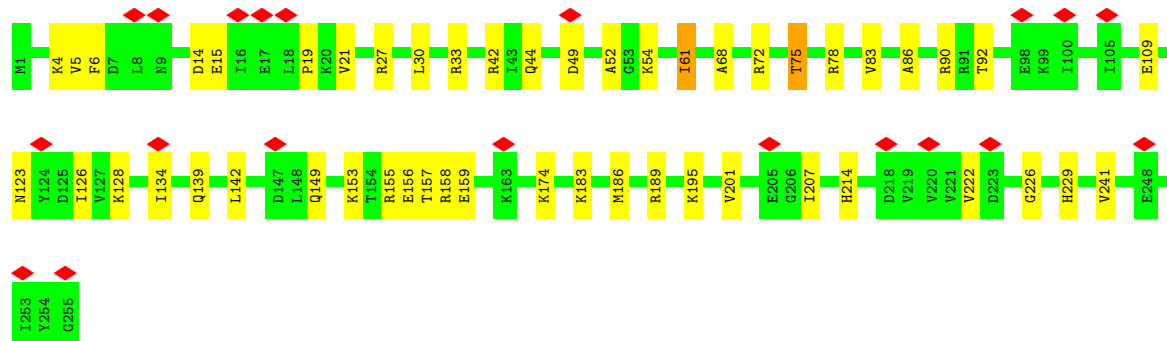
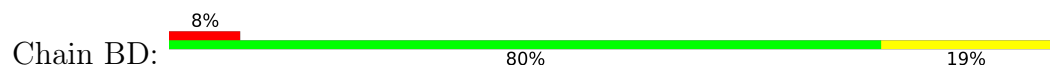
- Molecule 51: Large ribosomal subunit protein eL24



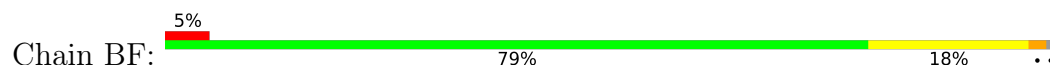
- Molecule 52: Large ribosomal subunit protein eL42



- Molecule 53: Large ribosomal subunit protein uL4

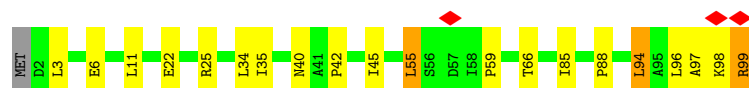
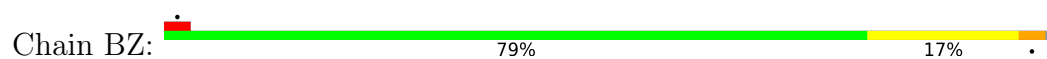


- Molecule 54: Large ribosomal subunit protein uL6

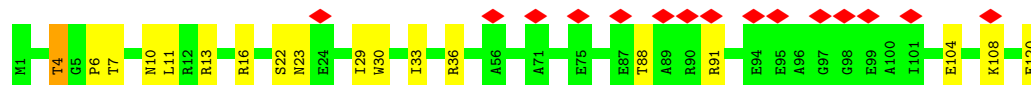
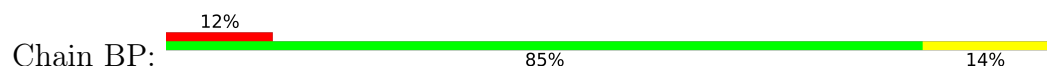




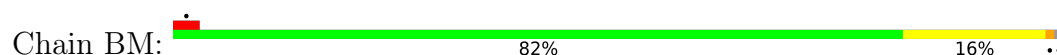
- Molecule 55: Large ribosomal subunit protein eL30



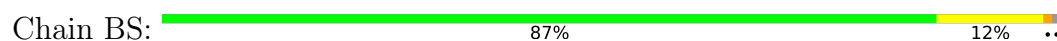
- Molecule 56: Large ribosomal subunit protein eL18



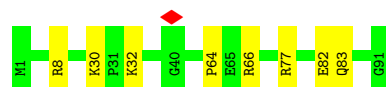
- Molecule 57: Large ribosomal subunit protein eL15



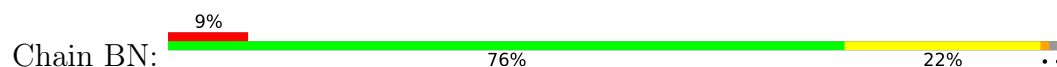
- Molecule 58: Large ribosomal subunit protein uL22

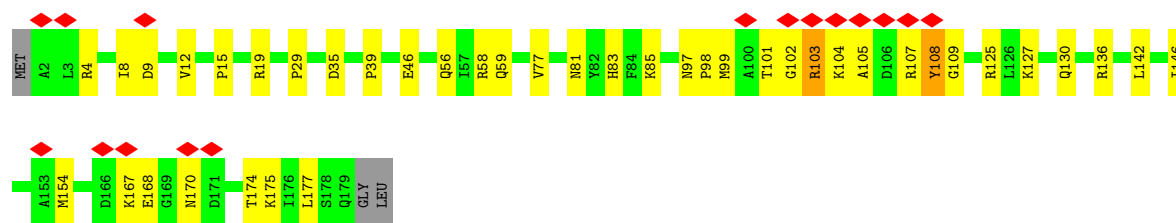


- Molecule 59: Large ribosomal subunit protein eL34

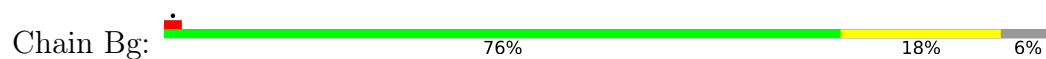


- Molecule 60: Large ribosomal subunit protein uL16

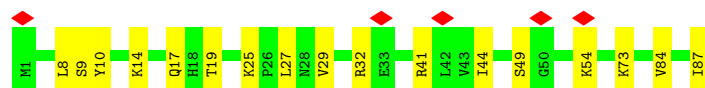
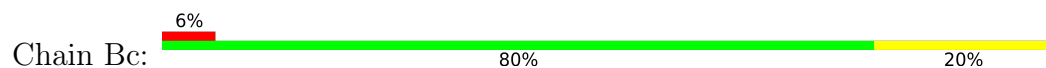




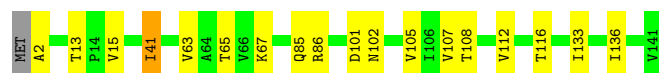
- Molecule 61: Large ribosomal subunit protein eL40



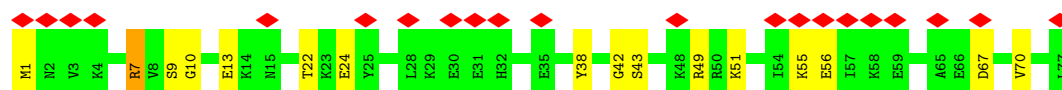
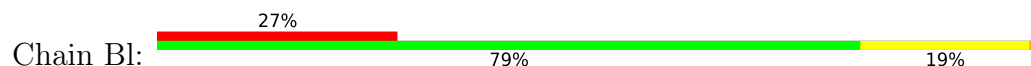
- Molecule 62: Large ribosomal subunit protein eL33



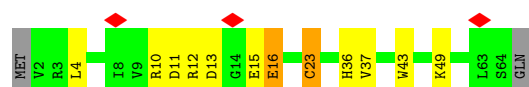
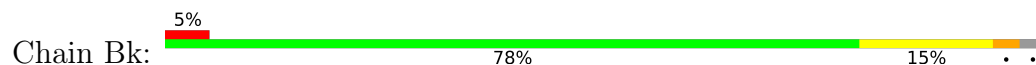
- Molecule 63: Large ribosomal subunit protein uL14



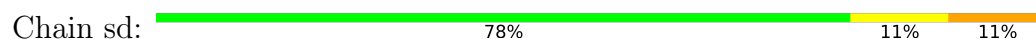
- Molecule 64: Large ribosomal subunit protein eL20



- Molecule 65: C2H2-type domain-containing protein



- Molecule 66: antiSD-RNA





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	182000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.891	Depositor
Minimum map value	-0.355	Depositor
Average map value	0.007	Depositor
Map value standard deviation	0.050	Depositor
Recommended contour level	0.07	Depositor
Map size (Å)	302.04, 302.04, 302.04	wwPDB
Map dimensions	360, 360, 360	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.83900005, 0.83900005, 0.83900005	Depositor

5 Model quality ⓘ

5.1 Standard geometry ⓘ

Bond lengths and bond angles in the following residue types are not validated in this section: UR3, 6MZ, LHH, ZN, 4AC, B8H, MA6, OMU, A2M, OMC, 5MC, MG, OMG

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	1	0.14	3/70601 (0.0%)	0.32	0/110066
2	2	0.18	0/34457	0.30	2/53715 (0.0%)
3	3	0.16	0/2995	0.34	0/4669
4	H	0.31	1/3120 (0.0%)	0.49	3/4209 (0.1%)
5	AA	0.13	0/1557	0.28	0/2087
6	AB	0.12	0/1602	0.26	0/2165
7	AC	0.11	0/463	0.27	0/628
8	AD	0.12	0/1476	0.23	0/1980
9	AE	0.13	0/2032	0.29	0/2742
10	AF	0.12	0/1838	0.25	0/2478
11	AG	0.12	0/993	0.29	0/1329
12	AH	0.21	0/1762	0.37	0/2366
13	AI	0.13	0/1055	0.25	0/1415
14	AJ	0.13	0/995	0.24	0/1327
15	AK	0.12	0/1089	0.26	0/1459
16	AL	0.11	0/817	0.28	0/1097
17	AM	0.14	0/973	0.28	0/1311
18	AN	0.12	0/1165	0.27	0/1547
19	AP	0.13	0/465	0.32	0/613
20	AQ	0.13	0/1333	0.26	0/1791
21	AR	0.13	0/907	0.27	0/1225
22	AS	0.12	0/548	0.28	0/725
23	AU	0.11	0/1253	0.26	0/1689
24	AV	0.11	0/826	0.24	0/1108
24	B6	0.15	0/808	0.35	0/1086
25	AW	0.12	0/476	0.27	0/641
26	AX	0.10	0/518	0.23	0/694
27	AY	0.09	0/412	0.25	0/549
28	AZ	0.10	0/1563	0.24	0/2099
29	A0	0.15	0/349	0.37	0/451
30	A3	0.11	0/945	0.31	0/1274
30	B4	0.17	0/945	0.36	0/1274

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
30	BG	0.18	0/944	0.33	0/1274
31	AT	0.14	0/1077	0.29	0/1439
32	AO	0.10	0/1135	0.28	0/1526
33	BB	0.12	0/1882	0.31	0/2538
34	BY	0.10	0/1254	0.26	0/1677
35	BO	0.17	0/1645	0.33	0/2209
36	BC	0.11	0/2933	0.30	0/3943
37	B5	0.14	0/619	0.30	0/830
37	BK	0.11	0/611	0.32	0/819
38	BL	0.13	0/1167	0.32	0/1552
39	Bf	0.11	0/443	0.26	0/589
40	BU	0.31	0/1027	0.45	1/1368 (0.1%)
41	Bb	0.10	0/1120	0.27	0/1494
42	Be	0.09	0/514	0.29	0/676
43	BE	0.21	0/1509	0.40	0/2022
44	Ba	0.10	0/793	0.28	0/1064
45	BT	0.11	0/705	0.28	0/945
46	BW	0.13	0/566	0.28	0/747
47	Bi	0.11	0/630	0.32	0/839
48	BI	0.08	0/1166	0.25	0/1559
49	BR	0.10	0/819	0.27	0/1098
50	BQ	0.09	0/1281	0.25	0/1689
51	BV	0.09	0/547	0.34	0/729
52	Bj	0.11	0/807	0.26	0/1070
53	BD	0.13	0/2069	0.31	0/2787
54	BF	0.15	0/1485	0.30	0/2003
55	BZ	0.14	0/759	0.32	0/1023
56	BP	0.09	0/984	0.25	0/1314
57	BM	0.11	0/1627	0.26	0/2170
58	BS	0.13	0/1275	0.32	0/1715
59	Bd	0.10	0/778	0.27	0/1036
60	BN	0.12	0/1483	0.30	0/1992
61	Bg	0.11	0/396	0.25	0/526
62	Bc	0.11	0/693	0.30	0/926
63	BJ	0.12	0/1080	0.30	0/1456
64	Bl	0.14	0/669	0.29	0/886
65	Bk	0.14	0/538	0.28	0/709
66	sd	0.24	0/223	0.49	0/347
All	All	0.15	4/179591 (0.0%)	0.31	6/264395 (0.0%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	H	377	ASP	C-N	14.46	1.54	1.33
1	1	94	G	O3'-P	-10.51	1.45	1.61
1	1	96	U	C1'-N1	5.73	1.57	1.48
1	1	2173	A2M	O3'-P	5.03	1.61	1.56

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	H	377	ASP	CA-C-N	-15.23	101.77	123.10
4	H	377	ASP	C-N-CA	-15.23	101.77	123.10
4	H	383	ARG	N-CA-C	6.30	120.58	113.01
40	BU	115	ILE	CB-CA-C	-5.83	104.27	112.14
2	2	1499	G	C2'-C3'-O3'	-5.44	105.54	113.70

There are no chirality outliers.

There are no planarity outliers.

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	1	65181	0	32950	1072	0
2	2	32311	0	16352	422	0
3	3	2703	0	1375	44	0
4	H	3074	0	3190	85	0
5	AA	1531	0	1623	28	0
6	AB	1571	0	1630	19	0
7	AC	449	0	435	8	0
8	AD	1452	0	1521	10	0
9	AE	1983	0	2060	28	0
10	AF	1808	0	1879	27	0
11	AG	977	0	1037	21	0
12	AH	1725	0	1780	35	0
13	AI	1034	0	1069	15	0
14	AJ	986	0	1070	9	0
15	AK	1073	0	1133	15	0
16	AL	809	0	859	17	0
17	AM	955	0	981	26	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
18	AN	1148	0	1248	14	0
19	AP	455	0	475	18	0
20	AQ	1305	0	1388	18	0
21	AR	884	0	906	11	0
22	AS	541	0	573	10	0
23	AU	1223	0	1263	18	0
24	AV	808	0	832	9	0
24	B6	790	0	806	16	0
25	AW	470	0	495	7	0
26	AX	516	0	544	9	0
27	AY	400	0	401	8	0
28	AZ	1541	0	1624	20	0
29	A0	343	0	407	7	0
30	A3	933	0	982	25	0
30	B4	933	0	982	30	0
30	BG	932	0	982	9	0
31	AT	1057	0	1131	26	0
32	AO	1116	0	1152	20	0
33	BB	1838	0	1903	30	0
34	BY	1235	0	1314	15	0
35	BO	1607	0	1638	30	0
36	BC	2870	0	3020	30	0
37	B5	614	0	671	13	0
37	BK	607	0	663	10	0
38	BL	1149	0	1218	23	0
39	Bf	435	0	494	9	0
40	BU	1011	0	1078	26	0
41	Bb	1095	0	1190	8	0
42	Be	503	0	533	12	0
43	BE	1485	0	1539	41	0
44	Ba	778	0	851	14	0
45	BT	696	0	758	7	0
46	BW	565	0	628	6	0
47	Bi	620	0	668	3	0
48	BI	1148	0	1237	15	0
49	BR	797	0	835	12	0
50	BQ	1265	0	1392	26	0
51	BV	532	0	523	6	0
52	Bj	789	0	842	7	0
53	BD	2026	0	2131	34	0
54	BF	1456	0	1514	20	0
55	BZ	749	0	796	14	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
56	BP	971	0	1039	10	0
57	BM	1588	0	1683	20	0
58	BS	1247	0	1280	10	0
59	Bd	759	0	825	6	0
60	BN	1452	0	1491	29	0
61	Bg	387	0	397	6	0
62	Bc	684	0	748	10	0
63	BJ	1066	0	1133	11	0
64	Bl	659	0	702	12	0
65	Bk	528	0	575	9	0
66	sd	199	0	98	7	0
67	1	166	0	0	0	0
67	2	69	0	0	0	0
67	3	1	0	0	0	0
67	AK	1	0	0	0	0
67	AN	1	0	0	0	0
67	BB	2	0	0	0	0
67	BD	1	0	0	0	0
67	BM	1	0	0	0	0
67	BS	1	0	0	0	0
67	Bb	1	0	0	0	0
68	AC	2	0	0	0	0
68	AF	1	0	0	0	0
68	AP	1	0	0	0	0
68	AR	1	0	0	0	0
68	AW	1	0	0	0	0
68	BV	1	0	0	0	0
68	Bd	1	0	0	0	0
68	Be	1	0	0	0	0
68	Bg	1	0	0	0	0
68	Bi	1	0	0	0	0
68	Bj	1	0	0	0	0
68	Bk	1	0	0	0	0
All	All	170684	0	124542	2351	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 8.

The worst 5 of 2351 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)
1:1:107:G:N2	1:1:112:A:N7	1.85	1.20
1:1:107:G:N2	1:1:112:A:C8	2.14	1.15
1:1:117:G:O2'	1:1:118:U:OP1	1.67	1.12
1:1:100:G:N2	1:1:119:C:O2	1.82	1.12
1:1:2887:A:HO2'	63:BJ:2:ALA:N	1.50	1.07

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
4	H	384/392 (98%)	382 (100%)	2 (0%)	0	100	100
5	AA	186/199 (94%)	181 (97%)	5 (3%)	0	100	100
6	AB	194/202 (96%)	193 (100%)	1 (0%)	0	100	100
7	AC	55/63 (87%)	53 (96%)	2 (4%)	0	100	100
8	AD	171/180 (95%)	169 (99%)	2 (1%)	0	100	100
9	AE	240/243 (99%)	236 (98%)	4 (2%)	0	100	100
10	AF	227/236 (96%)	222 (98%)	5 (2%)	0	100	100
11	AG	122/125 (98%)	117 (96%)	5 (4%)	0	100	100
12	AH	212/215 (99%)	210 (99%)	2 (1%)	0	100	100
13	AI	127/130 (98%)	125 (98%)	2 (2%)	0	100	100
14	AJ	123/127 (97%)	122 (99%)	1 (1%)	0	100	100
15	AK	133/135 (98%)	129 (97%)	4 (3%)	0	100	100
16	AL	98/102 (96%)	97 (99%)	1 (1%)	0	100	100
17	AM	125/137 (91%)	122 (98%)	3 (2%)	0	100	100
18	AN	144/147 (98%)	143 (99%)	1 (1%)	0	100	100
19	AP	53/56 (95%)	53 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
20	AQ	155/158 (98%)	152 (98%)	3 (2%)	0	100	100
21	AR	105/113 (93%)	101 (96%)	4 (4%)	0	100	100
22	AS	62/67 (92%)	58 (94%)	4 (6%)	0	100	100
23	AU	147/150 (98%)	147 (100%)	0	0	100	100
24	AV	94/99 (95%)	92 (98%)	2 (2%)	0	100	100
24	B6	92/99 (93%)	91 (99%)	1 (1%)	0	100	100
25	AW	59/65 (91%)	58 (98%)	1 (2%)	0	100	100
26	AX	63/71 (89%)	63 (100%)	0	0	100	100
27	AY	47/51 (92%)	41 (87%)	6 (13%)	0	100	100
28	AZ	194/210 (92%)	182 (94%)	12 (6%)	0	100	100
29	A0	34/37 (92%)	33 (97%)	1 (3%)	0	100	100
30	A3	120/123 (98%)	113 (94%)	7 (6%)	0	100	100
30	B4	120/123 (98%)	116 (97%)	4 (3%)	0	100	100
30	BG	120/123 (98%)	120 (100%)	0	0	100	100
31	AT	128/132 (97%)	127 (99%)	1 (1%)	0	100	100
32	AO	136/148 (92%)	128 (94%)	8 (6%)	0	100	100
33	BB	236/239 (99%)	230 (98%)	6 (2%)	0	100	100
34	BY	152/155 (98%)	152 (100%)	0	0	100	100
35	BO	196/203 (97%)	195 (100%)	1 (0%)	0	100	100
36	BC	358/361 (99%)	349 (98%)	9 (2%)	0	100	100
37	B5	79/82 (96%)	79 (100%)	0	0	100	100
37	BK	78/82 (95%)	75 (96%)	3 (4%)	0	100	100
38	BL	145/147 (99%)	139 (96%)	6 (4%)	0	100	100
39	Bf	48/51 (94%)	46 (96%)	2 (4%)	0	100	100
40	BU	119/121 (98%)	118 (99%)	1 (1%)	0	100	100
41	Bb	127/130 (98%)	127 (100%)	0	0	100	100
42	Be	59/62 (95%)	58 (98%)	1 (2%)	0	100	100
43	BE	183/188 (97%)	183 (100%)	0	0	100	100
44	Ba	92/94 (98%)	92 (100%)	0	0	100	100
45	BT	84/86 (98%)	83 (99%)	1 (1%)	0	100	100
46	BW	66/68 (97%)	66 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
47	Bi	80/83 (96%)	77 (96%)	3 (4%)	0	100	100
48	BI	140/142 (99%)	137 (98%)	3 (2%)	0	100	100
49	BR	94/97 (97%)	94 (100%)	0	0	100	100
50	BQ	148/151 (98%)	148 (100%)	0	0	100	100
51	BV	61/67 (91%)	60 (98%)	1 (2%)	0	100	100
52	Bj	92/94 (98%)	91 (99%)	1 (1%)	0	100	100
53	BD	253/255 (99%)	248 (98%)	5 (2%)	0	100	100
54	BF	181/184 (98%)	180 (99%)	1 (1%)	0	100	100
55	BZ	96/99 (97%)	95 (99%)	1 (1%)	0	100	100
56	BP	118/120 (98%)	118 (100%)	0	0	100	100
57	BM	191/194 (98%)	188 (98%)	3 (2%)	0	100	100
58	BS	152/155 (98%)	151 (99%)	1 (1%)	0	100	100
59	Bd	89/91 (98%)	89 (100%)	0	0	100	100
60	BN	176/181 (97%)	169 (96%)	7 (4%)	0	100	100
61	Bg	46/51 (90%)	46 (100%)	0	0	100	100
62	Bc	85/87 (98%)	84 (99%)	1 (1%)	0	100	100
63	BJ	138/141 (98%)	138 (100%)	0	0	100	100
64	Bl	75/77 (97%)	75 (100%)	0	0	100	100
65	Bk	61/65 (94%)	60 (98%)	1 (2%)	0	100	100
All	All	8568/8861 (97%)	8416 (98%)	152 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	
4	H	332/338 (98%)	306 (92%)	26 (8%)	11	13
5	AA	160/167 (96%)	151 (94%)	9 (6%)	19	24

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
6	AB	168/173 (97%)	163 (97%)	5 (3%)	36	49
7	AC	50/55 (91%)	50 (100%)	0	100	100
8	AD	156/160 (98%)	153 (98%)	3 (2%)	50	66
9	AE	213/214 (100%)	210 (99%)	3 (1%)	59	75
10	AF	192/198 (97%)	190 (99%)	2 (1%)	68	81
11	AG	107/108 (99%)	102 (95%)	5 (5%)	23	31
12	AH	183/184 (100%)	166 (91%)	17 (9%)	8	9
13	AI	106/107 (99%)	105 (99%)	1 (1%)	70	84
14	AJ	101/103 (98%)	97 (96%)	4 (4%)	28	38
15	AK	111/111 (100%)	109 (98%)	2 (2%)	51	68
16	AL	89/91 (98%)	84 (94%)	5 (6%)	19	24
17	AM	94/104 (90%)	91 (97%)	3 (3%)	34	47
18	AN	120/121 (99%)	119 (99%)	1 (1%)	73	85
19	AP	45/46 (98%)	41 (91%)	4 (9%)	9	10
20	AQ	142/143 (99%)	139 (98%)	3 (2%)	47	63
21	AR	96/102 (94%)	94 (98%)	2 (2%)	47	63
22	AS	58/61 (95%)	54 (93%)	4 (7%)	14	16
23	AU	126/127 (99%)	122 (97%)	4 (3%)	34	47
24	AV	88/90 (98%)	87 (99%)	1 (1%)	65	79
24	B6	86/90 (96%)	78 (91%)	8 (9%)	8	9
25	AW	53/56 (95%)	53 (100%)	0	100	100
26	AX	55/60 (92%)	49 (89%)	6 (11%)	6	6
27	AY	40/42 (95%)	38 (95%)	2 (5%)	22	28
28	AZ	155/168 (92%)	147 (95%)	8 (5%)	21	26
29	A0	34/35 (97%)	33 (97%)	1 (3%)	37	51
30	A3	98/99 (99%)	97 (99%)	1 (1%)	68	81
30	B4	98/99 (99%)	93 (95%)	5 (5%)	21	27
30	BG	98/99 (99%)	88 (90%)	10 (10%)	7	7
31	AT	113/114 (99%)	108 (96%)	5 (4%)	25	34
32	AO	115/123 (94%)	111 (96%)	4 (4%)	32	43
33	BB	188/189 (100%)	185 (98%)	3 (2%)	55	71

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
34	BY	132/133 (99%)	128 (97%)	4 (3%)	36	49
35	BO	167/170 (98%)	156 (93%)	11 (7%)	15	18
36	BC	306/307 (100%)	302 (99%)	4 (1%)	61	76
37	B5	65/66 (98%)	62 (95%)	3 (5%)	24	32
37	BK	64/66 (97%)	59 (92%)	5 (8%)	11	13
38	BL	118/118 (100%)	110 (93%)	8 (7%)	14	17
39	Bf	46/47 (98%)	44 (96%)	2 (4%)	26	35
40	BU	109/109 (100%)	105 (96%)	4 (4%)	30	41
41	Bb	117/118 (99%)	115 (98%)	2 (2%)	53	69
42	Be	52/53 (98%)	52 (100%)	0	100	100
43	BE	158/160 (99%)	134 (85%)	24 (15%)	3	2
44	Ba	84/84 (100%)	82 (98%)	2 (2%)	43	58
45	BT	77/77 (100%)	76 (99%)	1 (1%)	61	76
46	BW	63/63 (100%)	60 (95%)	3 (5%)	23	30
47	Bi	61/62 (98%)	59 (97%)	2 (3%)	33	45
48	BI	122/122 (100%)	122 (100%)	0	100	100
49	BR	86/87 (99%)	86 (100%)	0	100	100
50	BQ	132/133 (99%)	132 (100%)	0	100	100
51	BV	54/58 (93%)	53 (98%)	1 (2%)	50	66
52	Bj	83/84 (99%)	83 (100%)	0	100	100
53	BD	212/212 (100%)	208 (98%)	4 (2%)	50	66
54	BF	156/157 (99%)	145 (93%)	11 (7%)	13	16
55	BZ	79/80 (99%)	72 (91%)	7 (9%)	9	10
56	BP	102/102 (100%)	99 (97%)	3 (3%)	37	51
57	BM	160/161 (99%)	157 (98%)	3 (2%)	50	66
58	BS	130/131 (99%)	128 (98%)	2 (2%)	57	73
59	Bd	82/82 (100%)	80 (98%)	2 (2%)	43	58
60	BN	149/151 (99%)	145 (97%)	4 (3%)	39	53
61	Bg	37/39 (95%)	37 (100%)	0	100	100
62	Bc	74/74 (100%)	71 (96%)	3 (4%)	27	37
63	BJ	108/109 (99%)	105 (97%)	3 (3%)	38	52

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
64	Bl	72/72 (100%)	70 (97%)	2 (3%)	38 52
65	Bk	55/57 (96%)	52 (94%)	3 (6%)	19 24
All	All	7382/7521 (98%)	7102 (96%)	280 (4%)	30 40

5 of 280 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
54	BF	136	GLU
55	BZ	55	LEU
60	BN	103	ARG
22	AS	61	LYS
21	AR	48	THR

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 166 such sidechains are listed below:

Mol	Chain	Res	Type
49	BR	89	HIS
57	BM	174	ASN
50	BQ	75	HIS
54	BF	74	HIS
60	BN	59	GLN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	1	3016/3018 (99%)	472 (15%)	25 (0%)
2	2	1496/1512 (98%)	165 (11%)	3 (0%)
3	3	125/128 (97%)	16 (12%)	0
66	sd	8/9 (88%)	1 (12%)	0
All	All	4645/4667 (99%)	654 (14%)	28 (0%)

5 of 654 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	1	12	G
1	1	28	A
1	1	39	A
1	1	43	A
1	1	51	A

5 of 28 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	1	1288	U
2	2	1106	C
1	1	1364	C
1	1	3021	U
1	1	1363	U

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

159 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	$\# Z > 2$
1	4AC	1	60	1	21,24,25	1.01	2 (9%)	29,34,37	1.43	5 (17%)
1	4AC	1	694	1	21,24,25	1.02	1 (4%)	29,34,37	1.37	4 (13%)
2	4AC	2	319	2	21,24,25	1.01	2 (9%)	29,34,37	1.39	4 (13%)
2	OMG	2	680	2	23,26,27	1.21	3 (13%)	33,38,41	1.96	6 (18%)
1	4AC	1	2865	1	21,24,25	1.03	1 (4%)	29,34,37	1.39	4 (13%)
2	4AC	2	53	2	21,24,25	1.02	2 (9%)	29,34,37	1.38	4 (13%)
2	4AC	2	394	2	21,24,25	1.00	2 (9%)	29,34,37	1.38	4 (13%)
2	LHH	2	250	2	22,25,26	0.30	0	29,35,38	0.49	0
1	4AC	1	2001	1	21,24,25	1.07	2 (9%)	29,34,37	1.50	4 (13%)
2	OMU	2	1380	2	19,22,23	1.28	4 (21%)	26,31,34	1.71	4 (15%)
2	UR3	2	1467	2	19,22,23	0.97	0	26,32,35	1.47	2 (7%)
2	4AC	2	839	2	21,24,25	1.03	2 (9%)	29,34,37	1.53	5 (17%)
1	4AC	1	1550	1	21,24,25	1.02	1 (4%)	29,34,37	1.39	4 (13%)
1	LHH	1	616	1	22,25,26	0.30	0	29,35,38	0.46	0
2	5MC	2	535	2	18,22,23	0.93	2 (11%)	26,32,35	1.15	2 (7%)
2	4AC	2	636	2	21,24,25	1.01	2 (9%)	29,34,37	1.48	4 (13%)
1	4AC	1	2083	1	21,24,25	1.04	2 (9%)	29,34,37	1.39	5 (17%)
2	OMU	2	1177	2	19,22,23	1.22	3 (15%)	26,31,34	1.69	5 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	OMG	2	657	2	23,26,27	1.23	3 (13%)	33,38,41	1.97	6 (18%)
2	4AC	2	17	2	21,24,25	1.03	2 (9%)	29,34,37	1.36	4 (13%)
1	OMU	1	2784	1	19,22,23	1.18	2 (10%)	26,31,34	1.70	5 (19%)
2	MA6	2	1487	2	23,26,27	1.50	4 (17%)	34,38,41	2.18	11 (32%)
1	OMG	1	789	1	23,26,27	1.25	3 (13%)	33,38,41	2.00	7 (21%)
1	4AC	1	1554	1	21,24,25	1.02	1 (4%)	29,34,37	1.41	4 (13%)
1	4AC	1	1222	1,67	21,24,25	1.02	1 (4%)	29,34,37	1.44	4 (13%)
1	4AC	1	1667	1	21,24,25	1.02	1 (4%)	29,34,37	1.37	4 (13%)
1	4AC	1	1765	1	21,24,25	1.04	2 (9%)	29,34,37	1.38	4 (13%)
1	4AC	1	2585	1	21,24,25	1.05	2 (9%)	29,34,37	1.79	6 (20%)
2	4AC	2	868	2	21,24,25	1.02	2 (9%)	29,34,37	1.36	4 (13%)
2	5MC	2	1374	2	18,22,23	0.93	2 (11%)	26,32,35	1.10	2 (7%)
1	4AC	1	2992	1	21,24,25	1.06	1 (4%)	29,34,37	0.86	2 (6%)
2	4AC	2	479	2	21,24,25	1.01	2 (9%)	29,34,37	1.39	4 (13%)
2	OMC	2	1040	2	19,22,23	0.84	0	26,31,34	0.96	1 (3%)
1	OMC	1	1948	1	19,22,23	0.81	0	26,31,34	0.74	0
1	4AC	1	2027	1	21,24,25	1.03	2 (9%)	29,34,37	1.38	4 (13%)
2	OMG	2	913	2	23,26,27	1.19	3 (13%)	33,38,41	1.92	6 (18%)
1	5MC	1	2093	1	18,22,23	0.94	2 (11%)	26,32,35	1.12	2 (7%)
2	OMG	2	471	2	23,26,27	1.22	3 (13%)	33,38,41	1.94	6 (18%)
2	OMG	2	467	2	23,26,27	1.19	3 (13%)	33,38,41	1.94	6 (18%)
1	4AC	1	1617	1	21,24,25	1.02	1 (4%)	29,34,37	1.40	4 (13%)
1	4AC	1	1243	1	21,24,25	1.05	2 (9%)	29,34,37	1.44	4 (13%)
1	4AC	1	161	1	21,24,25	1.03	2 (9%)	29,34,37	1.40	4 (13%)
1	4AC	1	357	1	21,24,25	1.03	2 (9%)	29,34,37	1.39	4 (13%)
1	4AC	1	1885	1	21,24,25	1.01	2 (9%)	29,34,37	1.39	4 (13%)
2	4AC	2	1184	2	21,24,25	1.05	2 (9%)	29,34,37	1.56	4 (13%)
2	4AC	2	303	2	21,24,25	1.02	2 (9%)	29,34,37	1.37	4 (13%)
2	5MC	2	1202	2	18,22,23	0.98	2 (11%)	26,32,35	1.25	3 (11%)
1	4AC	1	22	1	21,24,25	1.04	2 (9%)	29,34,37	1.40	4 (13%)
1	4AC	1	2966	1	21,24,25	1.04	2 (9%)	29,34,37	1.49	4 (13%)
2	OMU	2	64	2	19,22,23	1.22	3 (15%)	26,31,34	1.70	5 (19%)
2	5MC	2	875	2	18,22,23	0.94	2 (11%)	26,32,35	1.17	3 (11%)
2	OMU	2	787	2	19,22,23	1.23	2 (10%)	26,31,34	1.80	5 (19%)
1	4AC	1	829	1	21,24,25	1.03	2 (9%)	29,34,37	1.37	4 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	4AC	2	851	2	21,24,25	1.03	2 (9%)	29,34,37	1.36	4 (13%)
2	4AC	2	286	2	21,24,25	1.03	2 (9%)	29,34,37	1.34	4 (13%)
1	4AC	1	1878	1	21,24,25	1.01	1 (4%)	29,34,37	1.53	4 (13%)
1	4AC	1	2548	1	21,24,25	1.02	2 (9%)	29,34,37	1.50	4 (13%)
1	4AC	1	922	1	21,24,25	1.03	2 (9%)	29,34,37	1.49	4 (13%)
2	4AC	2	1239	2	21,24,25	1.02	1 (4%)	29,34,37	1.35	4 (13%)
1	4AC	1	91	1	21,24,25	1.04	2 (9%)	29,34,37	1.58	4 (13%)
2	B8H	2	938	2	19,22,23	0.63	1 (5%)	22,32,35	0.71	1 (4%)
2	4AC	2	718	2	21,24,25	1.00	2 (9%)	29,34,37	1.39	4 (13%)
1	4AC	1	451	1	21,24,25	1.05	2 (9%)	29,34,37	1.47	4 (13%)
1	4AC	1	2608	1	21,24,25	1.07	2 (9%)	29,34,37	1.76	6 (20%)
1	4AC	1	314	1	21,24,25	1.03	1 (4%)	29,34,37	1.39	4 (13%)
2	4AC	2	731	2	21,24,25	1.02	2 (9%)	29,34,37	1.38	4 (13%)
2	5MC	2	1505	66,2	18,22,23	0.98	2 (11%)	26,32,35	1.17	3 (11%)
1	OMG	1	2144	1	23,26,27	1.21	3 (13%)	33,38,41	1.95	6 (18%)
1	4AC	1	1867	1	21,24,25	1.04	2 (9%)	29,34,37	1.37	4 (13%)
2	4AC	2	546	2	21,24,25	0.99	2 (9%)	29,34,37	1.38	4 (13%)
1	A2M	1	2173	1	22,25,26	1.46	4 (18%)	31,36,39	2.02	8 (25%)
1	4AC	1	2329	1	21,24,25	1.03	2 (9%)	29,34,37	1.39	4 (13%)
1	5MC	1	2203	1	18,22,23	0.95	2 (11%)	26,32,35	1.20	2 (7%)
1	4AC	1	3004	1	21,24,25	1.03	2 (9%)	29,34,37	1.35	4 (13%)
2	4AC	2	848	2	21,24,25	1.01	2 (9%)	29,34,37	1.37	4 (13%)
1	4AC	1	211	1	21,24,25	1.03	2 (9%)	29,34,37	1.41	4 (13%)
1	4AC	1	2570	1	21,24,25	1.02	1 (4%)	29,34,37	1.37	4 (13%)
3	4AC	3	32	3	21,24,25	1.04	1 (4%)	29,34,37	1.40	4 (13%)
1	4AC	1	136	1	21,24,25	1.03	1 (4%)	29,34,37	1.40	4 (13%)
1	4AC	1	2925	1	21,24,25	1.02	1 (4%)	29,34,37	1.42	4 (13%)
2	OMG	2	1069	2	23,26,27	1.21	3 (13%)	33,38,41	1.93	6 (18%)
2	4AC	2	1479	2	21,24,25	1.00	2 (9%)	29,34,37	1.43	4 (13%)
2	4AC	2	751	2	21,24,25	1.03	2 (9%)	29,34,37	1.34	4 (13%)
2	OMC	2	773	2	19,22,23	0.82	0	26,31,34	0.83	0
1	4AC	1	2495	1	21,24,25	1.04	2 (9%)	29,34,37	1.37	4 (13%)
1	4AC	1	533	1	21,24,25	1.03	1 (4%)	29,34,37	1.40	4 (13%)
1	A2M	1	1055	1	22,25,26	1.47	4 (18%)	31,36,39	2.11	10 (32%)
1	5MC	1	2198	1	18,22,23	0.95	2 (11%)	26,32,35	1.11	2 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	5MC	1	2733	1	18,22,23	0.98	1 (5%)	26,32,35	1.34	4 (15%)
2	LHH	2	1041	2	22,25,26	0.31	0	29,35,38	0.49	0
1	4AC	1	2136	1	21,24,25	1.04	1 (4%)	29,34,37	1.46	5 (17%)
1	4AC	1	1065	1	21,24,25	1.01	1 (4%)	29,34,37	1.41	4 (13%)
2	4AC	2	1147	2	21,24,25	1.04	1 (4%)	29,34,37	1.54	5 (17%)
1	4AC	1	2545	1	21,24,25	1.05	2 (9%)	29,34,37	1.46	4 (13%)
1	OMG	1	2138	1	23,26,27	1.20	3 (13%)	33,38,41	1.91	6 (18%)
1	4AC	1	2249	1	21,24,25	1.04	2 (9%)	29,34,37	1.41	4 (13%)
1	4AC	1	1293	1	21,24,25	1.04	2 (9%)	29,34,37	1.35	4 (13%)
2	5MC	2	1496	2	18,22,23	0.94	2 (11%)	26,32,35	1.18	2 (7%)
1	4AC	1	2908	1	21,24,25	1.04	2 (9%)	29,34,37	1.37	4 (13%)
1	OMG	1	328	1	23,26,27	1.21	3 (13%)	33,38,41	1.91	6 (18%)
1	4AC	1	1460	1	21,24,25	1.02	2 (9%)	29,34,37	1.40	4 (13%)
1	4AC	1	2287	1	21,24,25	1.01	1 (4%)	29,34,37	1.43	4 (13%)
1	4AC	1	641	1	21,24,25	1.04	2 (9%)	29,34,37	1.37	4 (13%)
1	4AC	1	2718	1	21,24,25	1.01	1 (4%)	29,34,37	1.43	4 (13%)
2	OMC	2	846	2	19,22,23	0.82	0	26,31,34	0.79	0
1	4AC	1	1938	1	21,24,25	1.04	2 (9%)	29,34,37	1.38	4 (13%)
2	5MC	2	693	2	18,22,23	0.93	2 (11%)	26,32,35	1.10	2 (7%)
1	4AC	1	341	1	21,24,25	1.04	2 (9%)	29,34,37	1.37	4 (13%)
2	4AC	2	1028	2	21,24,25	1.01	2 (9%)	29,34,37	1.34	4 (13%)
1	OMG	1	2678	1	23,26,27	1.20	3 (13%)	33,38,41	1.92	6 (18%)
1	4AC	1	243	1	21,24,25	1.02	1 (4%)	29,34,37	1.39	4 (13%)
2	4AC	2	590	2	21,24,25	1.02	2 (9%)	29,34,37	1.38	4 (13%)
2	A2M	2	373	2	22,25,26	1.46	4 (18%)	31,36,39	2.11	9 (29%)
1	4AC	1	1048	1	21,24,25	1.02	2 (9%)	29,34,37	1.44	4 (13%)
1	4AC	1	1962	1	21,24,25	1.03	2 (9%)	29,34,37	1.35	4 (13%)
2	4AC	2	828	2	21,24,25	1.00	2 (9%)	29,34,37	1.37	4 (13%)
2	4AC	2	1233	2	21,24,25	1.02	1 (4%)	29,34,37	1.42	4 (13%)
2	4AC	2	511	2	21,24,25	1.04	2 (9%)	29,34,37	1.34	4 (13%)
1	4AC	1	1755	1	21,24,25	1.01	1 (4%)	29,34,37	1.67	4 (13%)
1	4AC	1	1780	1	21,24,25	1.02	1 (4%)	29,34,37	1.62	4 (13%)
2	4AC	2	957	2	21,24,25	1.02	2 (9%)	29,34,37	1.35	4 (13%)
1	5MC	1	2162	1	18,22,23	0.96	2 (11%)	26,32,35	1.23	2 (7%)
1	5MC	1	2183	1	18,22,23	0.93	2 (11%)	26,32,35	1.17	2 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	OMG	1	923	1,67	23,26,27	1.21	3 (13%)	33,38,41	1.94	6 (18%)
1	5MC	1	2163	1	18,22,23	0.94	2 (11%)	26,32,35	1.26	2 (7%)
1	4AC	1	835	1	21,24,25	1.03	1 (4%)	29,34,37	1.40	4 (13%)
1	OMG	1	2147	1	23,26,27	1.20	3 (13%)	33,38,41	1.93	6 (18%)
2	OMU	2	774	2	19,22,23	1.22	3 (15%)	26,31,34	1.73	4 (15%)
2	OMG	2	934	2	23,26,27	1.22	3 (13%)	33,38,41	1.94	6 (18%)
1	4AC	1	2960	1	21,24,25	1.02	2 (9%)	29,34,37	1.36	4 (13%)
1	4AC	1	2062	1	21,24,25	1.03	2 (9%)	29,34,37	1.38	4 (13%)
2	5MC	2	1498	2	18,22,23	0.94	2 (11%)	26,32,35	1.25	4 (15%)
2	MA6	2	1488	2	23,26,27	1.45	4 (17%)	34,38,41	2.11	10 (29%)
1	4AC	1	1934	1	21,24,25	1.01	2 (9%)	29,34,37	1.39	4 (13%)
2	OMC	2	1376	2	19,22,23	0.81	0	26,31,34	0.76	0
1	4AC	1	1068	1	21,24,25	1.05	2 (9%)	29,34,37	1.54	4 (13%)
1	4AC	1	981	1	21,24,25	1.01	2 (9%)	29,34,37	1.41	4 (13%)
1	4AC	1	2642	1	21,24,25	1.03	2 (9%)	29,34,37	1.35	4 (13%)
1	4AC	1	1662	1	21,24,25	1.03	2 (9%)	29,34,37	1.40	4 (13%)
2	6MZ	2	1469	2,67	22,25,26	1.48	4 (18%)	30,36,39	2.22	10 (33%)
1	4AC	1	802	1	21,24,25	1.02	1 (4%)	29,34,37	1.45	4 (13%)
2	OMU	2	830	2	19,22,23	1.27	4 (21%)	26,31,34	1.69	5 (19%)
1	4AC	1	229	1	21,24,25	1.04	2 (9%)	29,34,37	1.40	4 (13%)
1	OMC	1	615	1	19,22,23	0.80	0	26,31,34	0.80	0
2	OMG	2	519	2	23,26,27	1.19	3 (13%)	33,38,41	1.96	6 (18%)
2	OMC	2	129	2	19,22,23	0.82	0	26,31,34	0.80	0
1	4AC	1	2937	1,67	21,24,25	1.02	2 (9%)	29,34,37	1.61	4 (13%)
2	4AC	2	379	2	21,24,25	1.01	2 (9%)	29,34,37	1.39	4 (13%)
1	4AC	1	1557	1	21,24,25	1.03	2 (9%)	29,34,37	1.34	4 (13%)
1	4AC	1	1621	1	21,24,25	1.03	1 (4%)	29,34,37	1.40	4 (13%)
2	OMG	2	873	2	23,26,27	1.21	3 (13%)	33,38,41	1.96	6 (18%)
1	4AC	1	1594	1,67	21,24,25	1.01	1 (4%)	29,34,37	1.43	4 (13%)
1	4AC	1	1873	1	21,24,25	1.02	1 (4%)	29,34,37	1.39	4 (13%)
2	4AC	2	626	2	21,24,25	1.02	2 (9%)	29,34,37	1.40	4 (13%)
2	4AC	2	703	2	21,24,25	1.01	2 (9%)	29,34,37	1.43	4 (13%)
2	5MC	2	1025	2	18,22,23	0.95	2 (11%)	26,32,35	1.11	2 (7%)
2	OMU	2	20	2	19,22,23	1.23	2 (10%)	26,31,34	1.81	5 (19%)
1	4AC	1	458	1	21,24,25	1.00	1 (4%)	29,34,37	1.40	4 (13%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	4AC	1	1695	1	21,24,25	1.03	2 (9%)	29,34,37	1.37	4 (13%)

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
1	4AC	1	60	1	-	0/11/29/30	0/2/2/2
1	4AC	1	694	1	-	0/11/29/30	0/2/2/2
2	4AC	2	319	2	-	0/11/29/30	0/2/2/2
2	OMG	2	680	2	-	0/9/27/28	0/3/3/3
1	4AC	1	2865	1	-	0/11/29/30	0/2/2/2
2	4AC	2	53	2	-	0/11/29/30	0/2/2/2
2	4AC	2	394	2	-	0/11/29/30	0/2/2/2
2	LHH	2	250	2	-	0/13/31/32	0/2/2/2
1	4AC	1	2001	1	-	0/11/29/30	0/2/2/2
2	OMU	2	1380	2	-	0/9/27/28	0/2/2/2
2	UR3	2	1467	2	-	0/7/25/26	0/2/2/2
2	4AC	2	839	2	-	0/11/29/30	0/2/2/2
1	4AC	1	1550	1	-	0/11/29/30	0/2/2/2
1	LHH	1	616	1	-	0/13/31/32	0/2/2/2
2	5MC	2	535	2	-	0/7/25/26	0/2/2/2
2	4AC	2	636	2	-	0/11/29/30	0/2/2/2
1	4AC	1	2083	1	-	0/11/29/30	0/2/2/2
2	OMU	2	1177	2	-	0/9/27/28	0/2/2/2
2	OMG	2	657	2	-	0/9/27/28	0/3/3/3
2	4AC	2	17	2	-	0/11/29/30	0/2/2/2
1	OMU	1	2784	1	-	0/9/27/28	0/2/2/2
2	MA6	2	1487	2	-	0/11/29/30	0/3/3/3
1	OMG	1	789	1	-	2/9/27/28	0/3/3/3
1	4AC	1	1554	1	-	0/11/29/30	0/2/2/2
1	4AC	1	1222	1,67	-	0/11/29/30	0/2/2/2
1	4AC	1	1667	1	-	0/11/29/30	0/2/2/2
1	4AC	1	1765	1	-	0/11/29/30	0/2/2/2
1	4AC	1	2585	1	-	3/11/29/30	0/2/2/2
2	4AC	2	868	2	-	0/11/29/30	0/2/2/2
2	5MC	2	1374	2	-	1/7/25/26	0/2/2/2
1	4AC	1	2992	1	-	2/11/29/30	0/2/2/2
2	4AC	2	479	2	-	0/11/29/30	0/2/2/2
2	OMC	2	1040	2	-	1/9/27/28	0/2/2/2
1	OMC	1	1948	1	-	0/9/27/28	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	4AC	1	2027	1	-	0/11/29/30	0/2/2/2
2	OMG	2	913	2	-	0/9/27/28	0/3/3/3
1	5MC	1	2093	1	-	0/7/25/26	0/2/2/2
2	OMG	2	471	2	-	0/9/27/28	0/3/3/3
2	OMG	2	467	2	-	0/9/27/28	0/3/3/3
1	4AC	1	1617	1	-	0/11/29/30	0/2/2/2
1	4AC	1	1243	1	-	0/11/29/30	0/2/2/2
1	4AC	1	161	1	-	0/11/29/30	0/2/2/2
1	4AC	1	357	1	-	0/11/29/30	0/2/2/2
1	4AC	1	1885	1	-	0/11/29/30	0/2/2/2
2	4AC	2	1184	2	-	0/11/29/30	0/2/2/2
2	4AC	2	303	2	-	0/11/29/30	0/2/2/2
2	5MC	2	1202	2	-	0/7/25/26	0/2/2/2
1	4AC	1	22	1	-	0/11/29/30	0/2/2/2
1	4AC	1	2966	1	-	0/11/29/30	0/2/2/2
2	OMU	2	64	2	-	0/9/27/28	0/2/2/2
2	5MC	2	875	2	-	0/7/25/26	0/2/2/2
2	OMU	2	787	2	-	4/9/27/28	0/2/2/2
1	4AC	1	829	1	-	0/11/29/30	0/2/2/2
2	4AC	2	851	2	-	0/11/29/30	0/2/2/2
2	4AC	2	286	2	-	0/11/29/30	0/2/2/2
1	4AC	1	1878	1	-	0/11/29/30	0/2/2/2
1	4AC	1	2548	1	-	0/11/29/30	0/2/2/2
1	4AC	1	922	1	-	0/11/29/30	0/2/2/2
2	4AC	2	1239	2	-	0/11/29/30	0/2/2/2
1	4AC	1	91	1	-	0/11/29/30	0/2/2/2
2	B8H	2	938	2	-	1/7/25/26	0/2/2/2
2	4AC	2	718	2	-	0/11/29/30	0/2/2/2
1	4AC	1	451	1	-	0/11/29/30	0/2/2/2
1	4AC	1	2608	1	-	3/11/29/30	0/2/2/2
1	4AC	1	314	1	-	0/11/29/30	0/2/2/2
2	4AC	2	731	2	-	0/11/29/30	0/2/2/2
2	5MC	2	1505	66,2	-	0/7/25/26	0/2/2/2
1	OMG	1	2144	1	-	2/9/27/28	0/3/3/3
1	4AC	1	1867	1	-	0/11/29/30	0/2/2/2
2	4AC	2	546	2	-	0/11/29/30	0/2/2/2
1	A2M	1	2173	1	-	0/9/27/28	0/3/3/3
1	4AC	1	2329	1	-	0/11/29/30	0/2/2/2
1	5MC	1	2203	1	-	0/7/25/26	0/2/2/2
1	4AC	1	3004	1	-	0/11/29/30	0/2/2/2
2	4AC	2	848	2	-	0/11/29/30	0/2/2/2
1	4AC	1	211	1	-	0/11/29/30	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	4AC	1	2570	1	-	0/11/29/30	0/2/2/2
3	4AC	3	32	3	-	1/11/29/30	0/2/2/2
1	4AC	1	136	1	-	0/11/29/30	0/2/2/2
1	4AC	1	2925	1	-	0/11/29/30	0/2/2/2
2	OMG	2	1069	2	-	2/9/27/28	0/3/3/3
2	4AC	2	1479	2	-	0/11/29/30	0/2/2/2
2	4AC	2	751	2	-	0/11/29/30	0/2/2/2
2	OMC	2	773	2	-	0/9/27/28	0/2/2/2
1	4AC	1	2495	1	-	0/11/29/30	0/2/2/2
1	4AC	1	533	1	-	0/11/29/30	0/2/2/2
1	A2M	1	1055	1	-	0/9/27/28	0/3/3/3
1	5MC	1	2198	1	-	0/7/25/26	0/2/2/2
1	5MC	1	2733	1	-	6/7/25/26	0/2/2/2
2	LHH	2	1041	2	-	0/13/31/32	0/2/2/2
1	4AC	1	2136	1	-	0/11/29/30	0/2/2/2
1	4AC	1	1065	1	-	0/11/29/30	0/2/2/2
2	4AC	2	1147	2	-	0/11/29/30	0/2/2/2
1	4AC	1	2545	1	-	0/11/29/30	0/2/2/2
1	OMG	1	2138	1	-	0/9/27/28	0/3/3/3
1	4AC	1	2249	1	-	0/11/29/30	0/2/2/2
1	4AC	1	1293	1	-	0/11/29/30	0/2/2/2
2	5MC	2	1496	2	-	0/7/25/26	0/2/2/2
1	4AC	1	2908	1	-	0/11/29/30	0/2/2/2
1	OMG	1	328	1	-	0/9/27/28	0/3/3/3
1	4AC	1	1460	1	-	0/11/29/30	0/2/2/2
1	4AC	1	2287	1	-	0/11/29/30	0/2/2/2
1	4AC	1	641	1	-	0/11/29/30	0/2/2/2
1	4AC	1	2718	1	-	0/11/29/30	0/2/2/2
2	OMC	2	846	2	-	0/9/27/28	0/2/2/2
1	4AC	1	1938	1	-	0/11/29/30	0/2/2/2
2	5MC	2	693	2	-	0/7/25/26	0/2/2/2
1	4AC	1	341	1	-	0/11/29/30	0/2/2/2
2	4AC	2	1028	2	-	0/11/29/30	0/2/2/2
1	OMG	1	2678	1	-	1/9/27/28	0/3/3/3
1	4AC	1	243	1	-	0/11/29/30	0/2/2/2
2	4AC	2	590	2	-	0/11/29/30	0/2/2/2
2	A2M	2	373	2	-	0/9/27/28	0/3/3/3
1	4AC	1	1048	1	-	0/11/29/30	0/2/2/2
1	4AC	1	1962	1	-	0/11/29/30	0/2/2/2
2	4AC	2	828	2	-	0/11/29/30	0/2/2/2
2	4AC	2	1233	2	-	0/11/29/30	0/2/2/2
2	4AC	2	511	2	-	0/11/29/30	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	4AC	1	1755	1	-	1/11/29/30	0/2/2/2
1	4AC	1	1780	1	-	0/11/29/30	0/2/2/2
2	4AC	2	957	2	-	0/11/29/30	0/2/2/2
1	5MC	1	2162	1	-	0/7/25/26	0/2/2/2
1	5MC	1	2183	1	-	0/7/25/26	0/2/2/2
1	OMG	1	923	1,67	-	2/9/27/28	0/3/3/3
1	5MC	1	2163	1	-	0/7/25/26	0/2/2/2
1	4AC	1	835	1	-	0/11/29/30	0/2/2/2
1	OMG	1	2147	1	-	2/9/27/28	0/3/3/3
2	OMU	2	774	2	-	0/9/27/28	0/2/2/2
2	OMG	2	934	2	-	1/9/27/28	0/3/3/3
1	4AC	1	2960	1	-	0/11/29/30	0/2/2/2
1	4AC	1	2062	1	-	0/11/29/30	0/2/2/2
2	5MC	2	1498	2	-	3/7/25/26	0/2/2/2
2	MA6	2	1488	2	-	2/11/29/30	0/3/3/3
1	4AC	1	1934	1	-	0/11/29/30	0/2/2/2
2	OMC	2	1376	2	-	0/9/27/28	0/2/2/2
1	4AC	1	1068	1	-	0/11/29/30	0/2/2/2
1	4AC	1	981	1	-	0/11/29/30	0/2/2/2
1	4AC	1	2642	1	-	0/11/29/30	0/2/2/2
1	4AC	1	1662	1	-	0/11/29/30	0/2/2/2
2	6MZ	2	1469	2,67	-	0/9/27/28	0/3/3/3
1	4AC	1	802	1	-	0/11/29/30	0/2/2/2
2	OMU	2	830	2	-	1/9/27/28	0/2/2/2
1	4AC	1	229	1	-	0/11/29/30	0/2/2/2
1	OMC	1	615	1	-	0/9/27/28	0/2/2/2
2	OMG	2	519	2	-	0/9/27/28	0/3/3/3
2	OMC	2	129	2	-	0/9/27/28	0/2/2/2
1	4AC	1	2937	1,67	-	1/11/29/30	0/2/2/2
2	4AC	2	379	2	-	0/11/29/30	0/2/2/2
1	4AC	1	1557	1	-	0/11/29/30	0/2/2/2
1	4AC	1	1621	1	-	0/11/29/30	0/2/2/2
2	OMG	2	873	2	-	1/9/27/28	0/3/3/3
1	4AC	1	1594	1,67	-	0/11/29/30	0/2/2/2
1	4AC	1	1873	1	-	0/11/29/30	0/2/2/2
2	4AC	2	626	2	-	0/11/29/30	0/2/2/2
2	4AC	2	703	2	-	0/11/29/30	0/2/2/2
2	5MC	2	1025	2	-	0/7/25/26	0/2/2/2
2	OMU	2	20	2	-	6/9/27/28	0/2/2/2
1	4AC	1	458	1	-	0/11/29/30	0/2/2/2
1	4AC	1	1695	1	-	0/11/29/30	0/2/2/2

The worst 5 of 298 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	2	1469	6MZ	C5-C4	4.58	1.47	1.39
2	2	1487	MA6	C5-C4	4.52	1.47	1.39
1	1	1055	A2M	C5-C4	4.52	1.47	1.39
1	1	2173	A2M	C5-C4	4.50	1.47	1.39
2	2	1488	MA6	C5-C4	4.41	1.47	1.39

The worst 5 of 647 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	2	657	OMG	C5-C4-N3	-6.59	117.77	128.46
2	2	471	OMG	C5-C4-N3	-6.38	118.12	128.46
2	2	680	OMG	C5-C4-N3	-6.37	118.13	128.46
1	1	789	OMG	C5-C4-N3	-6.33	118.18	128.46
1	1	2144	OMG	C5-C4-N3	-6.26	118.30	128.46

There are no chirality outliers.

5 of 49 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	1	789	OMG	C3'-C4'-C5'-O5'
1	1	2147	OMG	C3'-C4'-C5'-O5'
1	1	2585	4AC	O4'-C1'-N1-C2
1	1	2585	4AC	O4'-C1'-N1-C6
1	1	2608	4AC	O4'-C1'-N1-C2

There are no ring outliers.

107 monomers are involved in 177 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	1	694	4AC	1	0
2	2	319	4AC	1	0
2	2	680	OMG	1	0
1	1	2865	4AC	1	0
2	2	53	4AC	2	0
2	2	394	4AC	1	0
1	1	2001	4AC	4	0
2	2	1380	OMU	2	0
2	2	839	4AC	2	0
1	1	1550	4AC	2	0
2	2	636	4AC	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	1	2083	4AC	2	0
2	2	1177	OMU	2	0
1	1	789	OMG	2	0
1	1	1554	4AC	1	0
1	1	1222	4AC	2	0
1	1	1667	4AC	4	0
1	1	1765	4AC	2	0
1	1	2585	4AC	2	0
2	2	868	4AC	1	0
1	1	2992	4AC	1	0
2	2	479	4AC	2	0
1	1	2027	4AC	1	0
2	2	913	OMG	1	0
2	2	471	OMG	1	0
1	1	1617	4AC	4	0
1	1	1243	4AC	1	0
1	1	161	4AC	1	0
1	1	357	4AC	1	0
2	2	1184	4AC	2	0
2	2	64	OMU	1	0
1	1	829	4AC	1	0
1	1	1878	4AC	2	0
1	1	2548	4AC	2	0
1	1	922	4AC	2	0
2	2	1239	4AC	1	0
1	1	91	4AC	2	0
2	2	718	4AC	1	0
1	1	451	4AC	3	0
1	1	2608	4AC	3	0
1	1	314	4AC	1	0
2	2	731	4AC	2	0
1	1	1867	4AC	1	0
2	2	546	4AC	1	0
1	1	2329	4AC	1	0
1	1	3004	4AC	2	0
2	2	848	4AC	1	0
1	1	211	4AC	2	0
3	3	32	4AC	1	0
1	1	136	4AC	1	0
1	1	2925	4AC	2	0
2	2	1069	OMG	1	0
2	2	1479	4AC	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	1	2495	4AC	1	0
1	1	533	4AC	2	0
1	1	1055	A2M	1	0
1	1	2198	5MC	1	0
1	1	2733	5MC	2	0
1	1	2136	4AC	2	0
1	1	1065	4AC	3	0
2	2	1147	4AC	1	0
1	1	2545	4AC	3	0
1	1	2249	4AC	2	0
1	1	1293	4AC	1	0
2	2	1496	5MC	1	0
1	1	2908	4AC	2	0
1	1	1460	4AC	3	0
1	1	2287	4AC	5	0
1	1	2718	4AC	2	0
1	1	1938	4AC	1	0
1	1	341	4AC	1	0
1	1	2678	OMG	2	0
1	1	243	4AC	1	0
2	2	590	4AC	1	0
2	2	373	A2M	2	0
1	1	1048	4AC	2	0
2	2	828	4AC	1	0
2	2	1233	4AC	1	0
2	2	511	4AC	1	0
1	1	1755	4AC	3	0
1	1	1780	4AC	3	0
2	2	957	4AC	1	0
1	1	2162	5MC	2	0
1	1	2163	5MC	3	0
1	1	835	4AC	2	0
1	1	2960	4AC	2	0
1	1	2062	4AC	1	0
2	2	1488	MA6	1	0
1	1	1934	4AC	2	0
2	2	1376	OMC	2	0
1	1	1068	4AC	2	0
1	1	981	4AC	1	0
1	1	2642	4AC	1	0
1	1	1662	4AC	2	0
1	1	802	4AC	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	1	229	4AC	2	0
1	1	615	OMC	1	0
2	2	379	4AC	1	0
1	1	1557	4AC	1	0
1	1	1621	4AC	1	0
2	2	873	OMG	1	0
1	1	1594	4AC	1	0
2	2	626	4AC	2	0
2	2	703	4AC	4	0
2	2	1025	5MC	1	0
1	1	458	4AC	3	0
1	1	1695	4AC	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 257 ligands modelled in this entry, 257 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	1	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	1	107:G	O3'	112:A	P	11.57

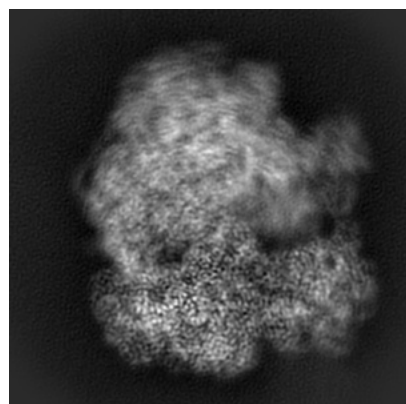
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-55135. These allow visual inspection of the internal detail of the map and identification of artifacts.

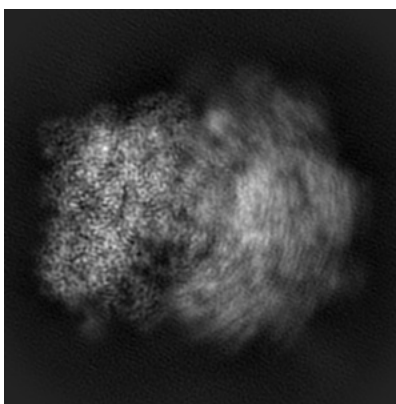
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

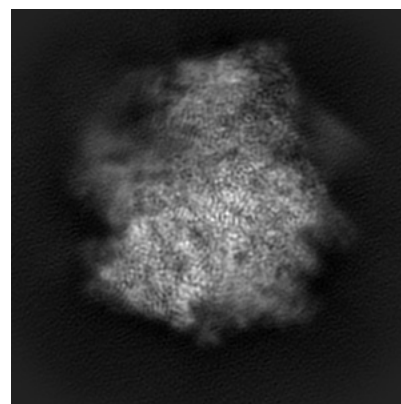
6.1.1 Primary map



X

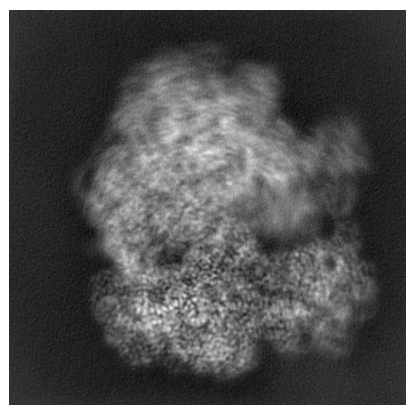


Y

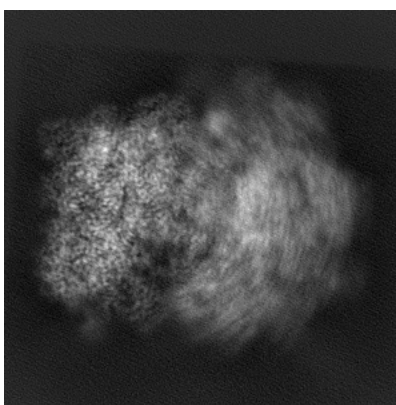


Z

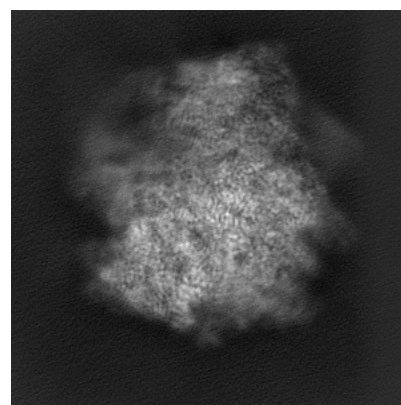
6.1.2 Raw map



X



Y

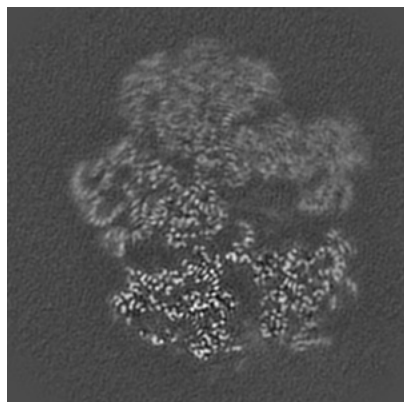


Z

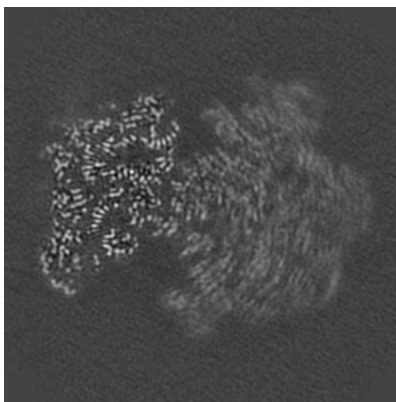
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

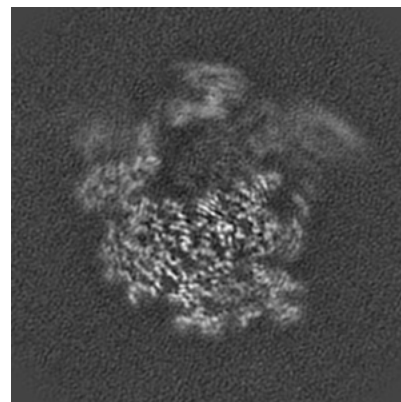
6.2.1 Primary map



X Index: 180

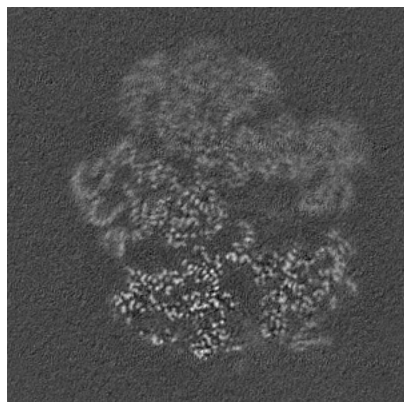


Y Index: 180

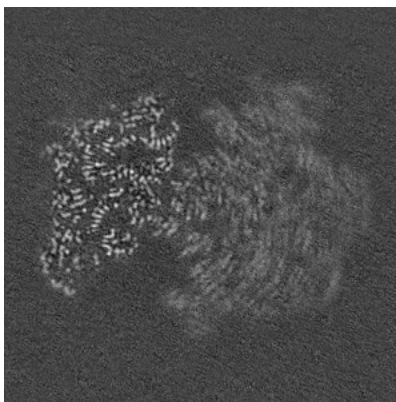


Z Index: 180

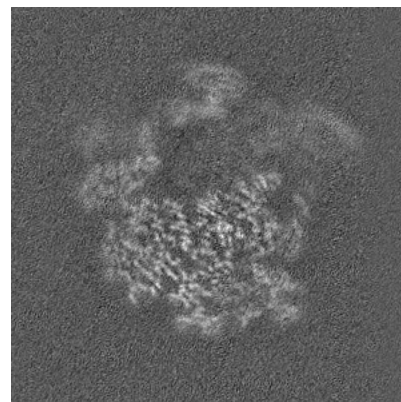
6.2.2 Raw map



X Index: 180



Y Index: 180

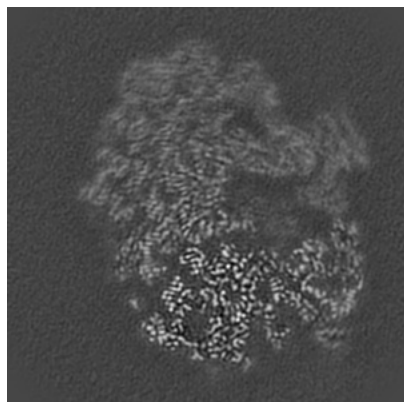


Z Index: 180

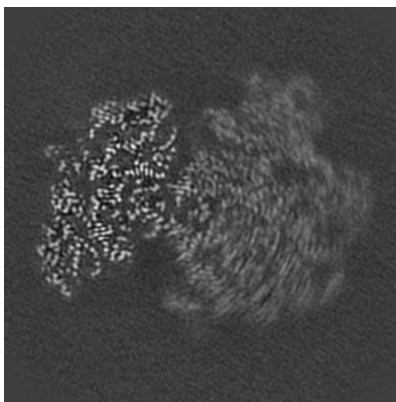
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

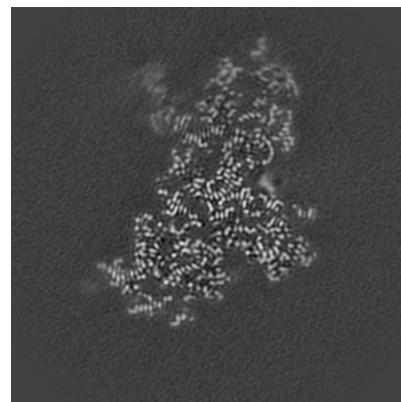
6.3.1 Primary map



X Index: 196

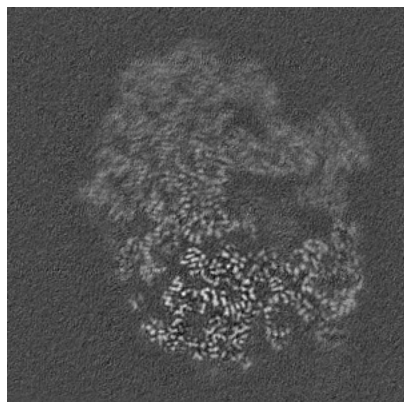


Y Index: 175

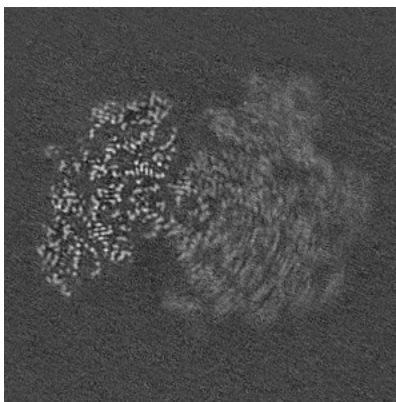


Z Index: 91

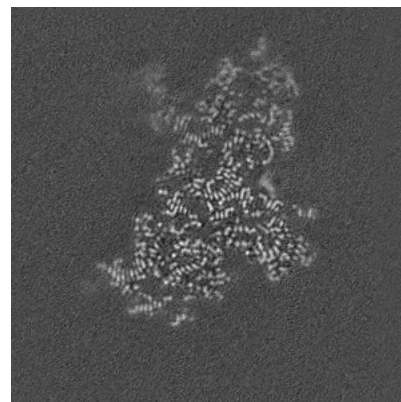
6.3.2 Raw map



X Index: 195



Y Index: 175

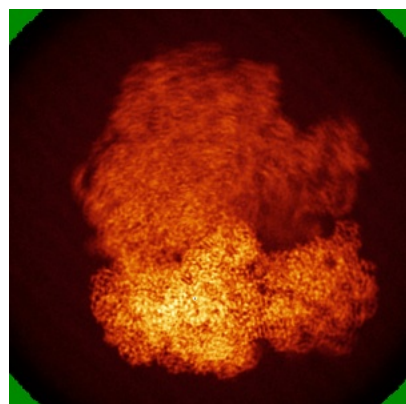


Z Index: 91

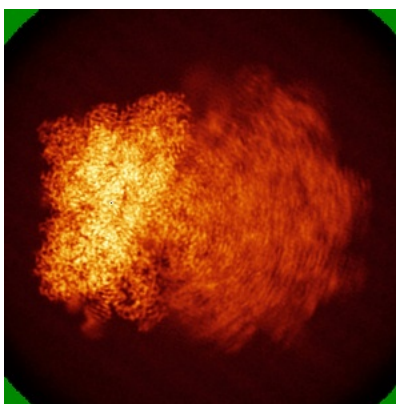
The images above show the largest variance slices of the map in three orthogonal directions.

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

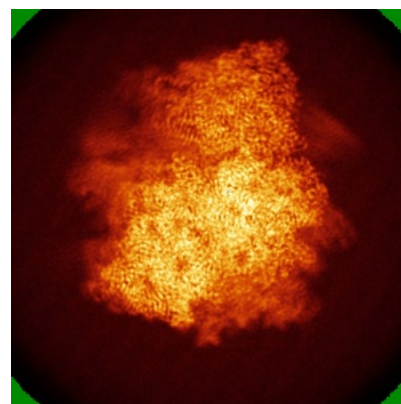
6.4.1 Primary map



X

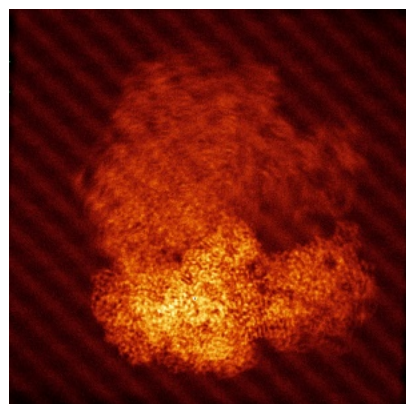


Y

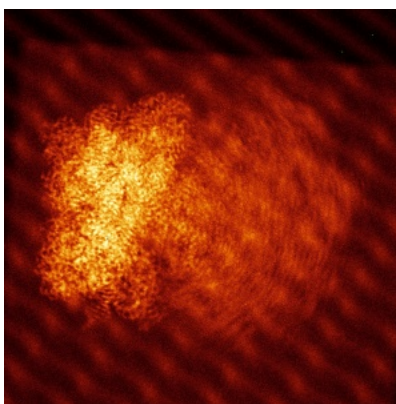


Z

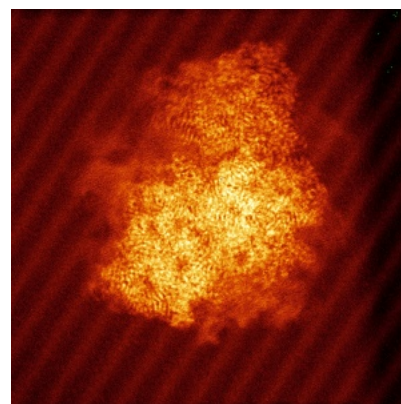
6.4.2 Raw map



X



Y



Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



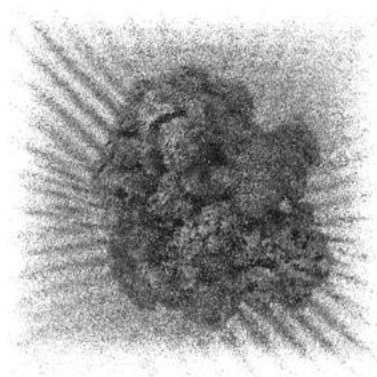
Y



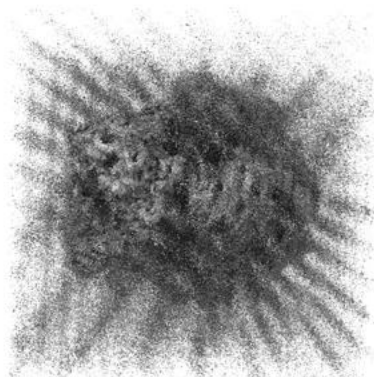
Z

The images above show the 3D surface view of the map at the recommended contour level 0.07. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

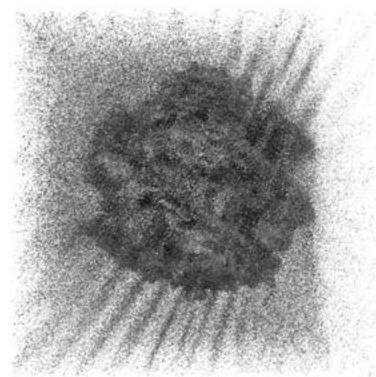
6.5.2 Raw map



X



Y



Z

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

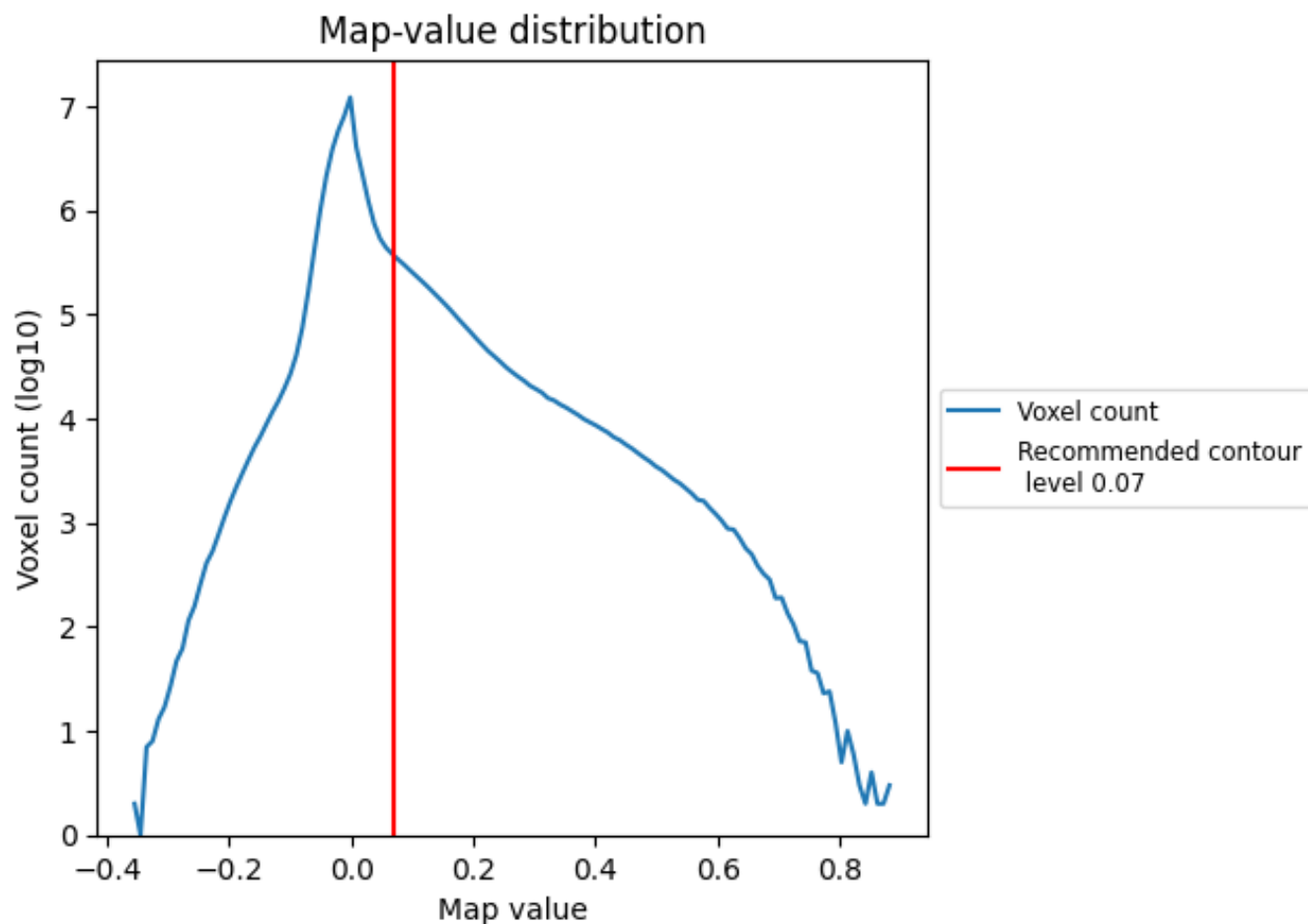
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

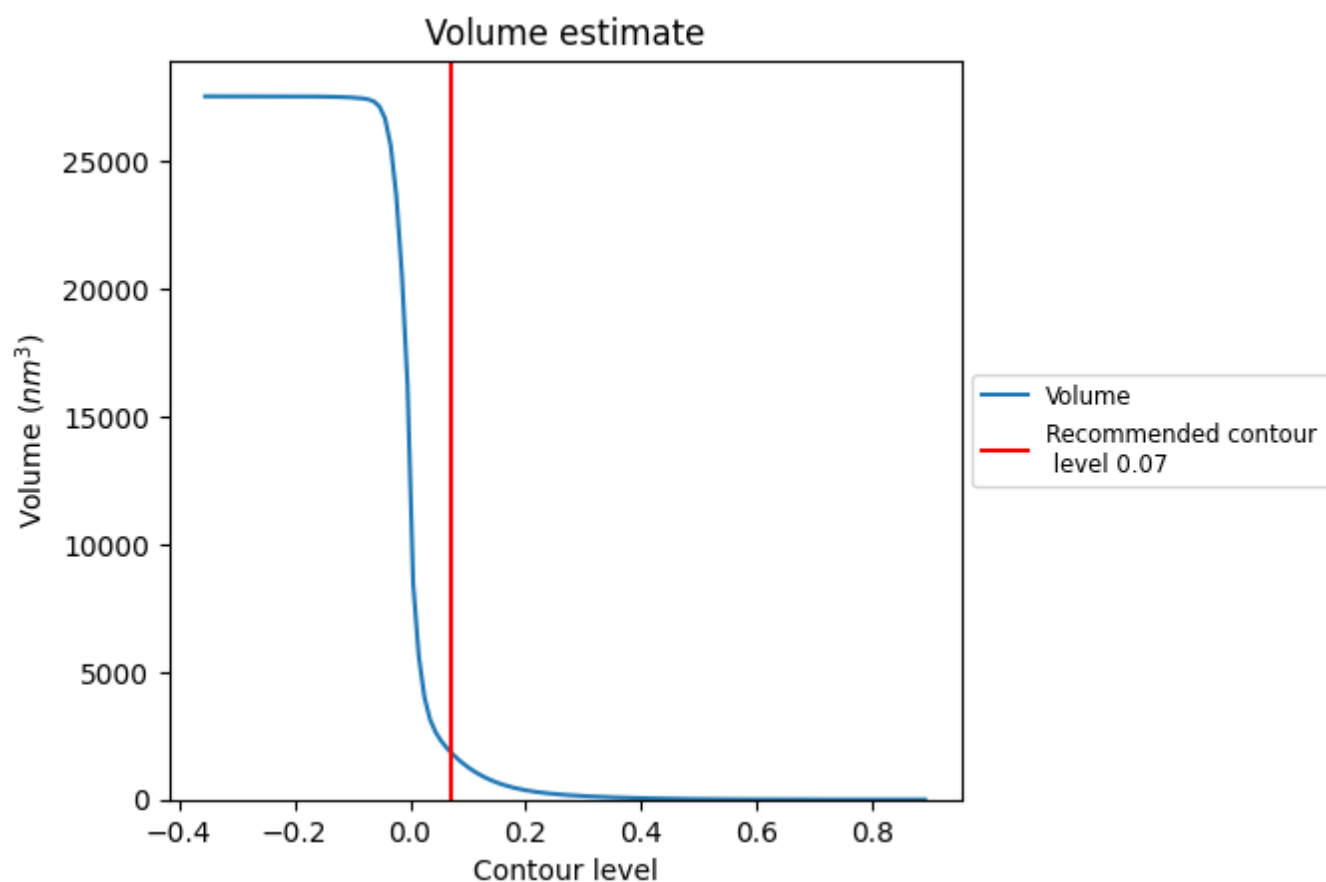
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

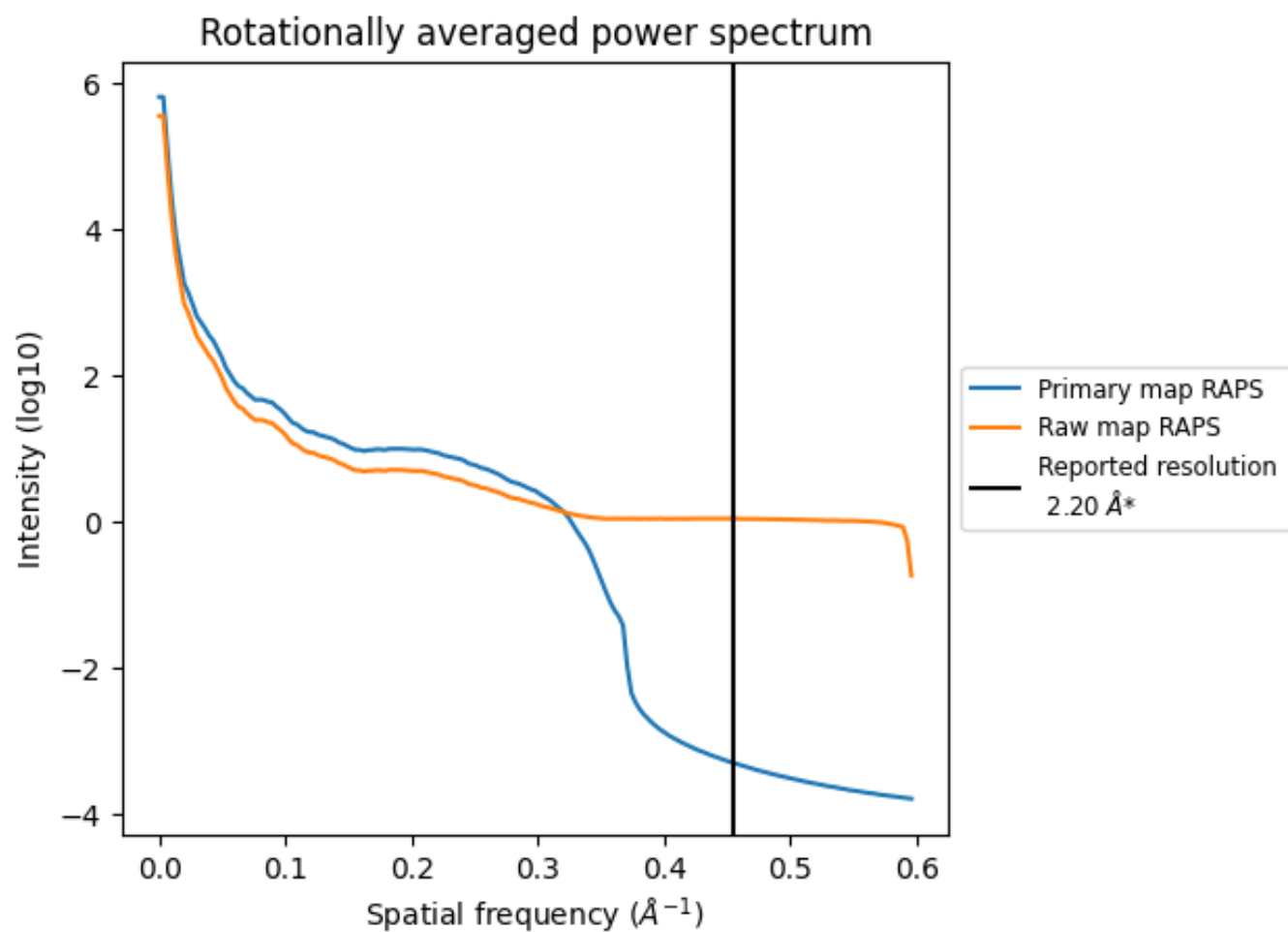
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1867 nm^3 ; this corresponds to an approximate mass of 1686 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 ⓘ

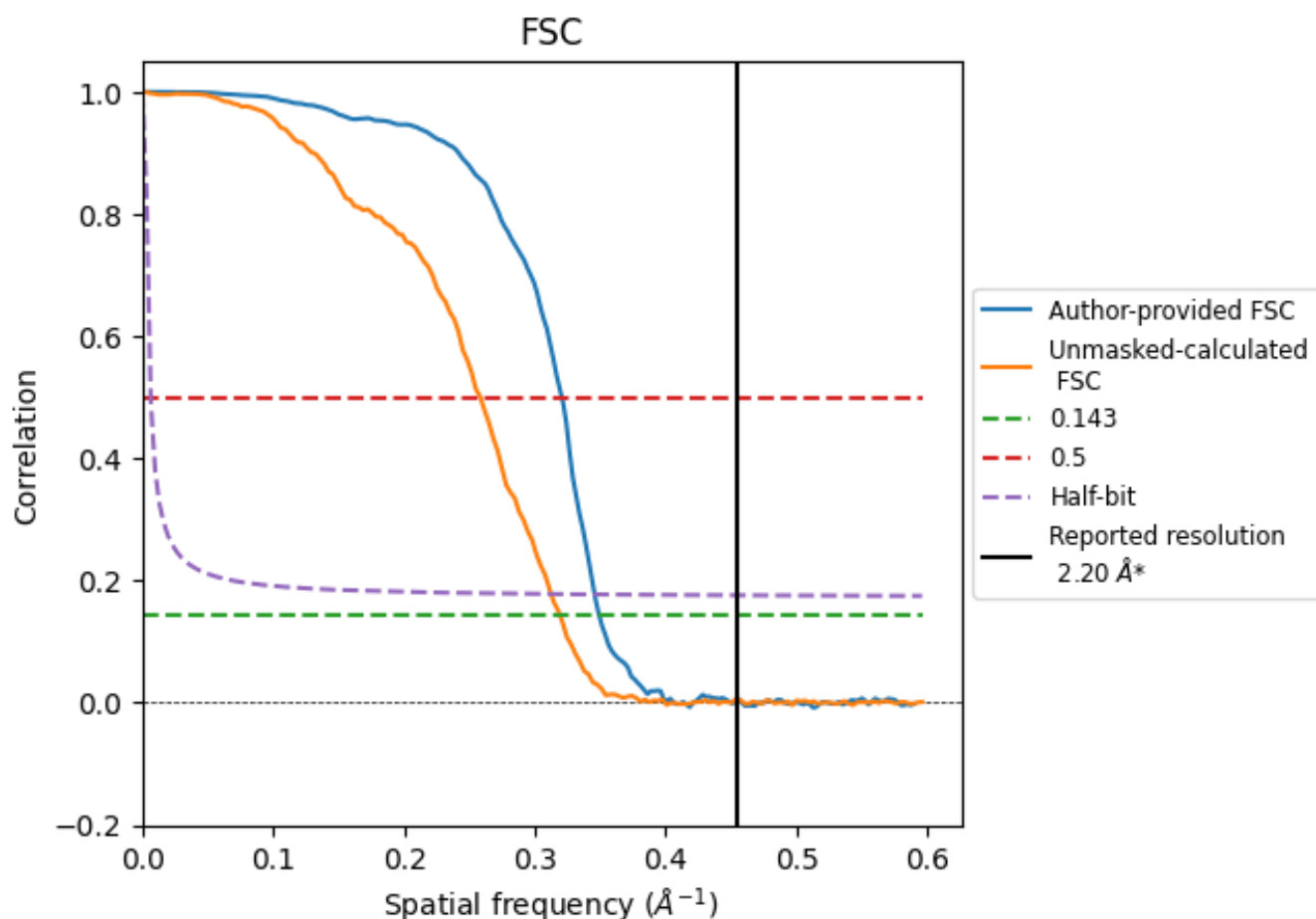


*Reported resolution corresponds to spatial frequency of $0.455 \text{ \AA}^{-1}$

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.455 Å⁻¹

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.20	-	-
Author-provided FSC curve	2.86	3.12	2.89
Unmasked-calculated*	3.13	3.87	3.20

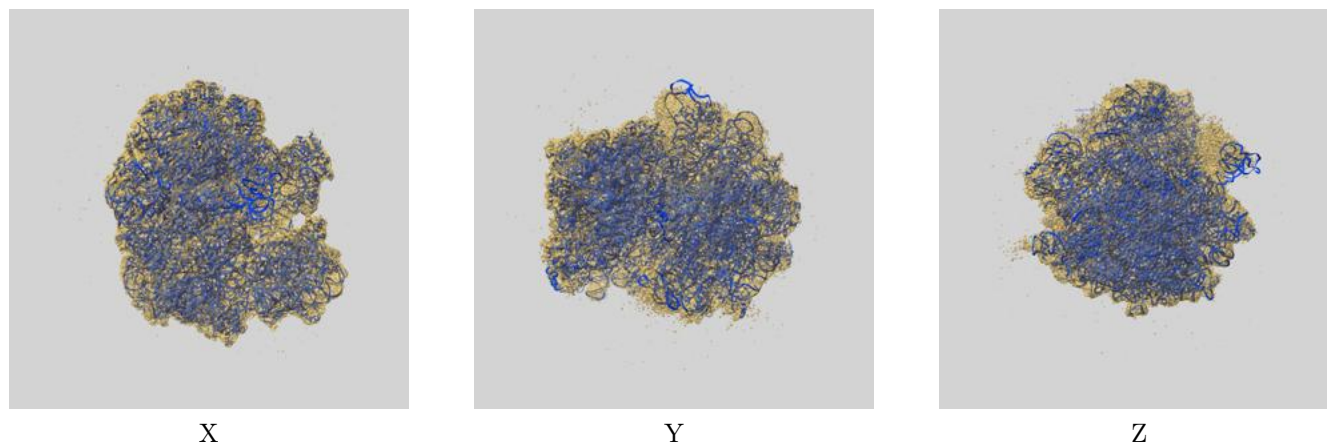
*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from author-provided FSC intersecting FSC 0.143 CUT-OFF 2.86 differs from the reported value 2.2 by more than 10 %

The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.13 differs from the reported value 2.2 by more than 10 %

9 Map-model fit [i](#)

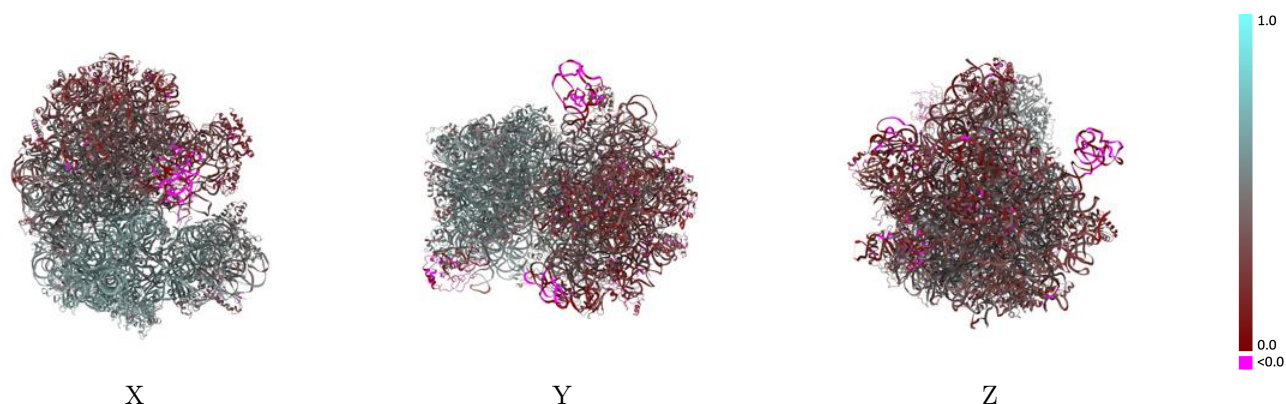
This section contains information regarding the fit between EMDB map EMD-55135 and PDB model 9SRA. Per-residue inclusion information can be found in [section 3](#) on [page 17](#).

9.1 Map-model overlay [i](#)



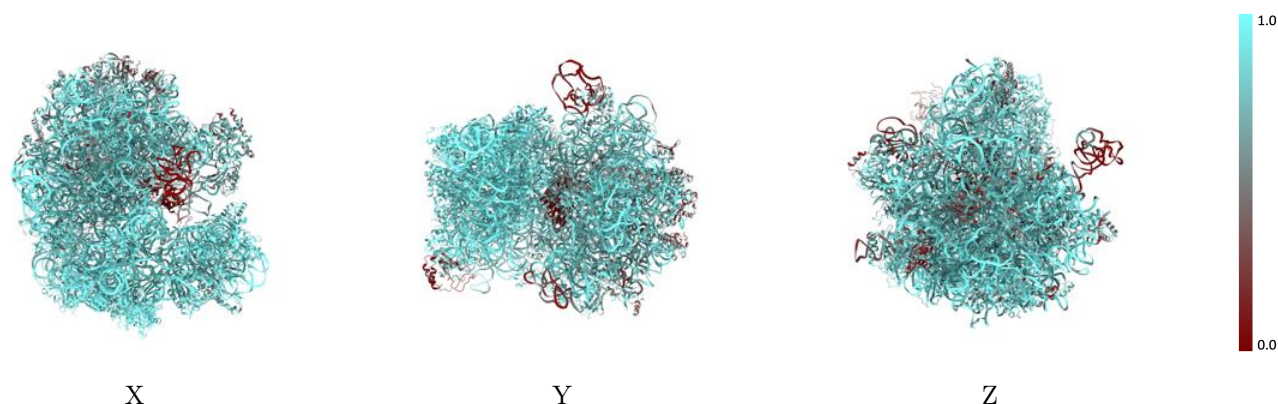
The images above show the 3D surface view of the map at the recommended contour level 0.07 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



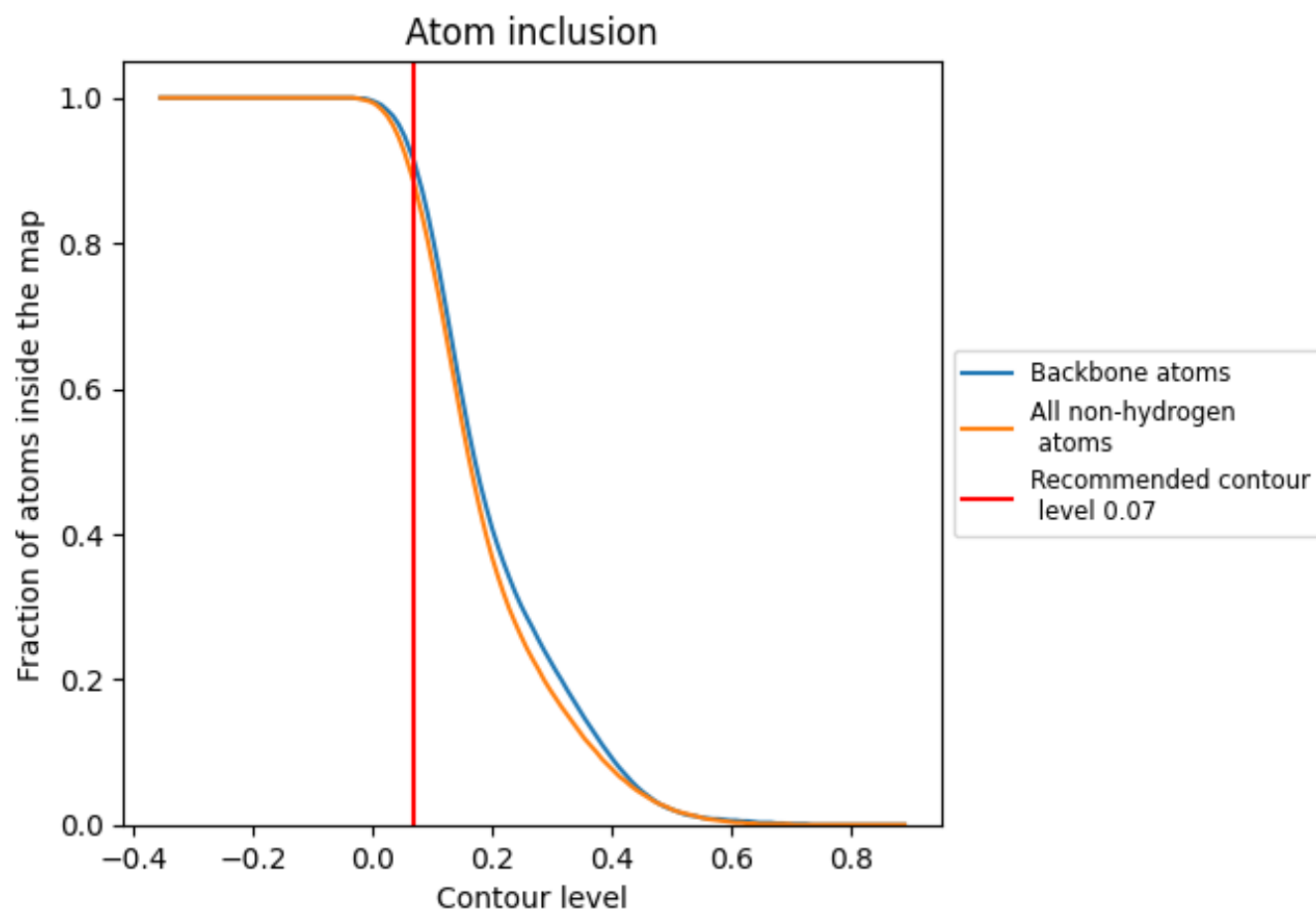
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.07).

9.4 Atom inclusion [i](#)



At the recommended contour level, 91% of all backbone atoms, 88% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.07) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8820	 0.4130
1	 0.9040	 0.3630
2	 0.9930	 0.5590
3	 0.9120	 0.2480
A0	 0.9500	 0.5960
A3	 0.3670	 0.0910
AA	 0.9400	 0.5430
AB	 0.9630	 0.5590
AC	 0.9660	 0.5740
AD	 0.9680	 0.5590
AE	 0.9680	 0.5650
AF	 0.9620	 0.5750
AG	 0.9380	 0.5040
AH	 0.9670	 0.5550
AI	 0.9790	 0.5980
AJ	 0.9540	 0.5500
AK	 0.9420	 0.5120
AL	 0.8550	 0.4280
AM	 0.9440	 0.5610
AN	 0.9540	 0.5760
AO	 0.9010	 0.4330
AP	 0.9290	 0.5180
AQ	 0.9590	 0.5670
AR	 0.9700	 0.5860
AS	 0.9330	 0.4910
AT	 0.8780	 0.4330
AU	 0.9380	 0.4730
AV	 0.9390	 0.5140
AW	 0.9740	 0.5280
AX	 0.8970	 0.4850
AY	 0.2330	 0.0680
AZ	 0.8930	 0.4620
B4	 0.3280	 0.1520
B5	 0.8530	 0.4050
B6	 0.3360	 0.1680



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Chain	Atom inclusion	Q-score
BB	 0.8710	 0.4790
BC	 0.8510	 0.3840
BD	 0.7510	 0.2670
BE	 0.7160	 0.1880
BF	 0.8040	 0.2580
BG	 0.6820	 0.2780
BI	 0.7730	 0.3000
BJ	 0.8560	 0.4420
BK	 0.7020	 0.1860
BL	 0.6410	 0.2590
BM	 0.8230	 0.3600
BN	 0.7610	 0.2940
BO	 0.6370	 0.2020
BP	 0.7400	 0.2220
BQ	 0.8810	 0.4570
BR	 0.7500	 0.2660
BS	 0.8760	 0.4130
BT	 0.7910	 0.3520
BU	 0.7700	 0.2700
BV	 0.9100	 0.4140
BW	 0.6870	 0.2640
BY	 0.7130	 0.2530
BZ	 0.8630	 0.4450
Ba	 0.8820	 0.4430
Bb	 0.7500	 0.2990
Bc	 0.7190	 0.2510
Bd	 0.9000	 0.4510
Be	 0.8310	 0.4000
Bf	 0.8380	 0.3500
Bg	 0.8730	 0.2910
Bi	 0.8790	 0.5000
Bj	 0.6700	 0.3110
Bk	 0.7950	 0.2740
Bl	 0.5770	 0.1920
H	 0.3280	 0.3540
sd	 0.9400	 0.4010