



## wwPDB EM Validation Summary Report ⓘ

Dec 11, 2022 – 04:42 pm GMT

PDB ID : 6SKG  
EMDB ID : EMD-10224  
Title : Cryo-EM Structure of *T. kodakarensis* 70S ribosome in TkNat10 deleted strain  
Authors : Matzov, D.; Sas-Chen, A.; Thomas, J.M.; Santangelo, T.; Meier, J.L.; Schwartz, S.; Shalev-Benami, M.  
Deposited on : 2019-08-15  
Resolution : 2.65 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](mailto: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>.

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

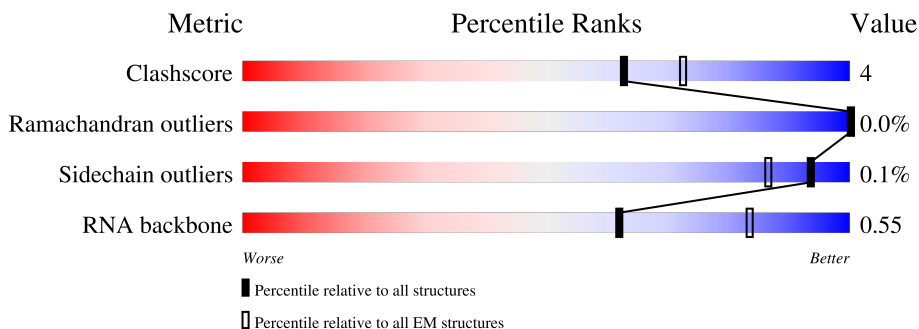
EMDB validation analysis : 0.0.1.dev43  
Mogul : 1.8.4, CSD as541be (2020)  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.3

# 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 2.65 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826
RNA backbone	4643	859

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  $\geq 3$ , 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	Aa	1498	8% (red), 78% (green), 18% (yellow), 2% (grey), 2% (grey)
2	Ab	201	20% (red), 98% (green), 2% (grey)
3	Ac	209	71% (red), 93% (green), 7% (grey)
4	Ad	200	25% (red), 95% (green), 5% (grey)
5	Ae	180	13% (red), 97% (green), 2% (grey)
6	Af	243	12% (red), 100% (green)
7	Ag	235	10% (red), 95% (green), 5% (grey)



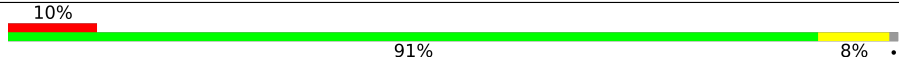
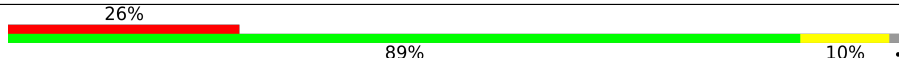
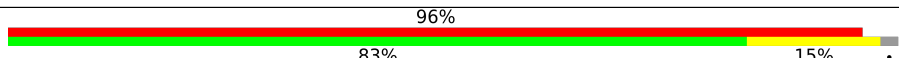
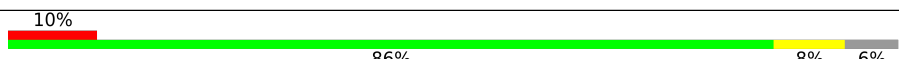
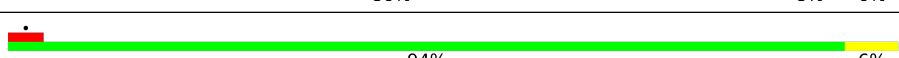
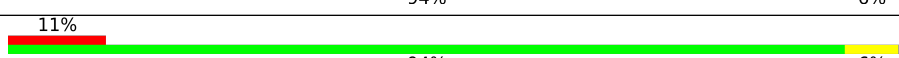
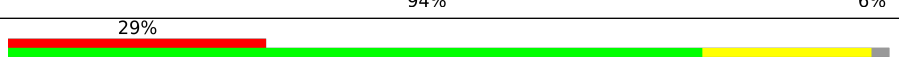

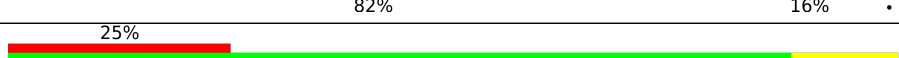

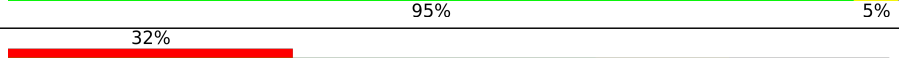



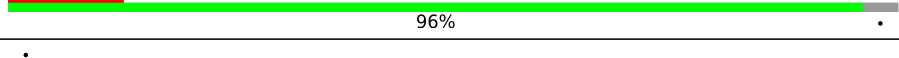
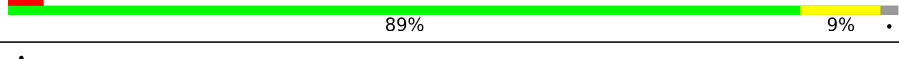
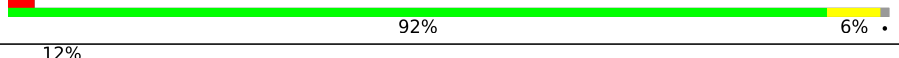
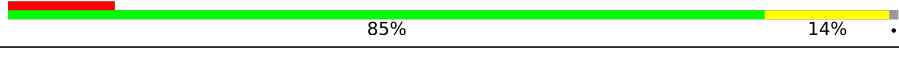
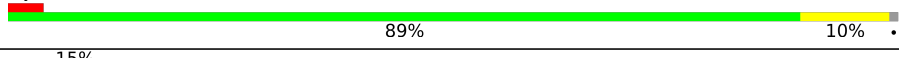
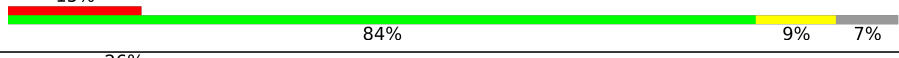
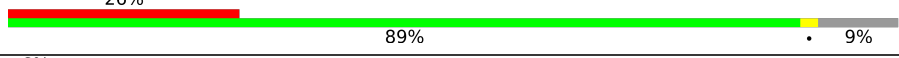
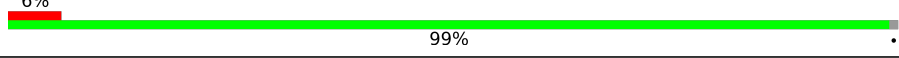
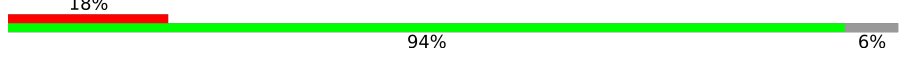
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Mol	Chain	Length	Quality of chain
8	Ah	125	49% 98%
9	Ai	215	35% 100%
10	Aj	130	5% 99%
11	Ak	130	11% 96%
12	Al	135	30% 99%
13	Am	102	70% 98%
14	An	140	17% 88% 9%
15	Ao	147	11% 97%
16	Ap	149	32% 91% 9%
17	Aq	151	13% 99%
18	Ar	56	38% 96%
19	As	114	11% 95% 5%
20	At	67	64% 96%
20	Bm	67	33% 94% 6%
21	Au	133	30% 87% 13%
22	Av	150	27% 99%
23	Aw	98	20% 97%
23	Bn	98	33% 94% 6%
24	Ax	65	28% 94% 6%
25	Ay	70	53% 91% 9%
26	Az	62	11% 89% 11%
27	BA	3037	9% 66% 24% 6%
28	BB	126	17% 72% 23%
29	BC	239	5% 92% 8%
30	BD	346	6% 90% 10%

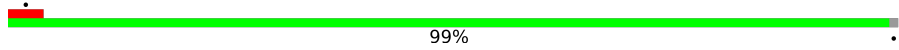
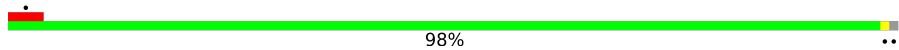

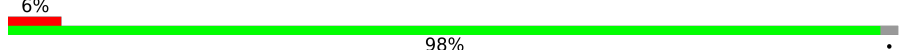

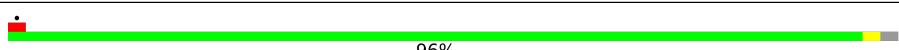
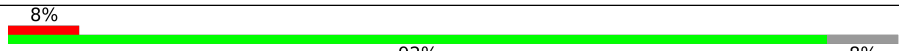
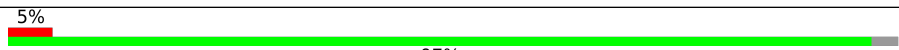
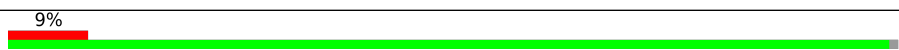
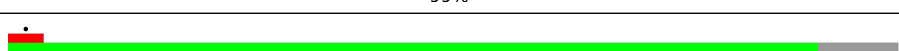
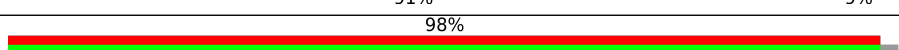
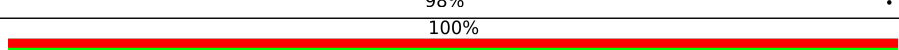
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Mol	Chain	Length	Quality of chain
31	BE	255	 87% 13%
32	BF	183	 83% 67% 23% 9%
33	BG	184	 10% 91% 8%
34	BH	123	 26% 89% 10%
34	BI	123	 96% 83% 15%
35	BJ	182	 10% 86% 8% 6%
36	BK	142	 94% 6%
37	BL	141	 11% 94% 6%
38	BM	83	 29% 78% 19%
38	BN	83	 31% 82% 16%
39	BO	148	 25% 88% 12%
40	BP	194	 95% 5%
41	BQ	201	 32% 66% 15% 18%
42	BR	121	 85% 14%
43	BS	150	 5% 88% 9%
44	BT	77	 13% 96%
45	BU	98	 89% 9%
46	BV	156	 92% 6%
47	BW	86	 12% 85% 14%
48	BX	121	 89% 10%
49	BY	67	 15% 84% 9% 7%
50	BZ	66	 26% 89% 9%
51	Ba	155	 6% 99%
52	Bb	102	 18% 94% 6%
53	Bc	90	 10% 99%

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Mol	Chain	Length	Quality of chain
54	Bd	125	 99%
55	Be	90	 98%
56	Bf	101	 85% 15%
57	Bg	86	 98%
58	Bh	63	 98%
59	Bi	51	 96%
60	Bj	51	 92% 8%
61	Bk	37	 97%
62	Bl	94	 99%
63	Bo	47	 91% 9%
64	Bp	236	 98%
65	Bq	34	 100%

## 2 Entry composition [i](#)

There are 67 unique types of molecules in this entry. The entry contains 167402 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 ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	P		
1	Aa	1463	31531	14063	5836	10169	1463	0	0

- Molecule 2 is a protein called 30S ribosomal protein S2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	Ab	196	1580	1021	272	284	3	0	0

- Molecule 3 is a protein called 30S ribosomal protein S3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	Ac	195	1536	979	283	271	3	0	0

- Molecule 4 is a protein called 30S ribosomal protein S3Ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	Ad	190	1550	997	277	273	3	0	0

- Molecule 5 is a protein called 30S ribosomal protein S4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
5	Ae	174	1454	912	287	253	2	0	0

- Molecule 6 is a protein called 30S ribosomal protein S4e.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	Af	242	1952	1261	349	337	5	0	0

- Molecule 7 is a protein called 30S ribosomal protein S5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
7	Ag	223	1753	1107	327	312	7	0	0

- Molecule 8 is a protein called 30S ribosomal protein S6e.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	Ah	123	952	599	179	172	2	0	0

- Molecule 9 is a protein called 30S ribosomal protein S7.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	Ai	214	1715	1087	317	303	8	0	0

- Molecule 10 is a protein called 30S ribosomal protein S8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	Aj	129	1020	662	176	180	2	0	0

- Molecule 11 is a protein called 30S ribosomal protein S8e.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
11	Ak	125	982	615	198	169	0	0

- Molecule 12 is a protein called 30S ribosomal protein S9.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
12	Al	133	1054	656	205	188	5	0	0

- Molecule 13 is a protein called 30S ribosomal protein S10.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
13	Am	100	803	498	154	148	3	0	0

- Molecule 14 is a protein called 30S ribosomal protein S11.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	An	127	Total	C	N	O	S	0	0
			953	587	190	173	3		

- Molecule 15 is a protein called 30S ribosomal protein S12.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	Ao	143	Total	C	N	O	S	0	0
			1122	712	216	192	2		

- Molecule 16 is a protein called 30S ribosomal protein S13.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	Ap	135	Total	C	N	O	S	0	0
			1090	691	214	181	4		

- Molecule 17 is a protein called 30S ribosomal protein S15.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	Aq	149	Total	C	N	O	S	0	0
			1217	776	233	206	2		

- Molecule 18 is a protein called 30S ribosomal protein S14 type Z.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	Ar	54	Total	C	N	O	S	0	0
			447	284	92	65	6		

- Molecule 19 is a protein called 30S ribosomal protein S17.

Mol	Chain	Residues	Atoms					AltConf	Trace
19	As	108	Total	C	N	O	S	0	0
			879	560	164	152	3		

- Molecule 20 is a protein called 30S ribosomal protein S17e.

Mol	Chain	Residues	Atoms					AltConf	Trace
20	At	64	Total	C	N	O	S	0	0
			538	338	103	95	2		
20	Bm	63	Total	C	N	O	S	0	0
			534	336	102	94	2		

- Molecule 21 is a protein called 30S ribosomal protein S19.



Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
21	Au	116	952	608	175	163	6	0	0

- Molecule 22 is a protein called 30S ribosomal protein S19e.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
22	Av	149	1216	788	215	213		0	0

- Molecule 23 is a protein called 30S ribosomal protein S24e.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
23	Aw	95	777	497	131	146	3	0	0
23	Bn	92	752	481	126	142	3	0	0

- Molecule 24 is a protein called 30S ribosomal protein S27e.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
24	Ax	61	461	294	83	79	5	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Ax	64	GLU	LEU	conflict	UNP Q5JE50

- Molecule 25 is a protein called 30S ribosomal protein S28e.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
25	Ay	64	499	307	99	93	0	0

- Molecule 26 is a protein called Predicted zinc-ribbon RNA-binding protein involved in translation.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
26	Az	55	426	269	74	75	8	0	0

- Molecule 27 is a RNA chain called 23S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	N	O	P			S
27	BA	2934	63209	28168	11754	20351	2934	2	0	0

- Molecule 28 is a RNA chain called 5S ribosomal RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace	
			Total	C	N	O	P			S
28	BB	125	2678	1191	492	870	125		0	0

- Molecule 29 is a protein called 50S ribosomal protein L2.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
29	BC	237	1820	1150	356	311	3	0	0

- Molecule 30 is a protein called 50S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
30	BD	344	2746	1765	505	469	7	0	0

- Molecule 31 is a protein called 50S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
31	BE	255	2026	1286	389	347	4	0	0

- Molecule 32 is a protein called 50S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
32	BF	166	1339	839	258	235	7	0	0

- Molecule 33 is a protein called 50S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
33	BG	183	1463	942	253	266	2	0	0

- Molecule 34 is a protein called 50S ribosomal protein L7Ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
34	BH	121	Total	C	N	O	S	0	0
			928	591	154	180	3		
34	BI	121	Total	C	N	O	S	0	0
			928	591	154	180	3		

- Molecule 35 is a protein called 50S ribosomal protein L10e.

Mol	Chain	Residues	Atoms					AltConf	Trace
35	BJ	171	Total	C	N	O	S	0	0
			1399	891	261	240	7		

- Molecule 36 is a protein called 50S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	BK	142	Total	C	N	O	S	0	0
			1146	731	213	198	4		

- Molecule 37 is a protein called 50S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	BL	140	Total	C	N	O	S	0	0
			1055	658	214	180	3		

- Molecule 38 is a protein called 50S ribosomal protein L14e.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	BM	81	Total	C	N	O	S	0	0
			610	382	119	108	1		
38	BN	81	Total	C	N	O	S	0	0
			610	382	119	108	1		

- Molecule 39 is a protein called 50S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	BO	148	Total	C	N	O	S	0	0
			1166	743	222	198	3		

- Molecule 40 is a protein called 50S ribosomal protein L15e.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	BP	193	Total	C	N	O	S	0	0
			1582	1010	316	250	6		

- Molecule 41 is a protein called 50S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
41	BQ	164	1314	841	243	228	2	0	0

- Molecule 42 is a protein called 50S ribosomal protein L18e.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
42	BR	120	959	601	187	169	2	0	0

- Molecule 43 is a protein called 50S ribosomal protein L19e.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
43	BS	146	1200	753	246	194	7	0	0

- Molecule 44 is a protein called 50S ribosomal protein L18Ae.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
44	BT	74	624	399	113	112	0	0

- Molecule 45 is a protein called 50S ribosomal protein L21e.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
45	BU	96	784	502	158	123	1	0	0

- Molecule 46 is a protein called 50S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
46	BV	154	1234	777	242	211	4	0	0

- Molecule 47 is a protein called 50S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
47	BW	85	683	438	119	123	3	0	0

- Molecule 48 is a protein called 50S ribosomal protein L24.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	BX	120	Total	C	N	O	S	0	0
			991	628	188	170	5		

- Molecule 49 is a protein called 50S ribosomal protein L24e.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	BY	62	Total	C	N	O	S	0	0
			524	334	99	85	6		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
50	BZ	60	Total	C	N	O	S	0	0
			506	314	98	90	4		

- Molecule 51 is a protein called 50S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	Ba	154	Total	C	N	O	S	0	0
			1242	788	235	214	5		

- Molecule 52 is a protein called 50S ribosomal protein L30e.

Mol	Chain	Residues	Atoms					AltConf	Trace
52	Bb	96	Total	C	N	O	S	0	0
			730	473	123	133	1		

- Molecule 53 is a protein called 50S ribosomal protein L31e.

Mol	Chain	Residues	Atoms				AltConf	Trace
53	Bc	89	Total	C	N	O	0	0
			721	463	140	118		

- Molecule 54 is a protein called 50S ribosomal protein L32e.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	Bd	124	Total	C	N	O	S	0	0
			1022	650	208	162	2		

- Molecule 55 is a protein called 50S ribosomal protein L34e.

Mol	Chain	Residues	Atoms					AltConf	Trace
55	Be	89	Total	C	N	O	S	0	0
			728	454	155	110	9		

- Molecule 56 is a protein called Ribosomal protein eL35A.

Mol	Chain	Residues	Atoms					AltConf	Trace
56	Bf	86	Total	C	N	O	S	0	0
			669	424	133	112			

- Molecule 57 is a protein called 50S ribosomal protein L37Ae.

Mol	Chain	Residues	Atoms					AltConf	Trace
57	Bg	84	Total	C	N	O	S	0	0
			630	392	132	101	5		

- Molecule 58 is a protein called 50S ribosomal protein L37e.

Mol	Chain	Residues	Atoms					AltConf	Trace
58	Bh	62	Total	C	N	O	S	0	0
			512	314	118	75	5		

- Molecule 59 is a protein called 50S ribosomal protein L39e.

Mol	Chain	Residues	Atoms				AltConf	Trace
59	Bi	50	Total	C	N	O	0	0
			432	276	97	59		

- Molecule 60 is a protein called 50S ribosomal protein L40e.

Mol	Chain	Residues	Atoms					AltConf	Trace
60	Bj	47	Total	C	N	O	S	0	0
			373	233	79	56	5		

- Molecule 61 is a protein called LSU ribosomal protein L41E.

Mol	Chain	Residues	Atoms					AltConf	Trace
61	Bk	36	Total	C	N	O	S	0	0
			345	218	86	39	2		

- Molecule 62 is a protein called 50S ribosomal protein L44e.

Mol	Chain	Residues	Atoms					AltConf	Trace
62	Bl	93	Total	C	N	O	S	0	0
			778	494	160	119	5		

- Molecule 63 is a protein called Nucleic acid-binding protein, containing C2H2 zinc-finger.

Mol	Chain	Residues	Atoms					AltConf	Trace
63	Bo	43	Total	C	N	O	S	0	0
			349	222	69	55	3		

- Molecule 64 is a protein called Predicted exosome subunit, UPF0023 family.

Mol	Chain	Residues	Atoms					AltConf	Trace
64	Bp	232	Total	C	N	O	S	0	0
			1862	1193	323	342	4		

- Molecule 65 is a protein called Unknown ribosomal protein.

Mol	Chain	Residues	Atoms				AltConf	Trace
65	Bq	34	Total	C	N	O	0	0
			170	102	34	34		

- Molecule 66 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
66	Ag	1	Total	Zn	0
			1	1	
66	Ar	1	Total	Zn	0
			1	1	
66	As	1	Total	Zn	0
			1	1	
66	Ax	1	Total	Zn	0
			1	1	
66	Az	2	Total	Zn	0
			2	2	
66	BY	1	Total	Zn	0
			1	1	
66	Be	2	Total	Zn	0
			2	2	
66	Bg	1	Total	Zn	0
			1	1	
66	Bh	1	Total	Zn	0
			1	1	

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Mol	Chain	Residues	Atoms		AltConf
66	Bj	1	Total 1	Zn 1	0
66	Bl	1	Total 1	Zn 1	0
66	Bo	1	Total 1	Zn 1	0

- Molecule 67 is water.

Mol	Chain	Residues	Atoms		AltConf
67	Aa	266	Total 266	O 266	0
67	Ad	1	Total 1	O 1	0
67	Af	4	Total 4	O 4	0
67	Ag	1	Total 1	O 1	0
67	Ai	1	Total 1	O 1	0
67	Aj	2	Total 2	O 2	0
67	Ak	6	Total 6	O 6	0
67	Al	2	Total 2	O 2	0
67	Am	2	Total 2	O 2	0
67	An	2	Total 2	O 2	0
67	Ao	1	Total 1	O 1	0
67	Ap	3	Total 3	O 3	0
67	Aq	5	Total 5	O 5	0
67	As	2	Total 2	O 2	0
67	Au	1	Total 1	O 1	0
67	Av	1	Total 1	O 1	0

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Mol	Chain	Residues	Atoms		AltConf
67	Az	1	Total 1	O 1	0
67	BA	2715	Total 2715	O 2715	0
67	BB	23	Total 23	O 23	0
67	BC	51	Total 51	O 51	0
67	BD	59	Total 59	O 59	0
67	BE	41	Total 41	O 41	0
67	BF	2	Total 2	O 2	0
67	BG	1	Total 1	O 1	0
67	BH	1	Total 1	O 1	0
67	BJ	13	Total 13	O 13	0
67	BK	18	Total 18	O 18	0
67	BL	15	Total 15	O 15	0
67	BM	3	Total 3	O 3	0
67	BN	3	Total 3	O 3	0
67	BO	30	Total 30	O 30	0
67	BP	36	Total 36	O 36	0
67	BQ	6	Total 6	O 6	0
67	BR	14	Total 14	O 14	0
67	BS	17	Total 17	O 17	0
67	BT	2	Total 2	O 2	0
67	BU	13	Total 13	O 13	0

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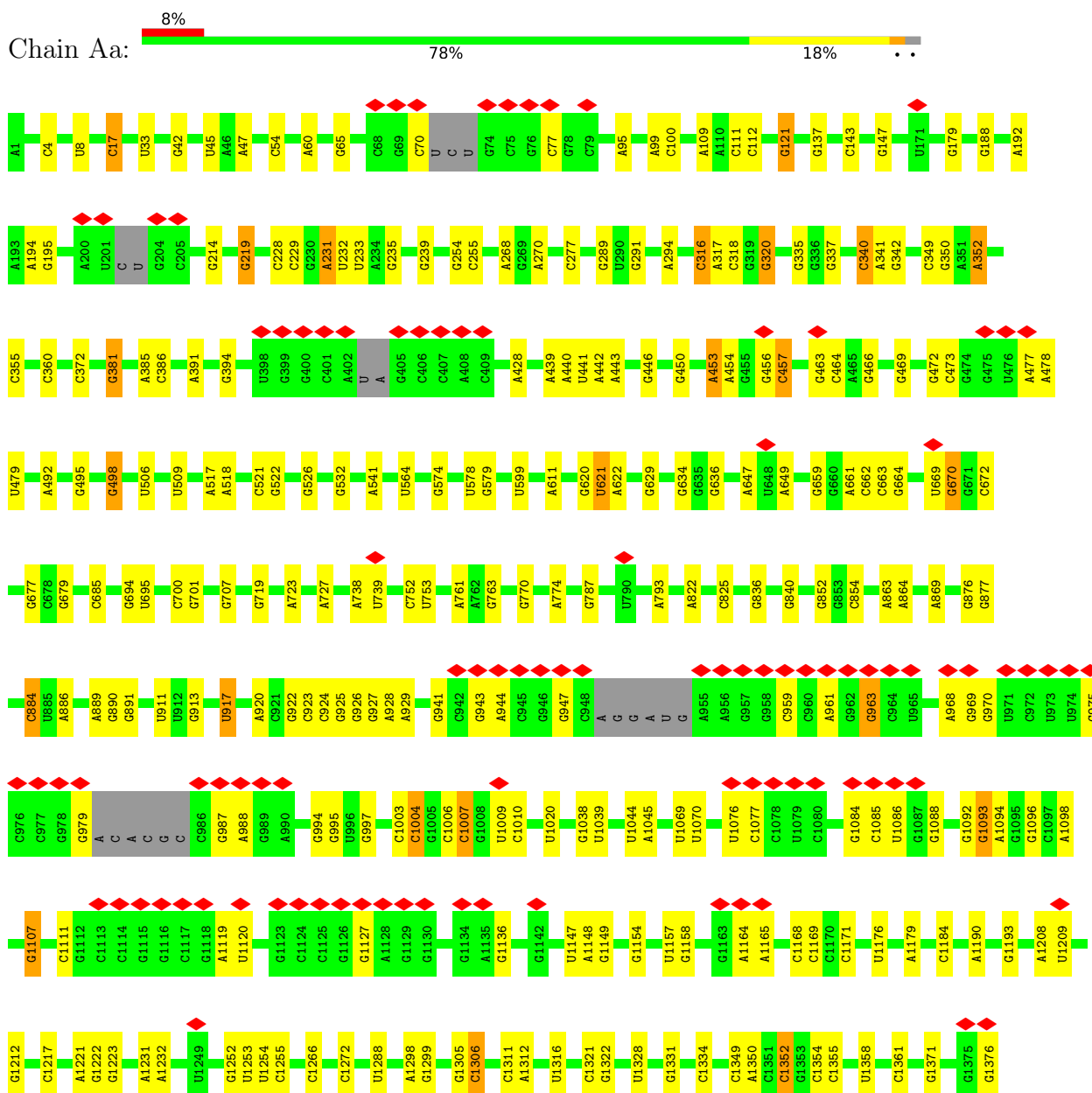
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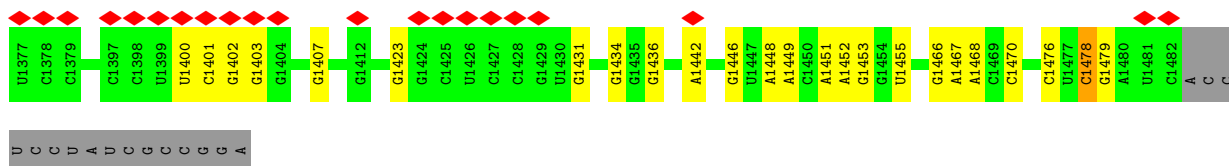
Mol	Chain	Residues	Atoms		AltConf
67	BV	22	Total 22	O 22	0
67	BW	3	Total 3	O 3	0
67	BX	10	Total 10	O 10	0
67	BY	12	Total 12	O 12	0
67	BZ	7	Total 7	O 7	0
67	Ba	14	Total 14	O 14	0
67	Bb	8	Total 8	O 8	0
67	Bc	10	Total 10	O 10	0
67	Bd	23	Total 23	O 23	0
67	Be	8	Total 8	O 8	0
67	Bf	7	Total 7	O 7	0
67	Bg	15	Total 15	O 15	0
67	Bh	23	Total 23	O 23	0
67	Bi	10	Total 10	O 10	0
67	Bj	3	Total 3	O 3	0
67	Bk	3	Total 3	O 3	0
67	Bl	8	Total 8	O 8	0
67	Bm	1	Total 1	O 1	0
67	Bn	1	Total 1	O 1	0
67	Bo	4	Total 4	O 4	0

### 3 Residue-property plots [i](#)

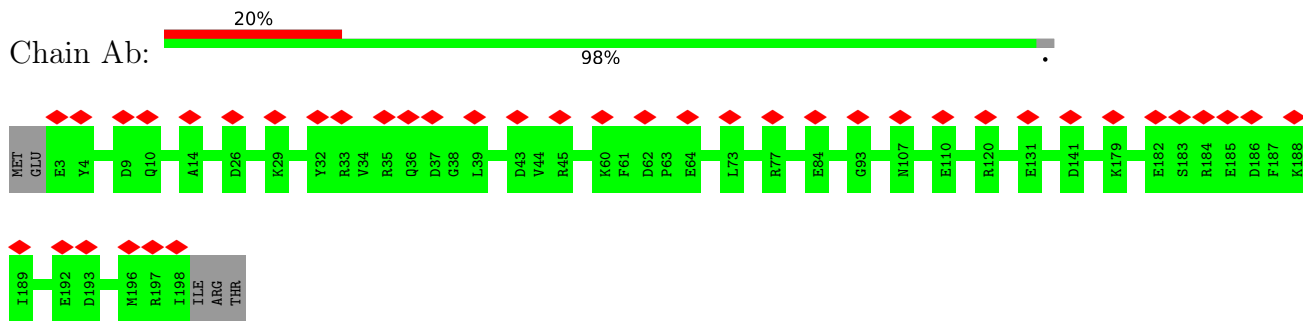
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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 ribosomal RNA

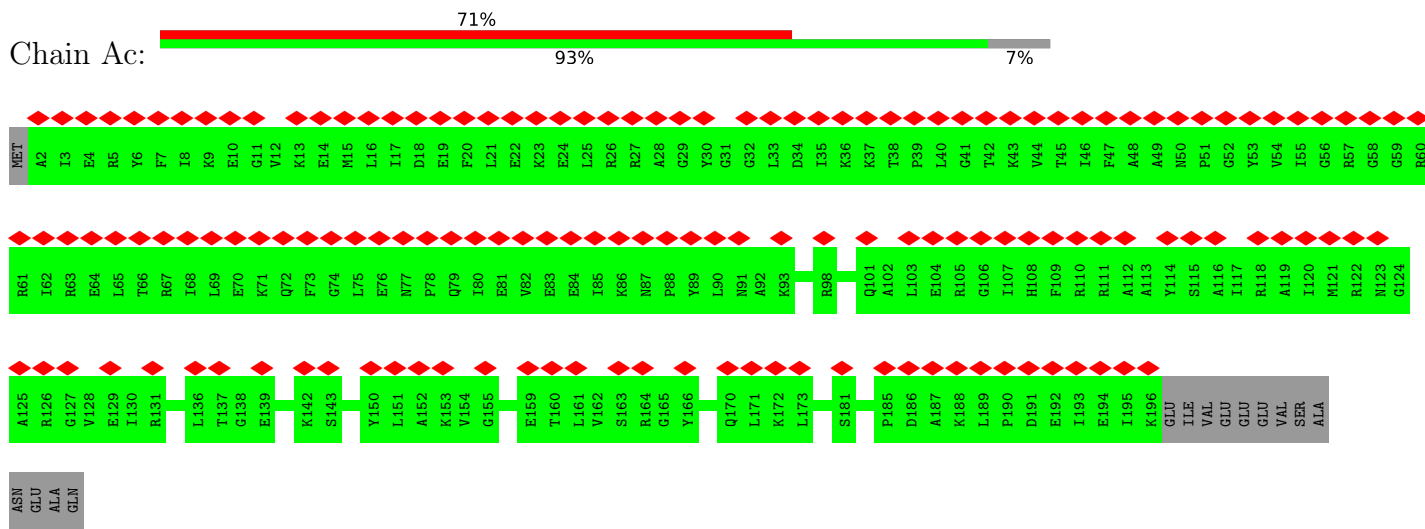




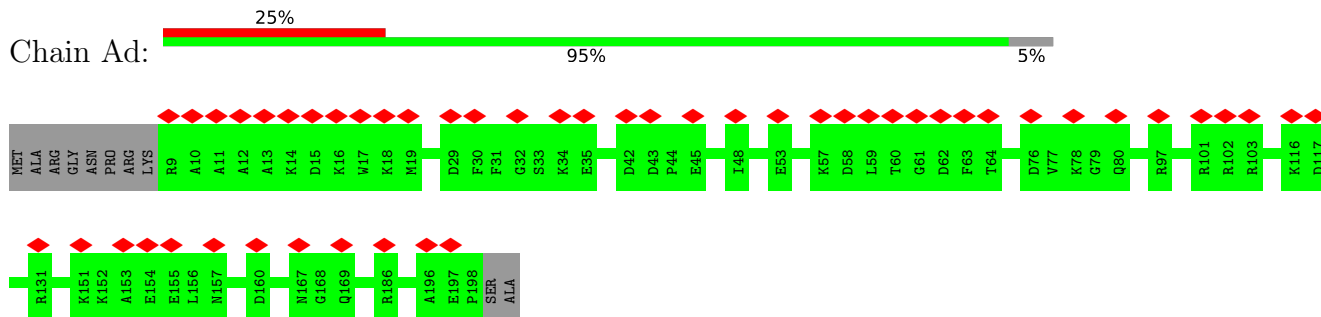
• Molecule 2: 30S ribosomal protein S2



• Molecule 3: 30S ribosomal protein S3



• Molecule 4: 30S ribosomal protein S3Ae

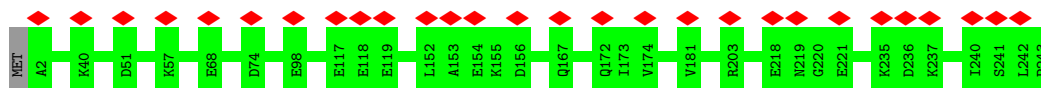


• Molecule 5: 30S ribosomal protein S4

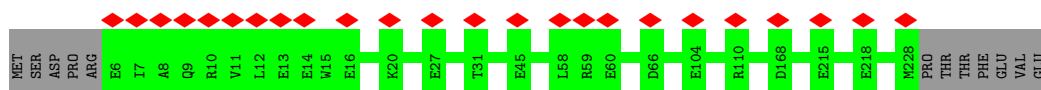




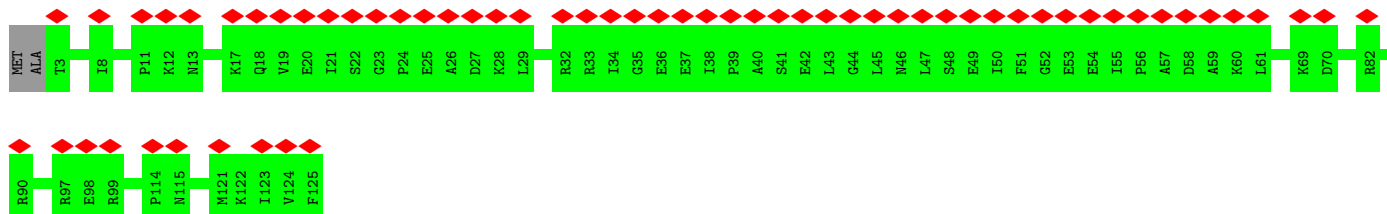
- Molecule 6: 30S ribosomal protein S4e



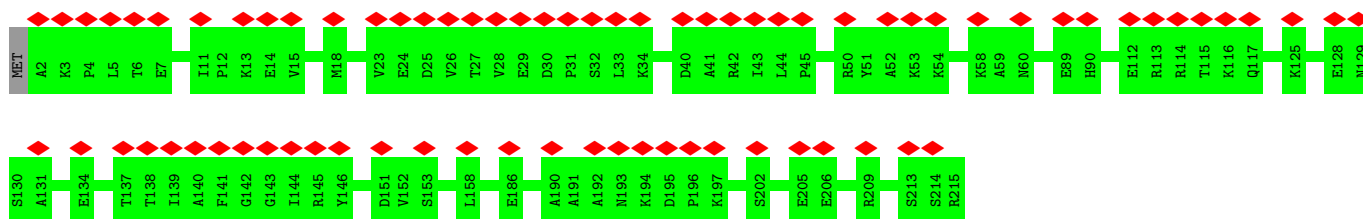
- Molecule 7: 30S ribosomal protein S5



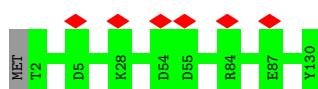
- Molecule 8: 30S ribosomal protein S6e



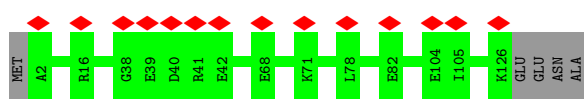
- Molecule 9: 30S ribosomal protein S7



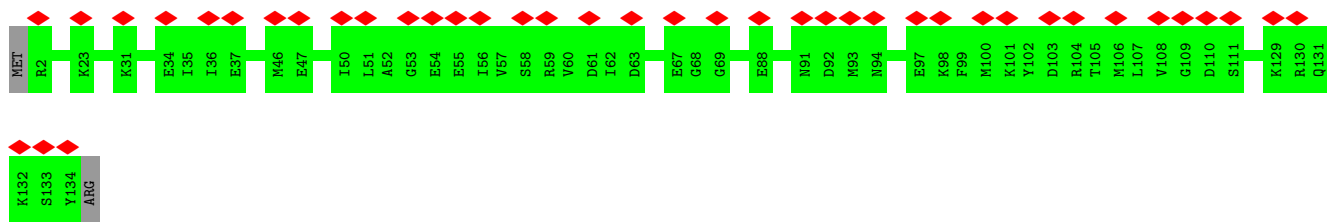
- Molecule 10: 30S ribosomal protein S8



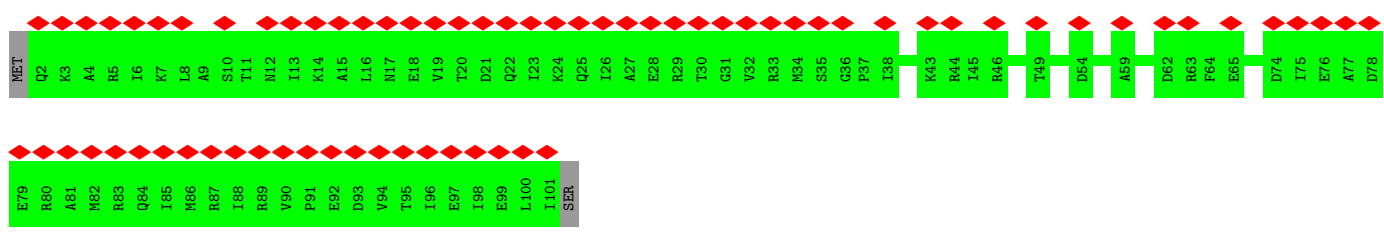
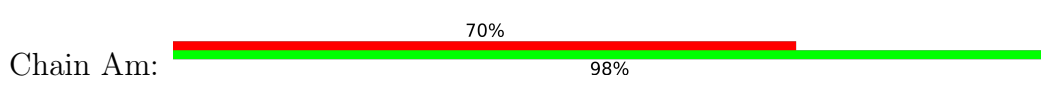
- Molecule 11: 30S ribosomal protein S8e



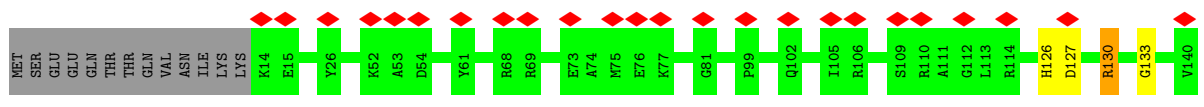
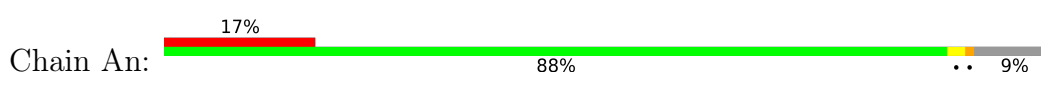
• Molecule 12: 30S ribosomal protein S9



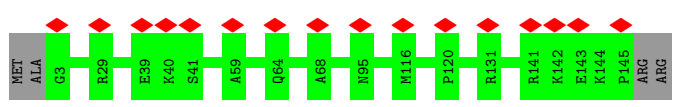
• Molecule 13: 30S ribosomal protein S10



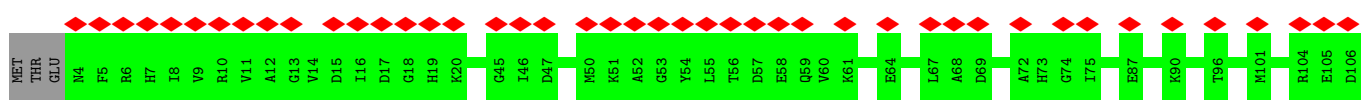
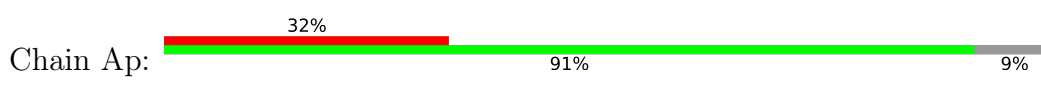
• Molecule 14: 30S ribosomal protein S11



• Molecule 15: 30S ribosomal protein S12

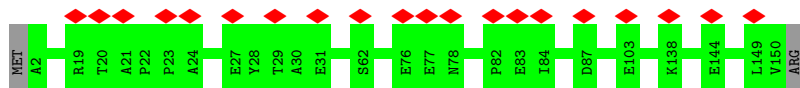


• Molecule 16: 30S ribosomal protein S13

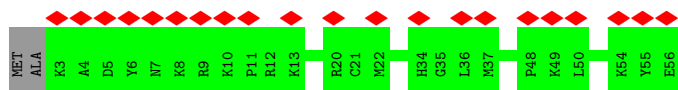
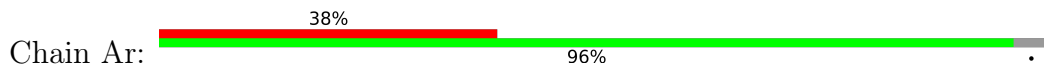




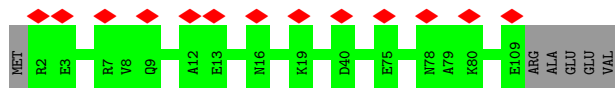
- Molecule 17: 30S ribosomal protein S15



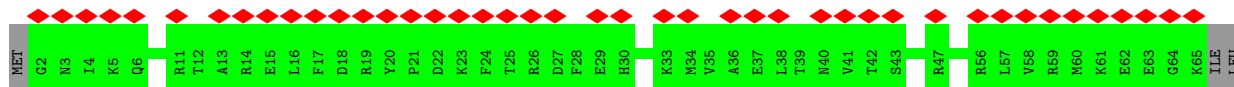
- Molecule 18: 30S ribosomal protein S14 type Z



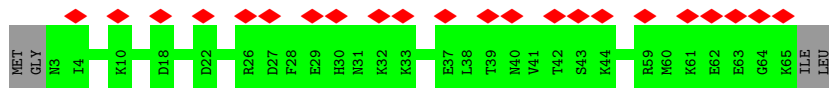
- Molecule 19: 30S ribosomal protein S17



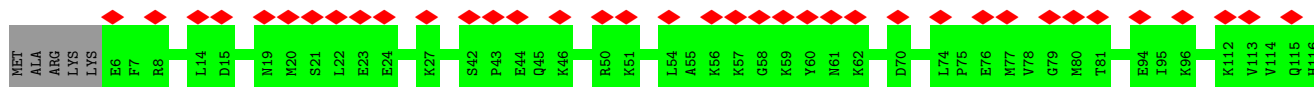
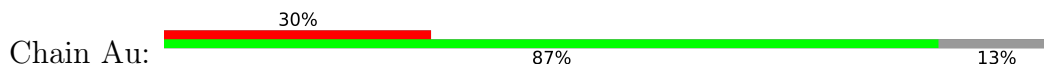
- Molecule 20: 30S ribosomal protein S17e

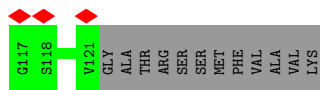


- Molecule 20: 30S ribosomal protein S17e

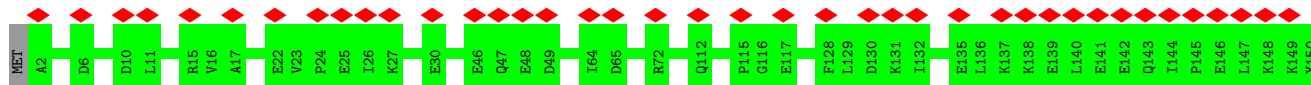


- Molecule 21: 30S ribosomal protein S19

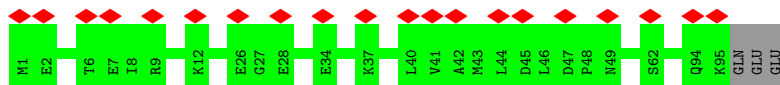




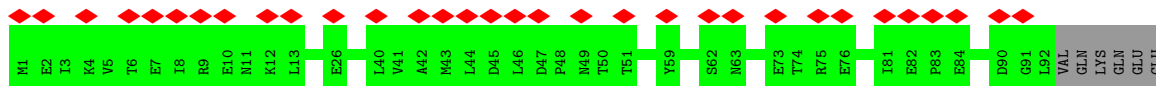
- Molecule 22: 30S ribosomal protein S19e



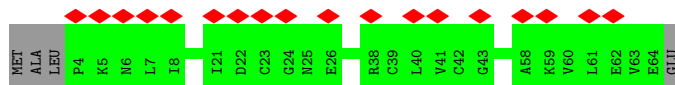
- Molecule 23: 30S ribosomal protein S24e



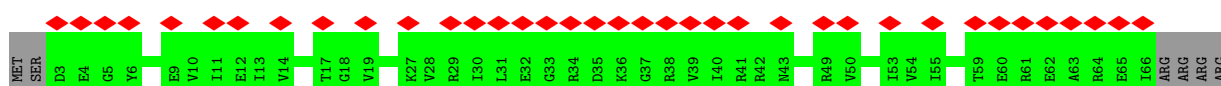
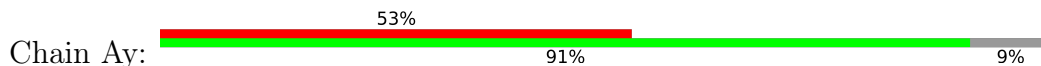
- Molecule 23: 30S ribosomal protein S24e



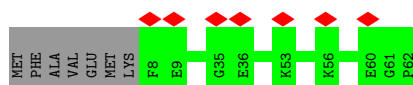
- Molecule 24: 30S ribosomal protein S27e



- Molecule 25: 30S ribosomal protein S28e

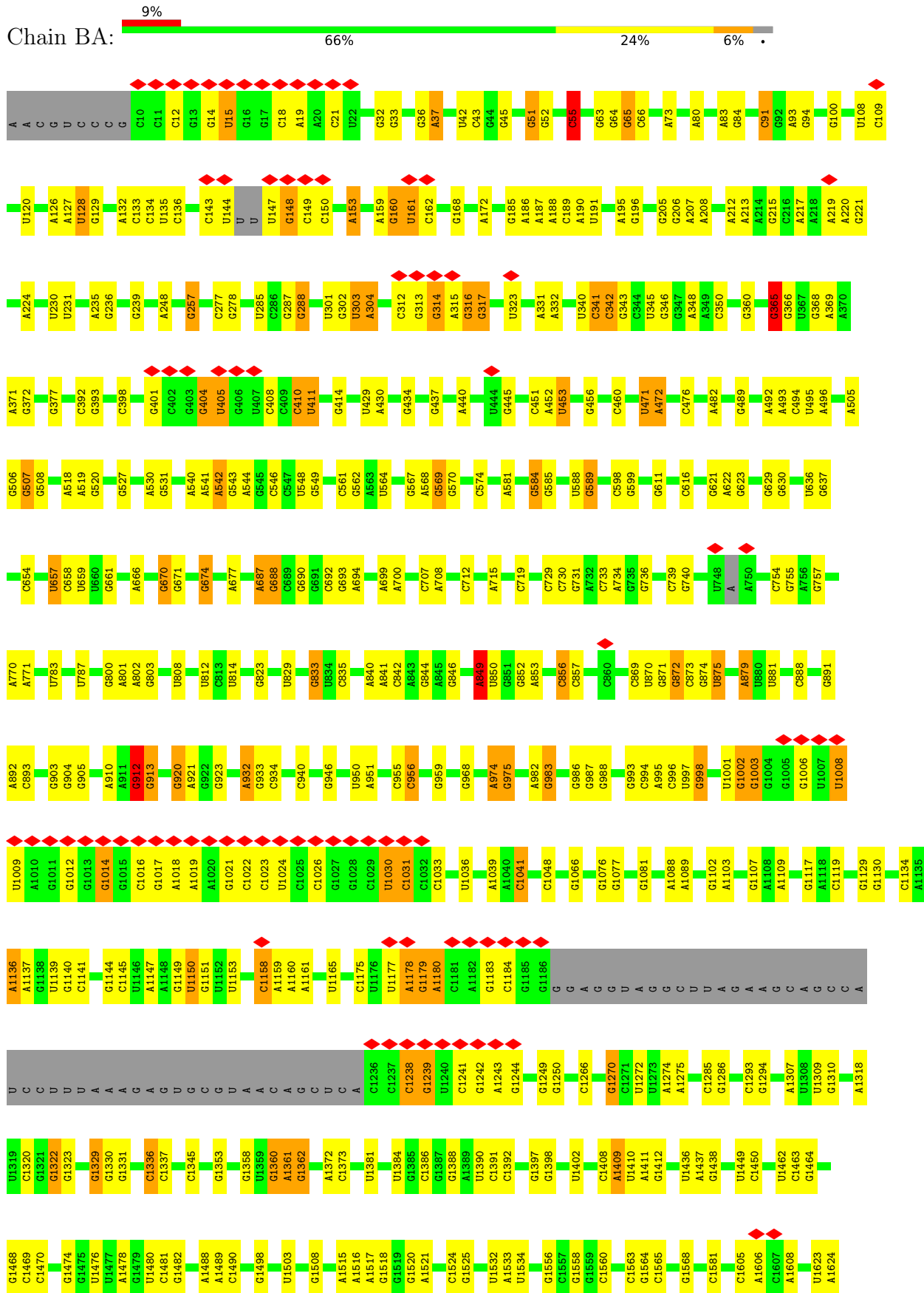


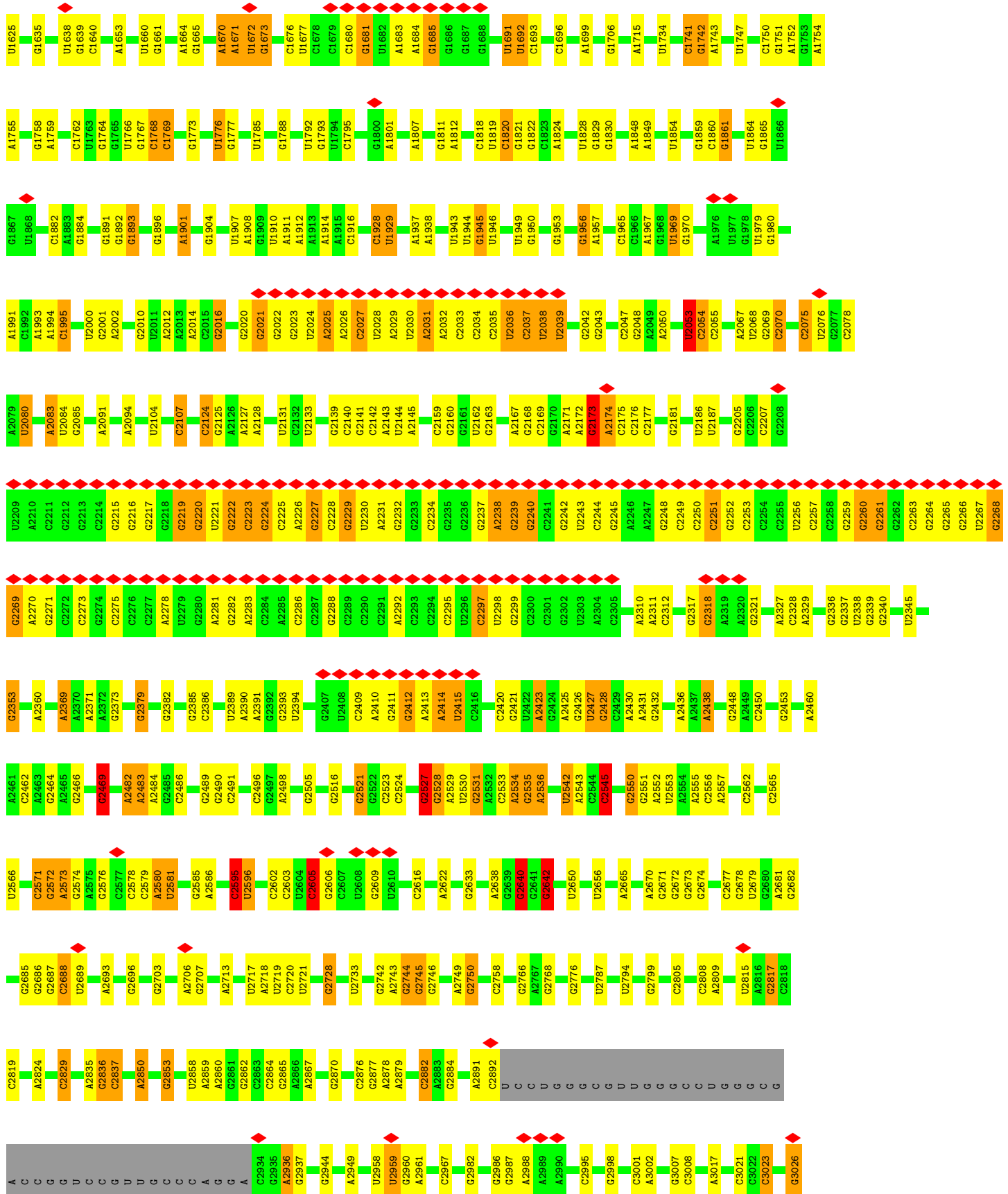
- Molecule 26: Predicted zinc-ribbon RNA-binding protein involved in translation

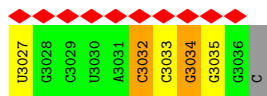


- Molecule 27: 23S ribosomal RNA

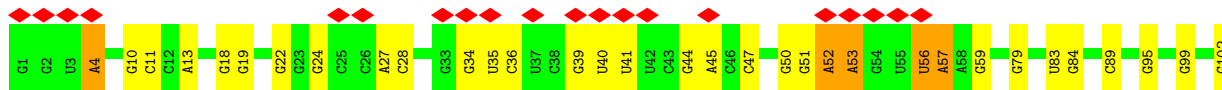
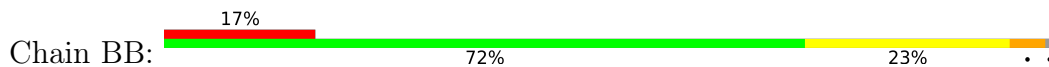








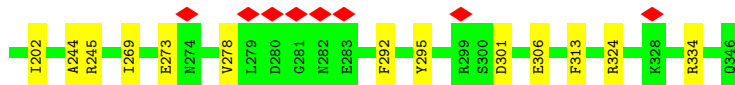
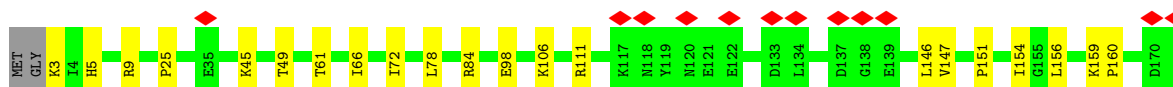
• Molecule 28: 5S ribosomal RNA



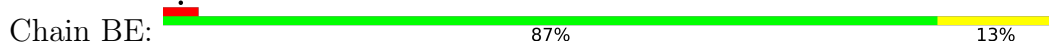
• Molecule 29: 50S ribosomal protein L2



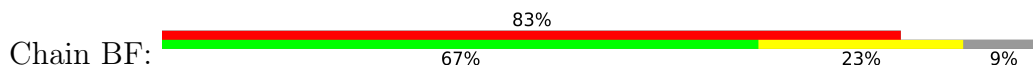
• Molecule 30: 50S ribosomal protein L3

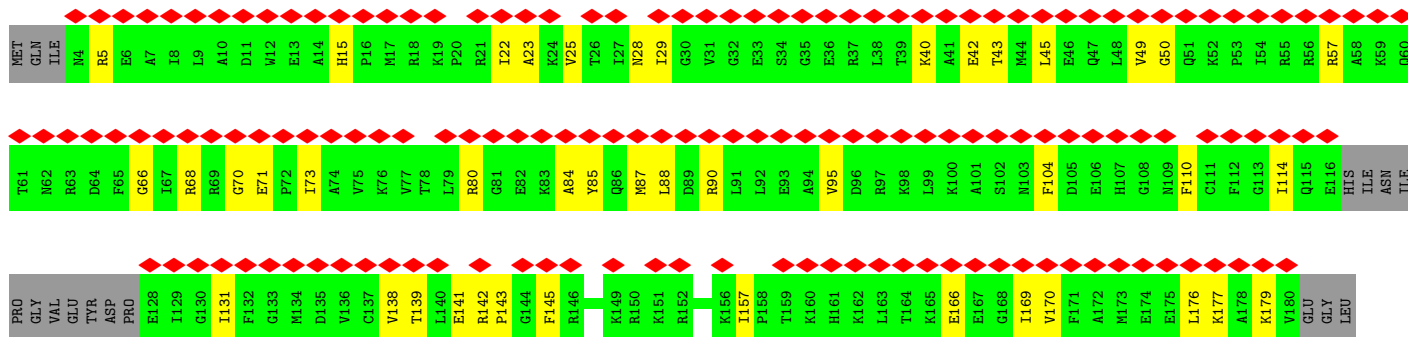


• Molecule 31: 50S ribosomal protein L4

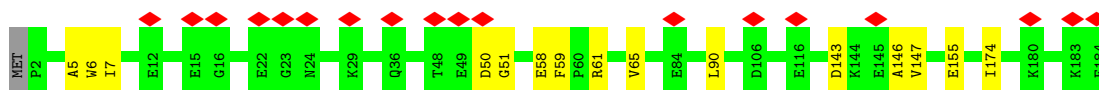


• Molecule 32: 50S ribosomal protein L5

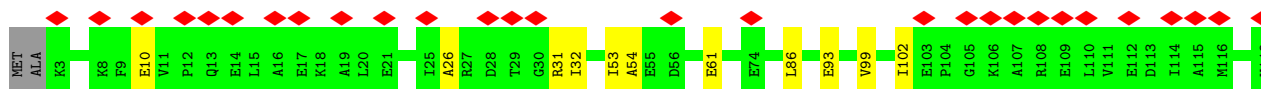
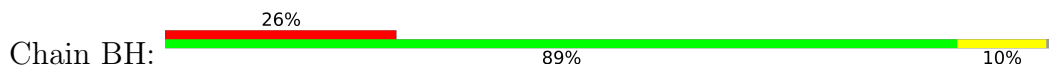




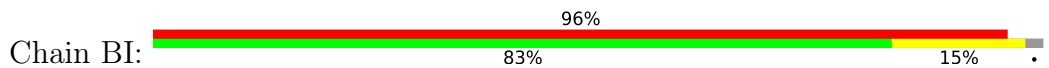
• Molecule 33: 50S ribosomal protein L6



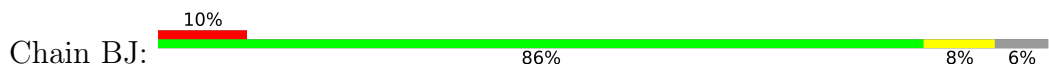
• Molecule 34: 50S ribosomal protein L7Ae

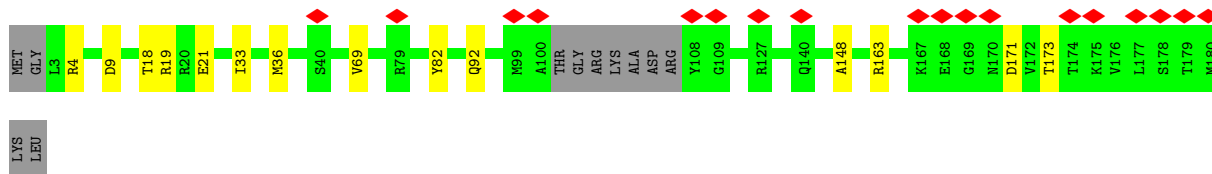


• Molecule 34: 50S ribosomal protein L7Ae

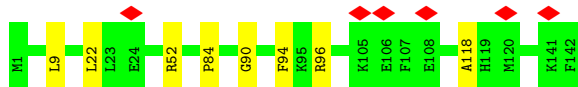


• Molecule 35: 50S ribosomal protein L10e

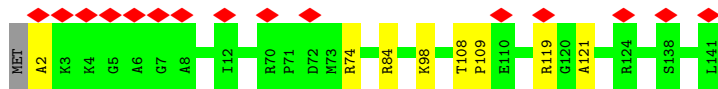




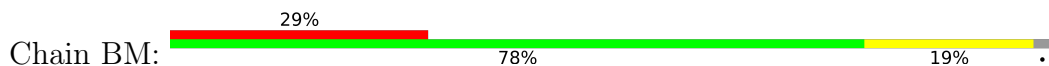
• Molecule 36: 50S ribosomal protein L13



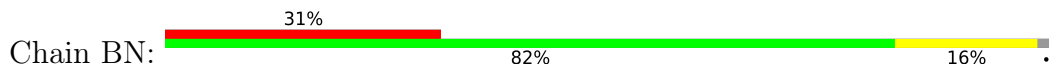
• Molecule 37: 50S ribosomal protein L14



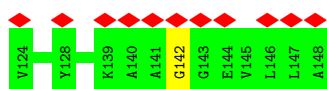
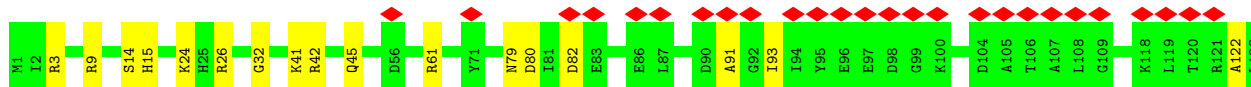
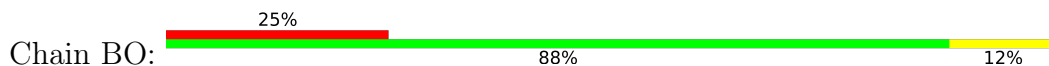
• Molecule 38: 50S ribosomal protein L14e



• Molecule 38: 50S ribosomal protein L14e



• Molecule 39: 50S ribosomal protein L15



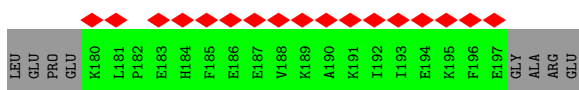
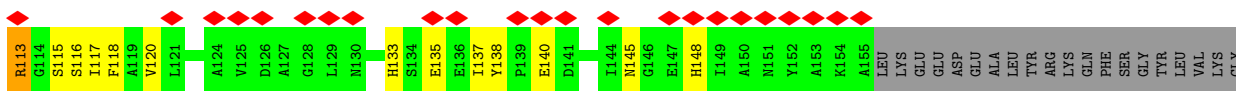
• Molecule 40: 50S ribosomal protein L15e

Chain BP:  95% 5%




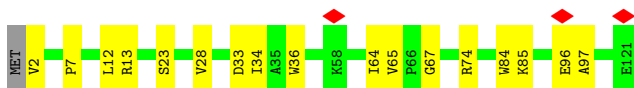
- Molecule 41: 50S ribosomal protein L18

Chain BQ:  32% 66% 15% 18%




- Molecule 42: 50S ribosomal protein L18e

Chain BR:  85% 14%



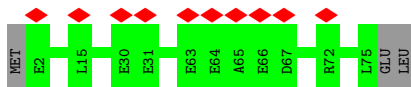
- Molecule 43: 50S ribosomal protein L19e

Chain BS:  5% 88% 9%




- Molecule 44: 50S ribosomal protein L18Ae

Chain BT:  13% 96%

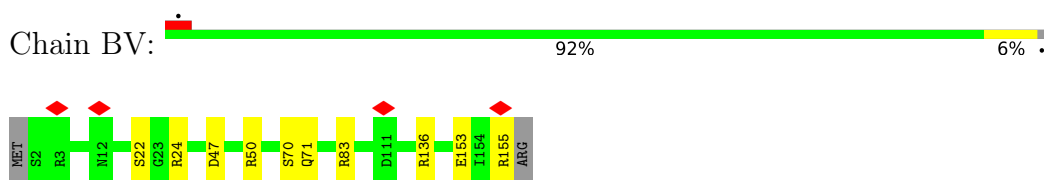


- Molecule 45: 50S ribosomal protein L21e

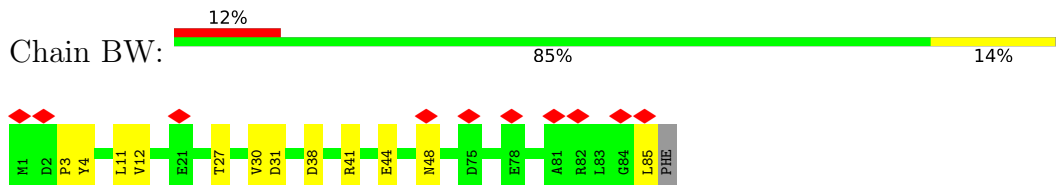
Chain BU:  89% 9%



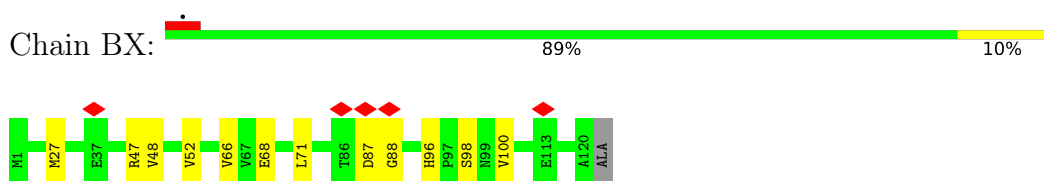
- Molecule 46: 50S ribosomal protein L22



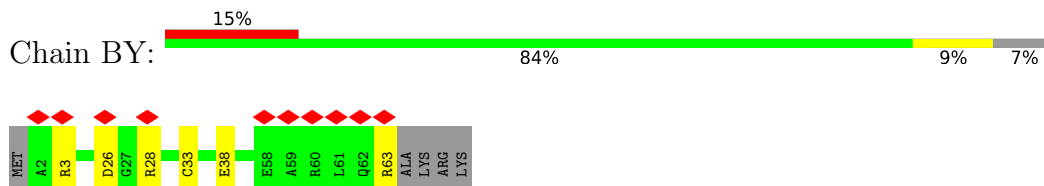
- Molecule 47: 50S ribosomal protein L23



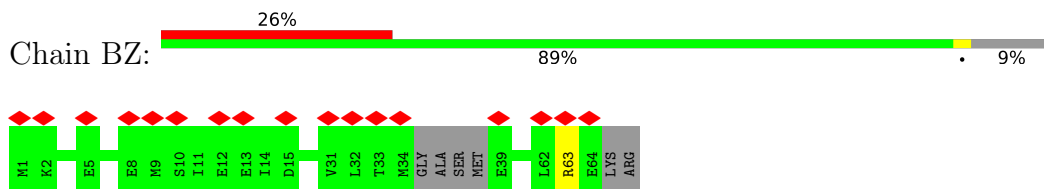
- Molecule 48: 50S ribosomal protein L24



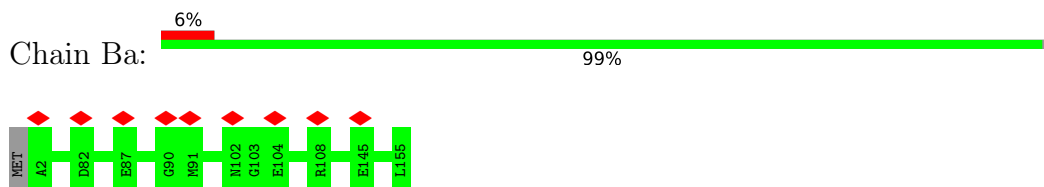
- Molecule 49: 50S ribosomal protein L24e



- Molecule 50: 50S ribosomal protein L29

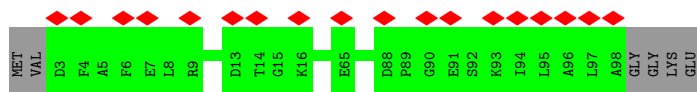


- Molecule 51: 50S ribosomal protein L30

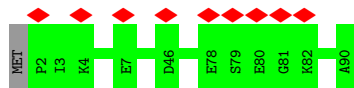


- Molecule 52: 50S ribosomal protein L30e

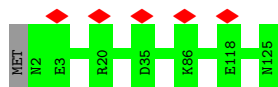




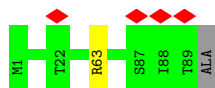
- Molecule 53: 50S ribosomal protein L31e



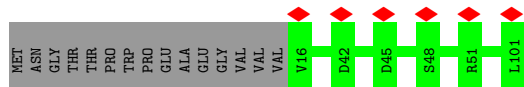
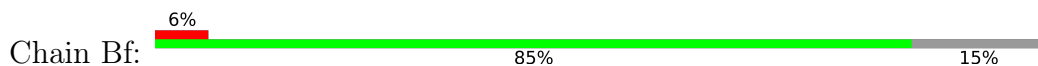
- Molecule 54: 50S ribosomal protein L32e



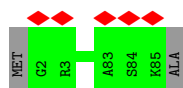
- Molecule 55: 50S ribosomal protein L34e



- Molecule 56: Ribosomal protein eL35A



- Molecule 57: 50S ribosomal protein L37Ae

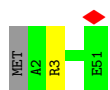


- Molecule 58: 50S ribosomal protein L37e

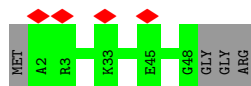


- Molecule 59: 50S ribosomal protein L39e

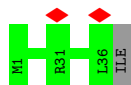




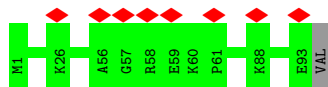
- Molecule 60: 50S ribosomal protein L40e



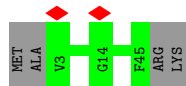
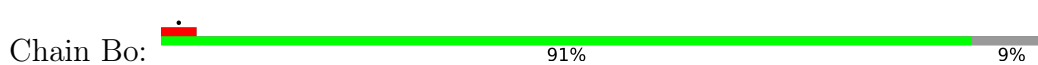
- Molecule 61: LSU ribosomal protein L41E



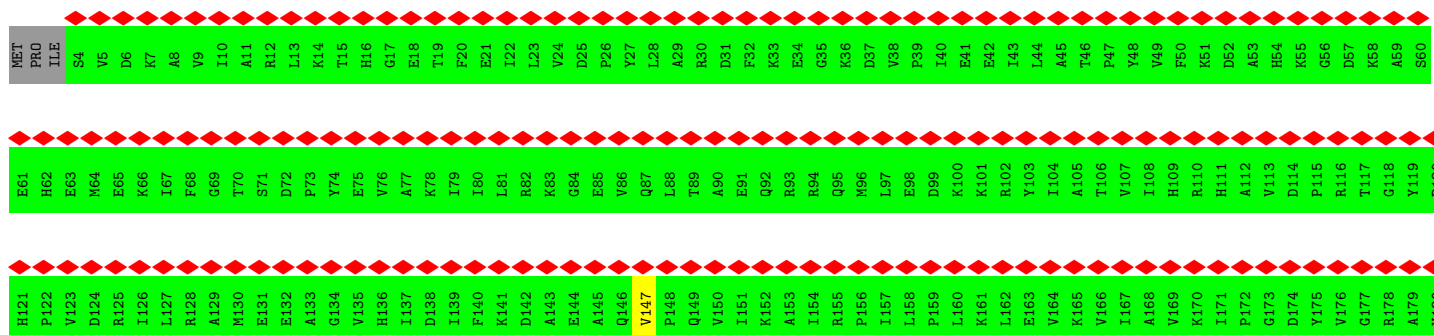
- Molecule 62: 50S ribosomal protein L44e

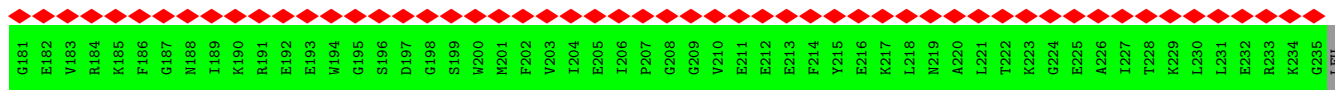


- Molecule 63: Nucleic acid-binding protein, containing C2H2 zinc-finger

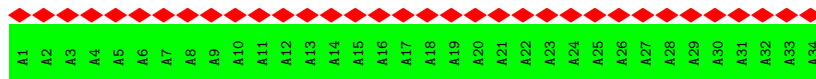


- Molecule 64: Predicted exosome subunit, UPF0023 family





• Molecule 65: Unknown ribosomal protein



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	116585	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	34	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	29000	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.497	Depositor
Minimum map value	-0.266	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.012	Depositor
Recommended contour level	0.05	Depositor
Map size (Å)	408.0, 408.0, 408.0	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.85, 0.85, 0.85	Depositor

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: A2M, 2MG, 4SU, MA6, 5MC, OMU, OMC, OMG, LV2, ZN, 6MZ, 5MU

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	Aa	0.38	0/33690	0.81	17/52572 (0.0%)
2	Ab	0.28	0/1611	0.45	0/2179
3	Ac	0.28	0/1558	0.44	0/2089
4	Ad	0.28	0/1578	0.48	0/2119
5	Ae	0.29	0/1477	0.48	0/1980
6	Af	0.30	0/1999	0.48	0/2700
7	Ag	0.29	0/1779	0.48	0/2396
8	Ah	0.31	0/968	0.49	0/1300
9	Ai	0.27	0/1748	0.45	0/2347
10	Aj	0.30	0/1039	0.48	0/1397
11	Ak	0.30	0/991	0.53	0/1322
12	Al	0.25	0/1068	0.50	0/1430
13	Am	0.25	0/810	0.45	0/1087
14	An	0.43	1/971 (0.1%)	0.53	0/1308
15	Ao	0.29	0/1139	0.46	0/1518
16	Ap	0.28	0/1113	0.48	0/1500
17	Aq	0.30	0/1241	0.45	0/1667
18	Ar	0.27	0/457	0.45	0/602
19	As	0.30	0/899	0.47	0/1214
20	At	0.26	0/545	0.42	0/725
20	Bm	0.28	0/541	0.45	0/720
21	Au	0.28	0/970	0.44	0/1295
22	Av	0.28	0/1249	0.43	0/1687
23	Aw	0.30	0/790	0.46	0/1063
23	Bn	0.29	0/765	0.46	0/1030
24	Ax	0.27	0/469	0.45	0/633
25	Ay	0.26	0/501	0.48	0/672
26	Az	0.30	0/440	0.45	0/599
27	BA	0.49	1/68729 (0.0%)	0.85	15/107278 (0.0%)
28	BB	0.33	0/2993	0.79	0/4668
29	BC	0.35	0/1860	0.53	0/2511
30	BD	0.33	0/2815	0.51	0/3795

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z  >5	RMSZ	# Z  >5
31	BE	0.32	0/2066	0.49	0/2785
32	BF	0.27	0/1356	0.52	0/1806
33	BG	0.30	0/1490	0.48	0/2006
34	BH	0.28	0/941	0.44	0/1272
34	BI	0.26	0/941	0.44	0/1272
35	BJ	0.32	0/1429	0.48	0/1921
36	BK	0.32	0/1165	0.54	0/1558
37	BL	0.30	0/1067	0.52	0/1435
38	BM	0.33	0/614	0.52	0/824
38	BN	0.28	0/614	0.53	0/824
39	BO	0.29	0/1188	0.52	0/1585
40	BP	0.36	0/1621	0.52	0/2163
41	BQ	0.29	0/1342	0.48	0/1804
42	BR	0.30	0/971	0.52	0/1301
43	BS	0.31	0/1216	0.53	0/1607
44	BT	0.29	0/636	0.46	0/852
45	BU	0.33	0/806	0.48	0/1080
46	BV	0.31	0/1259	0.49	0/1688
47	BW	0.30	0/690	0.48	0/925
48	BX	0.33	0/1007	0.50	0/1344
49	BY	0.36	0/538	0.47	0/716
50	BZ	0.28	0/506	0.49	0/668
51	Ba	0.30	0/1259	0.52	0/1680
52	Bb	0.28	0/742	0.45	0/1001
53	Bc	0.30	0/736	0.48	0/990
54	Bd	0.31	0/1044	0.49	0/1394
55	Be	0.32	0/746	0.50	0/997
56	Bf	0.33	0/681	0.55	0/919
57	Bg	0.35	0/640	0.53	0/855
58	Bh	0.38	0/524	0.53	0/692
59	Bi	0.31	0/441	0.51	0/588
60	Bj	0.34	0/381	0.56	0/505
61	Bk	0.49	0/351	0.63	0/454
62	Bl	0.33	0/796	0.53	0/1056
63	Bo	0.32	0/356	0.48	0/474
64	Bp	0.28	0/1900	0.50	1/2559 (0.0%)
65	Bq	0.30	0/169	0.44	0/235
All	All	0.40	2/173032 (0.0%)	0.73	33/255238 (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
14	An	0	2

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
14	An	130	ARG	C-N	8.62	1.50	1.34
27	BA	2173	G	O3'-P	6.46	1.69	1.61

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

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
27	BA	2173	G	P-O3'-C3'	9.17	130.70	119.70
1	Aa	1298	A	C2'-C3'-O3'	7.62	126.26	109.50
27	BA	2251	C	N3-C2-O2	-7.20	116.86	121.90
27	BA	549	G	O4'-C1'-N9	7.13	113.90	108.20
1	Aa	622	A	N9-C1'-C2'	-7.12	104.17	112.00

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
14	An	126	HIS	Peptide
14	An	127	ASP	Peptide

## 5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	Aa	31531	0	16007	0	0
2	Ab	1580	0	1634	0	0
3	Ac	1536	0	1621	0	0
4	Ad	1550	0	1618	0	0
5	Ae	1454	0	1521	0	0
6	Af	1952	0	2035	0	0
7	Ag	1753	0	1822	0	0
8	Ah	952	0	997	0	0
9	Ai	1715	0	1779	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
10	Aj	1020	0	1066	0	0
11	Ak	982	0	1071	0	0
12	Al	1054	0	1105	0	0
13	Am	803	0	854	0	0
14	An	953	0	973	0	0
15	Ao	1122	0	1212	0	0
16	Ap	1090	0	1121	0	0
17	Aq	1217	0	1302	0	0
18	Ar	447	0	469	0	0
19	As	879	0	907	0	0
20	At	538	0	560	0	0
20	Bm	534	0	557	0	0
21	Au	952	0	1000	0	0
22	Av	1216	0	1236	0	0
23	Aw	777	0	797	0	0
23	Bn	752	0	767	0	0
24	Ax	461	0	492	0	0
25	Ay	499	0	528	0	0
26	Az	426	0	402	0	0
27	BA	63209	0	32023	474	0
28	BB	2678	0	1361	18	0
29	BC	1820	0	1902	14	0
30	BD	2746	0	2840	27	0
31	BE	2026	0	2128	26	0
32	BF	1339	0	1403	31	0
33	BG	1463	0	1520	12	0
34	BH	928	0	970	7	0
34	BI	928	0	970	12	0
35	BJ	1399	0	1426	11	0
36	BK	1146	0	1213	5	0
37	BL	1055	0	1134	6	0
38	BM	610	0	660	16	0
38	BN	610	0	660	11	0
39	BO	1166	0	1229	12	0
40	BP	1582	0	1677	8	0
41	BQ	1314	0	1359	20	0
42	BR	959	0	1018	11	0
43	BS	1200	0	1298	10	0
44	BT	624	0	647	0	0
45	BU	784	0	821	9	0
46	BV	1234	0	1274	7	0
47	BW	683	0	743	8	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
48	BX	991	0	1053	7	0
49	BY	524	0	512	5	0
50	BZ	506	0	559	1	0
51	Ba	1242	0	1332	0	0
52	Bb	730	0	768	0	0
53	Bc	721	0	776	0	0
54	Bd	1022	0	1106	0	0
55	Be	728	0	786	0	0
56	Bf	669	0	714	0	0
57	Bg	630	0	675	0	0
58	Bh	512	0	527	0	0
59	Bi	432	0	488	0	0
60	Bj	373	0	390	0	0
61	Bk	345	0	407	0	0
62	Bl	778	0	831	0	0
63	Bo	349	0	358	0	0
64	Bp	1862	0	1900	0	0
65	Bq	170	0	172	0	0
66	Ag	1	0	0	0	0
66	Ar	1	0	0	0	0
66	As	1	0	0	0	0
66	Ax	1	0	0	0	0
66	Az	2	0	0	0	0
66	BY	1	0	0	0	0
66	Be	2	0	0	0	0
66	Bg	1	0	0	0	0
66	Bh	1	0	0	0	0
66	Bj	1	0	0	0	0
66	Bl	1	0	0	0	0
66	Bo	1	0	0	0	0
67	Aa	266	0	0	0	0
67	Ad	1	0	0	0	0
67	Af	4	0	0	0	0
67	Ag	1	0	0	0	0
67	Ai	1	0	0	0	0
67	Aj	2	0	0	0	0
67	Ak	6	0	0	0	0
67	Al	2	0	0	0	0
67	Am	2	0	0	0	0
67	An	2	0	0	0	0
67	Ao	1	0	0	0	0
67	Ap	3	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
67	Aq	5	0	0	0	0
67	As	2	0	0	0	0
67	Au	1	0	0	0	0
67	Av	1	0	0	0	0
67	Az	1	0	0	0	0
67	BA	2715	0	0	170	0
67	BB	23	0	0	3	0
67	BC	51	0	0	1	0
67	BD	59	0	0	1	0
67	BE	41	0	0	7	0
67	BF	2	0	0	1	0
67	BG	1	0	0	0	0
67	BH	1	0	0	0	0
67	BJ	13	0	0	0	0
67	BK	18	0	0	1	0
67	BL	15	0	0	1	0
67	BM	3	0	0	2	0
67	BN	3	0	0	1	0
67	BO	30	0	0	1	0
67	BP	36	0	0	1	0
67	BQ	6	0	0	1	0
67	BR	14	0	0	0	0
67	BS	17	0	0	0	0
67	BT	2	0	0	0	0
67	BU	13	0	0	3	0
67	BV	22	0	0	0	0
67	BW	3	0	0	0	0
67	BX	10	0	0	1	0
67	BY	12	0	0	1	0
67	BZ	7	0	0	0	0
67	Ba	14	0	0	0	0
67	Bb	8	0	0	0	0
67	Bc	10	0	0	0	0
67	Bd	23	0	0	0	0
67	Be	8	0	0	0	0
67	Bf	7	0	0	0	0
67	Bg	15	0	0	0	0
67	Bh	23	0	0	0	0
67	Bi	10	0	0	0	0
67	Bj	3	0	0	0	0
67	Bk	3	0	0	0	0
67	Bl	8	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
67	Bm	1	0	0	0	0
67	Bn	1	0	0	0	0
67	Bo	4	0	0	0	0
All	All	167402	0	119083	723	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 4.

The worst 5 of 723 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)
27:BA:2678:G:OP1	67:BA:3101:HOH:O	1.68	1.09
27:BA:1478:A:OP2	67:BA:3102:HOH:O	1.75	1.04
27:BA:1137:A:OP2	67:BA:3104:HOH:O	1.83	0.96
27:BA:1463:C:OP2	67:BA:3103:HOH:O	1.81	0.96
27:BA:168:G:N7	67:BA:3120:HOH:O	2.00	0.93

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	Ab	194/201 (96%)	181 (93%)	13 (7%)	0	100	100
3	Ac	193/209 (92%)	185 (96%)	8 (4%)	0	100	100
4	Ad	188/200 (94%)	175 (93%)	13 (7%)	0	100	100
5	Ae	172/180 (96%)	165 (96%)	7 (4%)	0	100	100
6	Af	240/243 (99%)	234 (98%)	6 (2%)	0	100	100
7	Ag	221/235 (94%)	209 (95%)	12 (5%)	0	100	100
8	Ah	121/125 (97%)	118 (98%)	3 (2%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
9	Ai	212/215 (99%)	205 (97%)	7 (3%)	0	100	100
10	Aj	127/130 (98%)	119 (94%)	8 (6%)	0	100	100
11	Ak	123/130 (95%)	116 (94%)	7 (6%)	0	100	100
12	Al	131/135 (97%)	123 (94%)	8 (6%)	0	100	100
13	Am	98/102 (96%)	94 (96%)	4 (4%)	0	100	100
14	An	125/140 (89%)	118 (94%)	6 (5%)	1 (1%)	19	29
15	Ao	141/147 (96%)	135 (96%)	6 (4%)	0	100	100
16	Ap	133/149 (89%)	118 (89%)	15 (11%)	0	100	100
17	Aq	147/151 (97%)	141 (96%)	6 (4%)	0	100	100
18	Ar	52/56 (93%)	46 (88%)	6 (12%)	0	100	100
19	As	106/114 (93%)	102 (96%)	4 (4%)	0	100	100
20	At	62/67 (92%)	60 (97%)	2 (3%)	0	100	100
20	Bm	61/67 (91%)	58 (95%)	3 (5%)	0	100	100
21	Au	114/133 (86%)	111 (97%)	3 (3%)	0	100	100
22	Av	147/150 (98%)	139 (95%)	8 (5%)	0	100	100
23	Aw	93/98 (95%)	87 (94%)	6 (6%)	0	100	100
23	Bn	90/98 (92%)	85 (94%)	5 (6%)	0	100	100
24	Ax	59/65 (91%)	56 (95%)	3 (5%)	0	100	100
25	Ay	62/70 (89%)	60 (97%)	2 (3%)	0	100	100
26	Az	53/62 (86%)	51 (96%)	2 (4%)	0	100	100
29	BC	235/239 (98%)	221 (94%)	14 (6%)	0	100	100
30	BD	342/346 (99%)	322 (94%)	20 (6%)	0	100	100
31	BE	253/255 (99%)	241 (95%)	12 (5%)	0	100	100
32	BF	162/183 (88%)	132 (82%)	30 (18%)	0	100	100
33	BG	181/184 (98%)	169 (93%)	12 (7%)	0	100	100
34	BH	119/123 (97%)	113 (95%)	6 (5%)	0	100	100
34	BI	119/123 (97%)	108 (91%)	11 (9%)	0	100	100
35	BJ	167/182 (92%)	157 (94%)	10 (6%)	0	100	100
36	BK	140/142 (99%)	136 (97%)	4 (3%)	0	100	100
37	BL	138/141 (98%)	128 (93%)	10 (7%)	0	100	100
38	BM	79/83 (95%)	70 (89%)	9 (11%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
38	BN	79/83 (95%)	72 (91%)	7 (9%)	0	100	100
39	BO	146/148 (99%)	134 (92%)	12 (8%)	0	100	100
40	BP	191/194 (98%)	182 (95%)	9 (5%)	0	100	100
41	BQ	158/201 (79%)	149 (94%)	9 (6%)	0	100	100
42	BR	118/121 (98%)	112 (95%)	6 (5%)	0	100	100
43	BS	144/150 (96%)	139 (96%)	5 (4%)	0	100	100
44	BT	72/77 (94%)	70 (97%)	2 (3%)	0	100	100
45	BU	94/98 (96%)	91 (97%)	3 (3%)	0	100	100
46	BV	152/156 (97%)	149 (98%)	3 (2%)	0	100	100
47	BW	83/86 (96%)	78 (94%)	5 (6%)	0	100	100
48	BX	118/121 (98%)	113 (96%)	5 (4%)	0	100	100
49	BY	60/67 (90%)	60 (100%)	0	0	100	100
50	BZ	56/66 (85%)	54 (96%)	2 (4%)	0	100	100
51	Ba	152/155 (98%)	147 (97%)	5 (3%)	0	100	100
52	Bb	94/102 (92%)	87 (93%)	7 (7%)	0	100	100
53	Bc	87/90 (97%)	80 (92%)	7 (8%)	0	100	100
54	Bd	122/125 (98%)	117 (96%)	5 (4%)	0	100	100
55	Be	87/90 (97%)	82 (94%)	5 (6%)	0	100	100
56	Bf	84/101 (83%)	80 (95%)	4 (5%)	0	100	100
57	Bg	82/86 (95%)	77 (94%)	5 (6%)	0	100	100
58	Bh	60/63 (95%)	57 (95%)	3 (5%)	0	100	100
59	Bi	48/51 (94%)	47 (98%)	1 (2%)	0	100	100
60	Bj	45/51 (88%)	43 (96%)	2 (4%)	0	100	100
61	Bk	34/37 (92%)	34 (100%)	0	0	100	100
62	Bl	91/94 (97%)	90 (99%)	1 (1%)	0	100	100
63	Bo	41/47 (87%)	38 (93%)	3 (7%)	0	100	100
64	Bp	230/236 (98%)	213 (93%)	17 (7%)	0	100	100
65	Bq	32/34 (94%)	30 (94%)	2 (6%)	0	100	100
All	All	8160/8603 (95%)	7713 (94%)	446 (6%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
14	An	133	GLY

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	Ab	170/175 (97%)	170 (100%)	0	100	100
3	Ac	156/168 (93%)	156 (100%)	0	100	100
4	Ad	162/169 (96%)	162 (100%)	0	100	100
5	Ae	153/156 (98%)	153 (100%)	0	100	100
6	Af	211/212 (100%)	211 (100%)	0	100	100
7	Ag	185/197 (94%)	185 (100%)	0	100	100
8	Ah	104/105 (99%)	104 (100%)	0	100	100
9	Ai	183/184 (100%)	183 (100%)	0	100	100
10	Aj	106/107 (99%)	106 (100%)	0	100	100
11	Ak	101/105 (96%)	101 (100%)	0	100	100
12	Al	110/112 (98%)	110 (100%)	0	100	100
13	Am	89/91 (98%)	89 (100%)	0	100	100
14	An	95/108 (88%)	94 (99%)	1 (1%)	73	85
15	Ao	117/120 (98%)	117 (100%)	0	100	100
16	Ap	111/123 (90%)	111 (100%)	0	100	100
17	Aq	129/131 (98%)	129 (100%)	0	100	100
18	Ar	45/46 (98%)	45 (100%)	0	100	100
19	As	96/101 (95%)	96 (100%)	0	100	100
20	At	58/61 (95%)	58 (100%)	0	100	100
20	Bm	58/61 (95%)	58 (100%)	0	100	100
21	Au	104/117 (89%)	104 (100%)	0	100	100
22	Av	125/126 (99%)	125 (100%)	0	100	100
23	Aw	85/88 (97%)	85 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
23	Bn	82/88 (93%)	82 (100%)	0	100	100
24	Ax	53/56 (95%)	53 (100%)	0	100	100
25	Ay	53/59 (90%)	53 (100%)	0	100	100
26	Az	49/55 (89%)	49 (100%)	0	100	100
29	BC	185/187 (99%)	185 (100%)	0	100	100
30	BD	288/289 (100%)	288 (100%)	0	100	100
31	BE	212/212 (100%)	212 (100%)	0	100	100
32	BF	140/155 (90%)	138 (99%)	2 (1%)	67	81
33	BG	158/159 (99%)	158 (100%)	0	100	100
34	BH	99/100 (99%)	99 (100%)	0	100	100
34	BI	99/100 (99%)	98 (99%)	1 (1%)	76	86
35	BJ	145/153 (95%)	145 (100%)	0	100	100
36	BK	120/120 (100%)	120 (100%)	0	100	100
37	BL	104/105 (99%)	104 (100%)	0	100	100
38	BM	63/65 (97%)	63 (100%)	0	100	100
38	BN	63/65 (97%)	63 (100%)	0	100	100
39	BO	116/116 (100%)	116 (100%)	0	100	100
40	BP	161/162 (99%)	161 (100%)	0	100	100
41	BQ	138/167 (83%)	137 (99%)	1 (1%)	84	91
42	BR	101/102 (99%)	101 (100%)	0	100	100
43	BS	120/124 (97%)	120 (100%)	0	100	100
44	BT	69/72 (96%)	69 (100%)	0	100	100
45	BU	84/85 (99%)	84 (100%)	0	100	100
46	BV	129/131 (98%)	129 (100%)	0	100	100
47	BW	75/76 (99%)	75 (100%)	0	100	100
48	BX	109/109 (100%)	109 (100%)	0	100	100
49	BY	54/58 (93%)	54 (100%)	0	100	100
50	BZ	57/61 (93%)	57 (100%)	0	100	100
51	Ba	132/133 (99%)	132 (100%)	0	100	100
52	Bb	76/80 (95%)	76 (100%)	0	100	100
53	Bc	75/76 (99%)	75 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
54	Bd	106/107 (99%)	106 (100%)	0	100	100
55	Be	80/80 (100%)	79 (99%)	1 (1%)	69	82
56	Bf	71/83 (86%)	71 (100%)	0	100	100
57	Bg	62/63 (98%)	62 (100%)	0	100	100
58	Bh	50/51 (98%)	50 (100%)	0	100	100
59	Bi	45/46 (98%)	44 (98%)	1 (2%)	52	70
60	Bj	36/38 (95%)	36 (100%)	0	100	100
61	Bk	34/35 (97%)	34 (100%)	0	100	100
62	Bl	83/84 (99%)	83 (100%)	0	100	100
63	Bo	36/39 (92%)	36 (100%)	0	100	100
64	Bp	196/200 (98%)	196 (100%)	0	100	100
All	All	6961/7209 (97%)	6954 (100%)	7 (0%)	93	97

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
34	BI	67	PRO
41	BQ	113	ARG
59	Bi	3	ARG
55	Be	63	ARG
32	BF	179	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 13 such sidechains are listed below:

Mol	Chain	Res	Type
35	BJ	141	HIS
41	BQ	42	ASN
48	BX	17	ASN
45	BU	96	GLN
46	BV	12	ASN

### 5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	Aa	1453/1498 (96%)	240 (16%)	0
27	BA	2929/3037 (96%)	488 (16%)	36 (1%)

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Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
28	BB	124/126 (98%)	15 (12%)	3 (2%)
All	All	4506/4661 (96%)	743 (16%)	39 (0%)

5 of 743 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	Aa	4	C
1	Aa	17	5MC
1	Aa	33	U
1	Aa	42	G
1	Aa	45	U

5 of 39 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
27	BA	2482	A
27	BA	3023	C
27	BA	2534	A
27	BA	2595	OMC
28	BB	52	A

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

145 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
27	5MC	BA	2688	27	18,22,23	3.55	7 (38%)	26,32,35	1.04	1 (3%)
27	A2M	BA	2329	27	18,25,26	0.99	1 (5%)	18,36,39	1.18	2 (11%)
1	OMU	Aa	8	1	19,22,23	3.02	8 (42%)	26,31,34	1.71	5 (19%)
1	OMU	Aa	753	1	19,22,23	3.05	8 (42%)	26,31,34	1.70	4 (15%)
1	OMG	Aa	219	1	18,26,27	2.48	8 (44%)	19,38,41	1.48	4 (21%)
1	2MG	Aa	995	1	18,26,27	0.89	1 (5%)	16,38,41	1.14	2 (12%)



Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
27	OMU	BA	1150	27	19,22,23	3.03	8 (42%)	26,31,34	1.70	4 (15%)
27	OMG	BA	2728	27	18,26,27	2.48	8 (44%)	19,38,41	1.50	4 (21%)
1	OMC	Aa	825	1	19,22,23	2.97	8 (42%)	26,31,34	0.73	0
1	OMG	Aa	636	1	18,26,27	2.50	8 (44%)	19,38,41	1.48	4 (21%)
1	5MC	Aa	1321	1	18,22,23	3.51	7 (38%)	26,32,35	1.15	2 (7%)
1	OMG	Aa	121	1	18,26,27	2.48	8 (44%)	19,38,41	1.50	4 (21%)
27	5MC	BA	2075	27	18,22,23	3.51	7 (38%)	26,32,35	1.04	2 (7%)
27	OMC	BA	2595	27	19,22,23	3.03	8 (42%)	26,31,34	0.92	1 (3%)
1	5MC	Aa	17	1	18,22,23	3.51	7 (38%)	26,32,35	1.01	2 (7%)
27	A2M	BA	879	27	18,25,26	4.24	7 (38%)	18,36,39	2.99	3 (16%)
27	OMG	BA	2353	27	18,26,27	1.05	1 (5%)	19,38,41	1.03	2 (10%)
27	OMG	BA	1956	27	18,26,27	2.49	8 (44%)	19,38,41	1.52	4 (21%)
1	OMG	Aa	994	1	18,26,27	2.51	8 (44%)	19,38,41	1.51	4 (21%)
1	OMG	Aa	495	1	18,26,27	2.50	8 (44%)	19,38,41	1.52	4 (21%)
1	MA6	Aa	1467	1	19,26,27	1.03	2 (10%)	18,38,41	3.61	2 (11%)
1	OMG	Aa	963	1	18,26,27	2.53	8 (44%)	19,38,41	1.47	4 (21%)
1	OMG	Aa	1107	1	18,26,27	2.51	8 (44%)	19,38,41	1.50	4 (21%)
1	5MC	Aa	1004	1	18,22,23	3.59	7 (38%)	26,32,35	1.01	2 (7%)
1	5MC	Aa	1007	1	18,22,23	3.57	7 (38%)	26,32,35	1.05	1 (3%)
27	OMG	BA	1508	27	18,26,27	2.50	8 (44%)	19,38,41	1.51	4 (21%)
27	OMC	BA	2107	27	19,22,23	2.99	8 (42%)	26,31,34	0.77	0
1	5MC	Aa	1306	1	18,22,23	3.55	7 (38%)	26,32,35	1.01	2 (7%)
27	OMU	BA	1480	27	19,22,23	3.04	8 (42%)	26,31,34	1.70	4 (15%)
27	OMG	BA	2521	27	18,26,27	2.49	8 (44%)	19,38,41	1.56	4 (21%)
27	OMG	BA	2528	27	18,26,27	2.46	8 (44%)	19,38,41	1.48	4 (21%)
27	OMG	BA	833	27	18,26,27	2.49	8 (44%)	19,38,41	1.50	4 (21%)
27	OMC	BA	869	27	19,22,23	2.97	8 (42%)	26,31,34	0.70	0
1	OMG	Aa	450	1	18,26,27	1.05	1 (5%)	19,38,41	1.11	2 (10%)
1	OMC	Aa	1354	1	19,22,23	2.97	8 (42%)	26,31,34	0.77	0
1	OMG	Aa	446	1	18,26,27	2.48	8 (44%)	19,38,41	1.56	4 (21%)
27	A2M	BA	932	27	18,25,26	4.28	7 (38%)	18,36,39	2.98	3 (16%)
27	OMU	BA	2656	27	19,22,23	3.03	8 (42%)	26,31,34	1.70	4 (15%)
1	5MC	Aa	923	1	18,22,23	3.54	7 (38%)	26,32,35	1.08	2 (7%)
27	OMG	BA	912	27	18,26,27	2.47	8 (44%)	19,38,41	1.50	4 (21%)
27	OMU	BA	2542	27	19,22,23	3.01	8 (42%)	26,31,34	1.66	4 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
27	OMC	BA	342	27	19,22,23	0.84	0	26,31,34	0.98	1 (3%)
1	OMG	Aa	498	1	18,26,27	2.50	8 (44%)	19,38,41	1.55	4 (21%)
1	5MC	Aa	1006	1	18,22,23	3.57	7 (38%)	26,32,35	1.04	2 (7%)
27	OMG	BA	2527	27	18,26,27	2.48	8 (44%)	19,38,41	1.48	4 (21%)
27	4SU	BA	2581	27	18,21,22	4.24	8 (44%)	26,30,33	2.23	5 (19%)
27	OMG	BA	1525	27	18,26,27	2.46	8 (44%)	19,38,41	1.50	4 (21%)
27	A2M	BA	849	27	18,25,26	4.19	7 (38%)	18,36,39	3.00	3 (16%)
27	OMG	BA	2672	27	18,26,27	2.47	8 (44%)	19,38,41	1.52	4 (21%)
1	5MU	Aa	917	1	19,22,23	1.36	4 (21%)	28,32,35	2.15	6 (21%)
1	5MC	Aa	1266	1	18,22,23	3.56	7 (38%)	26,32,35	1.02	2 (7%)
27	OMC	BA	2545	27	19,22,23	3.02	8 (42%)	26,31,34	0.76	0
1	5MC	Aa	672	1	18,22,23	3.54	7 (38%)	26,32,35	1.02	2 (7%)
27	OMC	BA	1820	27	19,22,23	3.00	8 (42%)	26,31,34	0.71	0
1	5MC	Aa	1003	1	18,22,23	3.57	7 (38%)	26,32,35	1.01	2 (7%)
1	OMC	Aa	229	1	19,22,23	3.00	8 (42%)	26,31,34	0.78	0
27	4SU	BA	2553	27	18,21,22	1.75	4 (22%)	26,30,33	2.19	5 (19%)
27	OMG	BA	674	27	18,26,27	2.47	8 (44%)	19,38,41	1.54	4 (21%)
27	5MU	BA	1854	27	19,22,23	4.88	7 (36%)	28,32,35	3.73	9 (32%)
1	5MC	Aa	1478	1	18,22,23	3.58	7 (38%)	26,32,35	1.11	2 (7%)
27	OMG	BA	288	27	18,26,27	2.51	8 (44%)	19,38,41	1.53	4 (21%)
1	5MC	Aa	1476	1	18,22,23	3.52	7 (38%)	26,32,35	0.99	2 (7%)
27	5MC	BA	2605	27	18,22,23	3.55	7 (38%)	26,32,35	1.12	2 (7%)
27	OMC	BA	2124	27	19,22,23	2.97	8 (42%)	26,31,34	0.66	0
27	A2M	BA	2091	27	18,25,26	0.92	1 (5%)	18,36,39	1.19	2 (11%)
27	OMG	BA	65	27	18,26,27	2.46	8 (44%)	19,38,41	1.51	4 (21%)
27	OMG	BA	2640	27	18,26,27	2.50	8 (44%)	19,38,41	1.52	4 (21%)
1	OMU	Aa	1020	1	19,22,23	3.04	8 (42%)	26,31,34	1.67	5 (19%)
1	OMU	Aa	1358	1	19,22,23	3.00	8 (42%)	26,31,34	1.71	5 (19%)
27	OMG	BA	1329	27	18,26,27	2.47	8 (44%)	19,38,41	1.59	4 (21%)
27	A2M	BA	2936	27	18,25,26	4.23	7 (38%)	18,36,39	2.92	3 (16%)
1	5MC	Aa	1352	1	18,22,23	3.59	7 (38%)	26,32,35	0.98	1 (3%)
1	OMG	Aa	147	1	18,26,27	2.50	8 (44%)	19,38,41	1.51	4 (21%)
1	OMG	Aa	913	1	18,26,27	2.51	8 (44%)	19,38,41	1.49	4 (21%)
27	5MU	BA	1766	27	19,22,23	4.88	7 (36%)	28,32,35	3.67	9 (32%)
1	OMG	Aa	770	1	18,26,27	2.49	8 (44%)	19,38,41	1.55	4 (21%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
27	OMG	BA	584	27	18,26,27	2.50	8 (44%)	19,38,41	1.52	4 (21%)
27	OMG	BA	2010	27	18,26,27	2.46	8 (44%)	19,38,41	1.53	4 (21%)
27	OMC	BA	2047	27	19,22,23	2.96	8 (42%)	26,31,34	0.74	0
1	5MC	Aa	685	1	18,22,23	3.53	7 (38%)	26,32,35	1.04	2 (7%)
1	5MC	Aa	457	1	18,22,23	3.58	7 (38%)	26,32,35	1.01	2 (7%)
27	5MU	BA	2427	27	19,22,23	4.92	7 (36%)	28,32,35	3.62	9 (32%)
27	OMG	BA	365	27	18,26,27	2.46	8 (44%)	19,38,41	1.50	4 (21%)
27	OMG	BA	2550	27	18,26,27	1.08	1 (5%)	19,38,41	0.99	2 (10%)
1	OMG	Aa	670	1	18,26,27	2.51	8 (44%)	19,38,41	1.56	4 (21%)
27	OMG	BA	2745	27	18,26,27	2.45	8 (44%)	19,38,41	1.56	3 (15%)
1	MA6	Aa	1468	1	19,26,27	1.04	2 (10%)	18,38,41	3.72	2 (11%)
27	OMG	BA	670	27	18,26,27	2.47	8 (44%)	19,38,41	1.65	5 (26%)
27	OMG	BA	2168	27	18,26,27	2.46	8 (44%)	19,38,41	1.50	4 (21%)
27	5MU	BA	875	27	19,22,23	4.92	7 (36%)	28,32,35	3.63	10 (35%)
27	OMC	BA	2967	27	19,22,23	3.00	8 (42%)	26,31,34	0.73	0
27	OMG	BA	800	27	18,26,27	2.47	8 (44%)	19,38,41	1.52	4 (21%)
27	OMU	BA	1776	27	19,22,23	3.02	8 (42%)	26,31,34	1.65	4 (15%)
1	OMC	Aa	924	1	19,22,23	3.01	8 (42%)	26,31,34	0.73	0
27	OMG	BA	1953	27	18,26,27	2.45	8 (44%)	19,38,41	1.48	4 (21%)
27	OMG	BA	2642	27	18,26,27	2.48	8 (44%)	19,38,41	1.64	4 (21%)
1	OMG	Aa	659	1	18,26,27	2.48	8 (44%)	19,38,41	1.47	4 (21%)
27	OMG	BA	872	27	18,26,27	2.44	8 (44%)	19,38,41	1.47	3 (15%)
1	OMG	Aa	532	1	18,26,27	2.49	8 (44%)	19,38,41	1.51	4 (21%)
27	OMU	BA	453	27	19,22,23	3.03	8 (42%)	26,31,34	1.72	4 (15%)
27	5MC	BA	2055	27	18,22,23	3.57	7 (38%)	26,32,35	1.00	1 (3%)
27	OMC	BA	2169	27	19,22,23	2.98	8 (42%)	26,31,34	0.75	0
1	5MC	Aa	473	1	18,22,23	3.54	7 (38%)	26,32,35	1.07	2 (7%)
27	OMU	BA	1969	27	19,22,23	3.05	8 (42%)	26,31,34	1.73	5 (19%)
27	OMC	BA	66	27	19,22,23	2.97	8 (42%)	26,31,34	0.76	0
27	OMU	BA	2389	27	19,22,23	1.33	4 (21%)	26,31,34	2.01	7 (26%)
1	OMG	Aa	320	1	18,26,27	2.49	8 (44%)	19,38,41	1.54	4 (21%)
1	6MZ	Aa	1449	1	18,25,26	0.88	1 (5%)	16,36,39	1.93	3 (18%)
27	OMC	BA	91	27	19,22,23	2.99	8 (42%)	26,31,34	0.69	0
27	5MC	BA	1965	27	18,22,23	3.55	7 (38%)	26,32,35	1.04	2 (7%)
27	A2M	BA	505	27	18,25,26	4.22	7 (38%)	18,36,39	2.93	3 (16%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
27	OMG	BA	507	27	18,26,27	2.47	8 (44%)	19,38,41	1.53	3 (15%)
1	OMG	Aa	852	1	18,26,27	0.99	1 (5%)	19,38,41	1.06	2 (10%)
1	5MC	Aa	854	1	18,22,23	3.55	7 (38%)	26,32,35	1.05	2 (7%)
27	OMG	BA	1893	27	18,26,27	2.49	8 (44%)	19,38,41	1.56	4 (21%)
1	5MC	Aa	464	1	18,22,23	3.59	7 (38%)	26,32,35	1.03	2 (7%)
1	OMG	Aa	574	1	18,26,27	2.52	8 (44%)	19,38,41	1.50	4 (21%)
27	OMG	BA	63	27	18,26,27	2.48	8 (44%)	19,38,41	1.57	4 (21%)
1	LV2	Aa	918	1	20,23,24	0.89	0	26,33,36	0.81	0
1	OMC	Aa	752	1	19,22,23	3.01	8 (42%)	26,31,34	0.71	0
1	OMC	Aa	112	1	19,22,23	2.99	8 (42%)	26,31,34	0.74	0
27	5MC	BA	1345	27	18,22,23	3.55	7 (38%)	26,32,35	1.02	1 (3%)
1	5MC	Aa	1184	1	18,22,23	3.56	7 (38%)	26,32,35	1.01	2 (7%)
27	5MC	BA	2070	27	18,22,23	3.50	7 (38%)	26,32,35	1.01	2 (7%)
27	OMG	BA	2466	27	18,26,27	2.46	8 (44%)	19,38,41	1.47	3 (15%)
1	OMG	Aa	836	1	18,26,27	2.51	8 (44%)	19,38,41	1.55	4 (21%)
27	OMG	BA	2870	27	18,26,27	2.45	8 (44%)	19,38,41	1.48	4 (21%)
27	OMU	BA	657	27	19,22,23	3.01	8 (42%)	26,31,34	1.70	4 (15%)
27	5MC	BA	55	27	18,22,23	3.48	7 (38%)	26,32,35	1.17	3 (11%)
1	OMU	Aa	479	1	19,22,23	3.07	8 (42%)	26,31,34	1.70	4 (15%)
1	OMC	Aa	1361	1	19,22,23	3.02	8 (42%)	26,31,34	0.74	0
1	OMG	Aa	891	1	18,26,27	2.48	8 (44%)	19,38,41	1.51	4 (21%)
27	OMG	BA	2469	27	18,26,27	2.46	8 (44%)	19,38,41	1.47	4 (21%)
27	5MU	BA	829	27	19,22,23	4.86	7 (36%)	28,32,35	3.68	10 (35%)
1	A2M	Aa	352	1	18,25,26	4.20	7 (38%)	18,36,39	3.03	3 (16%)
1	OMG	Aa	1223	1	18,26,27	2.52	8 (44%)	19,38,41	1.51	4 (21%)
27	OMG	BA	215	27	18,26,27	2.47	8 (44%)	19,38,41	1.55	4 (21%)
1	5MC	Aa	349	1	18,22,23	3.54	7 (38%)	26,32,35	1.04	2 (7%)
27	OMG	BA	2379	27	18,26,27	2.48	8 (44%)	19,38,41	1.53	4 (21%)
1	OMG	Aa	629	1	18,26,27	2.51	8 (44%)	19,38,41	1.54	4 (21%)
1	OMG	Aa	381	1	18,26,27	2.52	8 (44%)	19,38,41	1.53	4 (21%)
27	OMC	BA	1995	27	19,22,23	3.00	8 (42%)	26,31,34	0.74	0
1	5MC	Aa	340	1	18,22,23	3.56	7 (38%)	26,32,35	1.11	2 (7%)
27	OMU	BA	1402	27	19,22,23	3.06	8 (42%)	26,31,34	1.67	4 (15%)
27	OMG	BA	2016	27	18,26,27	2.47	8 (44%)	19,38,41	1.48	4 (21%)

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
27	5MC	BA	2688	27	-	0/7/25/26	0/2/2/2
27	A2M	BA	2329	27	-	2/5/27/28	0/3/3/3
1	OMU	Aa	8	1	-	6/9/27/28	0/2/2/2
1	OMU	Aa	753	1	-	1/9/27/28	0/2/2/2
1	OMG	Aa	219	1	-	2/5/27/28	0/3/3/3
1	2MG	Aa	995	1	-	0/5/27/28	0/3/3/3
27	OMU	BA	1150	27	-	2/9/27/28	0/2/2/2
27	OMG	BA	2728	27	-	1/5/27/28	0/3/3/3
1	OMC	Aa	825	1	-	0/9/27/28	0/2/2/2
1	OMG	Aa	636	1	-	1/5/27/28	0/3/3/3
1	5MC	Aa	1321	1	-	0/7/25/26	0/2/2/2
1	OMG	Aa	121	1	-	2/5/27/28	0/3/3/3
27	5MC	BA	2075	27	-	0/7/25/26	0/2/2/2
27	OMC	BA	2595	27	-	1/9/27/28	0/2/2/2
1	5MC	Aa	17	1	-	2/7/25/26	0/2/2/2
27	A2M	BA	879	27	-	1/5/27/28	0/3/3/3
27	OMG	BA	2353	27	-	2/5/27/28	0/3/3/3
27	OMG	BA	1956	27	-	2/5/27/28	0/3/3/3
1	OMG	Aa	994	1	-	0/5/27/28	0/3/3/3
1	OMG	Aa	495	1	-	2/5/27/28	0/3/3/3
1	MA6	Aa	1467	1	-	0/7/29/30	0/3/3/3
1	OMG	Aa	963	1	-	2/5/27/28	0/3/3/3
1	OMG	Aa	1107	1	-	2/5/27/28	0/3/3/3
1	5MC	Aa	1004	1	-	3/7/25/26	0/2/2/2
1	5MC	Aa	1007	1	-	2/7/25/26	0/2/2/2
27	OMG	BA	1508	27	-	0/5/27/28	0/3/3/3
27	OMC	BA	2107	27	-	0/9/27/28	0/2/2/2
1	5MC	Aa	1306	1	-	3/7/25/26	0/2/2/2
27	OMU	BA	1480	27	-	0/9/27/28	0/2/2/2
27	OMG	BA	2521	27	-	1/5/27/28	0/3/3/3
27	OMG	BA	2528	27	-	0/5/27/28	0/3/3/3
27	OMG	BA	833	27	-	0/5/27/28	0/3/3/3
27	OMC	BA	869	27	-	0/9/27/28	0/2/2/2
1	OMG	Aa	450	1	-	0/5/27/28	0/3/3/3
1	OMC	Aa	1354	1	-	0/9/27/28	0/2/2/2
1	OMG	Aa	446	1	-	0/5/27/28	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
27	A2M	BA	932	27	-	0/5/27/28	0/3/3/3
27	OMU	BA	2656	27	-	0/9/27/28	0/2/2/2
1	5MC	Aa	923	1	-	0/7/25/26	0/2/2/2
27	OMG	BA	912	27	-	1/5/27/28	0/3/3/3
27	OMU	BA	2542	27	-	2/9/27/28	0/2/2/2
27	OMC	BA	342	27	-	2/9/27/28	0/2/2/2
1	OMG	Aa	498	1	-	2/5/27/28	0/3/3/3
1	5MC	Aa	1006	1	-	0/7/25/26	0/2/2/2
27	OMG	BA	2527	27	-	2/5/27/28	0/3/3/3
27	4SU	BA	2581	27	-	1/7/25/26	0/2/2/2
27	OMG	BA	1525	27	-	2/5/27/28	0/3/3/3
27	A2M	BA	849	27	-	2/5/27/28	0/3/3/3
27	OMG	BA	2672	27	-	0/5/27/28	0/3/3/3
1	5MU	Aa	917	1	-	2/7/25/26	0/2/2/2
1	5MC	Aa	1266	1	-	0/7/25/26	0/2/2/2
27	OMC	BA	2545	27	-	3/9/27/28	0/2/2/2
1	5MC	Aa	672	1	-	0/7/25/26	0/2/2/2
27	OMC	BA	1820	27	-	1/9/27/28	0/2/2/2
1	5MC	Aa	1003	1	-	0/7/25/26	0/2/2/2
1	OMC	Aa	229	1	-	1/9/27/28	0/2/2/2
27	4SU	BA	2553	27	-	0/7/25/26	0/2/2/2
27	OMG	BA	674	27	-	2/5/27/28	0/3/3/3
27	5MU	BA	1854	27	-	0/7/25/26	0/2/2/2
1	5MC	Aa	1478	1	-	5/7/25/26	0/2/2/2
27	OMG	BA	288	27	-	1/5/27/28	0/3/3/3
1	5MC	Aa	1476	1	-	0/7/25/26	0/2/2/2
27	5MC	BA	2605	27	-	4/7/25/26	0/2/2/2
27	OMC	BA	2124	27	-	1/9/27/28	0/2/2/2
27	A2M	BA	2091	27	-	0/5/27/28	0/3/3/3
27	OMG	BA	65	27	-	1/5/27/28	0/3/3/3
27	OMG	BA	2640	27	-	3/5/27/28	0/3/3/3
1	OMU	Aa	1020	1	-	1/9/27/28	0/2/2/2
1	OMU	Aa	1358	1	-	0/9/27/28	0/2/2/2
27	OMG	BA	1329	27	-	0/5/27/28	0/3/3/3
27	A2M	BA	2936	27	-	1/5/27/28	0/3/3/3
1	5MC	Aa	1352	1	-	2/7/25/26	0/2/2/2
1	OMG	Aa	147	1	-	1/5/27/28	0/3/3/3
1	OMG	Aa	913	1	-	2/5/27/28	0/3/3/3
27	5MU	BA	1766	27	-	0/7/25/26	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	OMG	Aa	770	1	-	0/5/27/28	0/3/3/3
27	OMG	BA	584	27	-	0/5/27/28	0/3/3/3
27	OMG	BA	2010	27	-	0/5/27/28	0/3/3/3
27	OMC	BA	2047	27	-	3/9/27/28	0/2/2/2
1	5MC	Aa	685	1	-	0/7/25/26	0/2/2/2
1	5MC	Aa	457	1	-	2/7/25/26	0/2/2/2
27	5MU	BA	2427	27	-	0/7/25/26	0/2/2/2
27	OMG	BA	365	27	-	3/5/27/28	0/3/3/3
27	OMG	BA	2550	27	-	0/5/27/28	0/3/3/3
1	OMG	Aa	670	1	-	2/5/27/28	0/3/3/3
27	OMG	BA	2745	27	-	0/5/27/28	0/3/3/3
1	MA6	Aa	1468	1	-	1/7/29/30	0/3/3/3
27	OMG	BA	670	27	-	1/5/27/28	0/3/3/3
27	OMG	BA	2168	27	-	0/5/27/28	0/3/3/3
27	5MU	BA	875	27	-	0/7/25/26	0/2/2/2
27	OMC	BA	2967	27	-	0/9/27/28	0/2/2/2
27	OMG	BA	800	27	-	1/5/27/28	0/3/3/3
27	OMU	BA	1776	27	-	1/9/27/28	0/2/2/2
1	OMC	Aa	924	1	-	1/9/27/28	0/2/2/2
27	OMG	BA	1953	27	-	1/5/27/28	0/3/3/3
27	OMG	BA	2642	27	-	3/5/27/28	0/3/3/3
1	OMG	Aa	659	1	-	1/5/27/28	0/3/3/3
27	OMG	BA	872	27	-	0/5/27/28	0/3/3/3
1	OMG	Aa	532	1	-	0/5/27/28	0/3/3/3
27	OMU	BA	453	27	-	2/9/27/28	0/2/2/2
27	5MC	BA	2055	27	-	0/7/25/26	0/2/2/2
27	OMC	BA	2169	27	-	0/9/27/28	0/2/2/2
1	5MC	Aa	473	1	-	0/7/25/26	0/2/2/2
27	OMU	BA	1969	27	-	0/9/27/28	0/2/2/2
27	OMC	BA	66	27	-	1/9/27/28	0/2/2/2
27	OMU	BA	2389	27	-	2/9/27/28	0/2/2/2
1	OMG	Aa	320	1	-	3/5/27/28	0/3/3/3
1	6MZ	Aa	1449	1	-	2/5/27/28	0/3/3/3
27	OMC	BA	91	27	-	1/9/27/28	0/2/2/2
27	5MC	BA	1965	27	-	0/7/25/26	0/2/2/2
27	A2M	BA	505	27	-	0/5/27/28	0/3/3/3
27	OMG	BA	507	27	-	0/5/27/28	0/3/3/3
1	OMG	Aa	852	1	-	0/5/27/28	0/3/3/3
1	5MC	Aa	854	1	-	0/7/25/26	0/2/2/2
27	OMG	BA	1893	27	-	2/5/27/28	0/3/3/3

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
1	5MC	Aa	464	1	-	0/7/25/26	0/2/2/2
1	OMG	Aa	574	1	-	1/5/27/28	0/3/3/3
27	OMG	BA	63	27	-	0/5/27/28	0/3/3/3
1	LV2	Aa	918	1	-	0/9/29/30	0/2/2/2
1	OMC	Aa	752	1	-	0/9/27/28	0/2/2/2
1	OMC	Aa	112	1	-	0/9/27/28	0/2/2/2
27	5MC	BA	1345	27	-	1/7/25/26	0/2/2/2
1	5MC	Aa	1184	1	-	0/7/25/26	0/2/2/2
27	5MC	BA	2070	27	-	0/7/25/26	0/2/2/2
27	OMG	BA	2466	27	-	0/5/27/28	0/3/3/3
1	OMG	Aa	836	1	-	0/5/27/28	0/3/3/3
27	OMG	BA	2870	27	-	0/5/27/28	0/3/3/3
27	OMU	BA	657	27	-	1/9/27/28	0/2/2/2
27	5MC	BA	55	27	-	3/7/25/26	0/2/2/2
1	OMU	Aa	479	1	-	0/9/27/28	0/2/2/2
1	OMC	Aa	1361	1	-	1/9/27/28	0/2/2/2
1	OMG	Aa	891	1	-	1/5/27/28	0/3/3/3
27	OMG	BA	2469	27	-	4/5/27/28	0/3/3/3
27	5MU	BA	829	27	-	0/7/25/26	0/2/2/2
1	A2M	Aa	352	1	-	2/5/27/28	0/3/3/3
1	OMG	Aa	1223	1	-	0/5/27/28	0/3/3/3
27	OMG	BA	215	27	-	0/5/27/28	0/3/3/3
1	5MC	Aa	349	1	-	0/7/25/26	0/2/2/2
27	OMG	BA	2379	27	-	2/5/27/28	0/3/3/3
1	OMG	Aa	629	1	-	0/5/27/28	0/3/3/3
1	OMG	Aa	381	1	-	2/5/27/28	0/3/3/3
27	OMC	BA	1995	27	-	1/9/27/28	0/2/2/2
1	5MC	Aa	340	1	-	3/7/25/26	0/2/2/2
27	OMU	BA	1402	27	-	2/9/27/28	0/2/2/2
27	OMG	BA	2016	27	-	2/5/27/28	0/3/3/3

The worst 5 of 1024 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
27	BA	932	A2M	O4'-C1'	15.41	1.62	1.41
27	BA	879	A2M	O4'-C1'	15.29	1.62	1.41
27	BA	2936	A2M	O4'-C1'	15.28	1.62	1.41
27	BA	505	A2M	O4'-C1'	15.18	1.62	1.41
1	Aa	352	A2M	O4'-C1'	15.09	1.62	1.41

The worst 5 of 446 bond angle outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	Aa	1468	MA6	N1-C6-N6	-14.53	101.77	117.06
1	Aa	1467	MA6	N1-C6-N6	-14.09	102.23	117.06
27	BA	1854	5MU	C5-C4-N3	12.41	125.90	115.31
27	BA	1766	5MU	C5-C4-N3	12.21	125.73	115.31
27	BA	829	5MU	C5-C4-N3	12.09	125.63	115.31

There are no chirality outliers.

5 of 144 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	Aa	147	OMG	C1'-C2'-O2'-CM2
1	Aa	320	OMG	C3'-C4'-C5'-O5'
1	Aa	352	A2M	C1'-C2'-O2'-CM'
1	Aa	381	OMG	O4'-C4'-C5'-O5'
1	Aa	381	OMG	C3'-C4'-C5'-O5'

There are no ring outliers.

43 monomers are involved in 65 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
27	BA	2688	5MC	1	0
27	BA	1150	OMU	1	0
27	BA	2728	OMG	1	0
27	BA	2075	5MC	1	0
27	BA	2595	OMC	2	0
27	BA	879	A2M	2	0
27	BA	1956	OMG	2	0
27	BA	2107	OMC	1	0
27	BA	2521	OMG	1	0
27	BA	2528	OMG	1	0
27	BA	833	OMG	1	0
27	BA	932	A2M	1	0
27	BA	912	OMG	2	0
27	BA	2527	OMG	1	0
27	BA	2581	4SU	1	0
27	BA	849	A2M	1	0
27	BA	2545	OMC	2	0
27	BA	1820	OMC	1	0
27	BA	288	OMG	1	0
27	BA	2605	5MC	2	0
27	BA	2124	OMC	3	0
27	BA	65	OMG	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
27	BA	2640	OMG	2	0
27	BA	1329	OMG	1	0
27	BA	2936	A2M	1	0
27	BA	2427	5MU	3	0
27	BA	365	OMG	1	0
27	BA	2550	OMG	2	0
27	BA	2745	OMG	1	0
27	BA	670	OMG	5	0
27	BA	1776	OMU	2	0
27	BA	2642	OMG	5	0
27	BA	872	OMG	1	0
27	BA	1969	OMU	1	0
27	BA	91	OMC	1	0
27	BA	507	OMG	2	0
27	BA	1893	OMG	1	0
27	BA	2070	5MC	1	0
27	BA	657	OMU	2	0
27	BA	55	5MC	3	0
27	BA	2469	OMG	1	0
27	BA	2379	OMG	1	0
27	BA	1995	OMC	1	0

## 5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

Of 14 ligands modelled in this entry, 14 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](#)

There are no chain breaks in this entry.

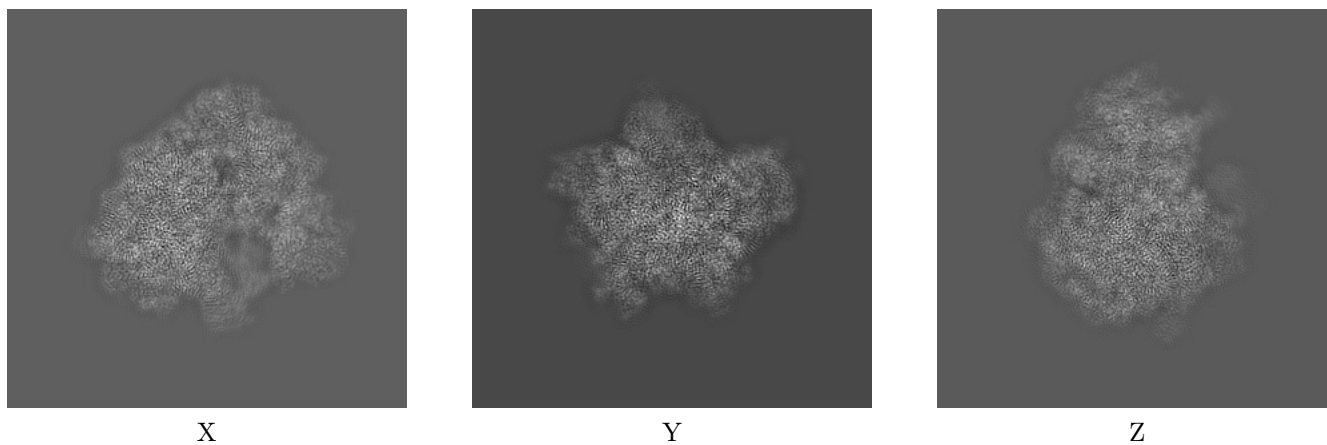
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-10224. These allow visual inspection of the internal detail of the map and identification of artifacts.

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

### 6.1 Orthogonal projections [i](#)

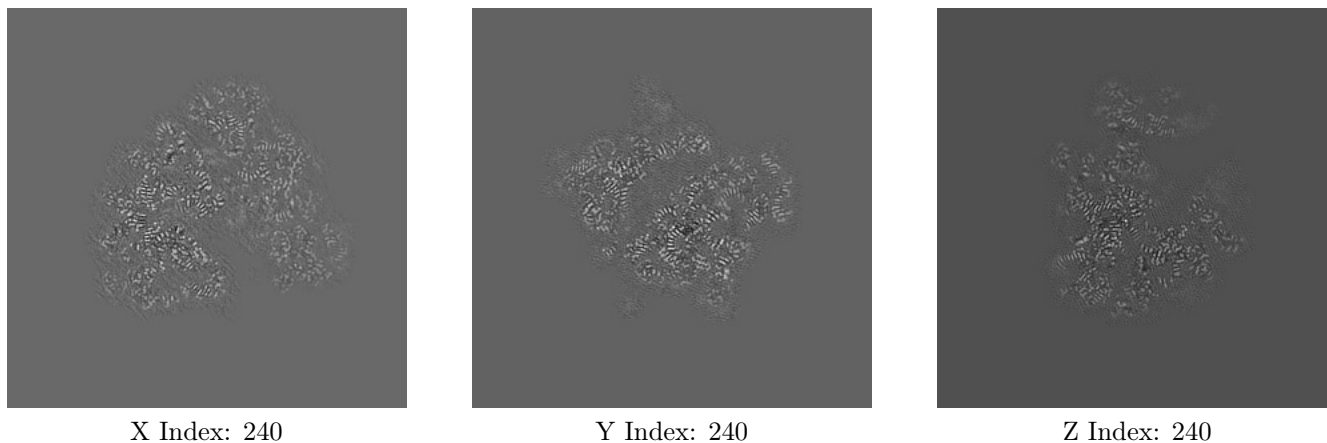
#### 6.1.1 Primary map



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

### 6.2 Central slices [i](#)

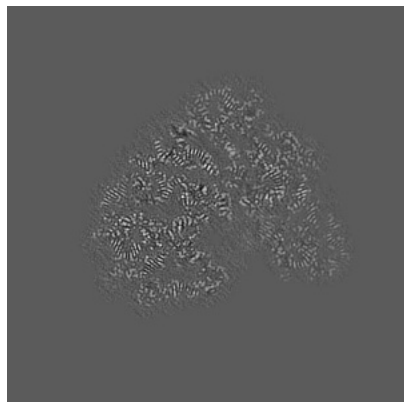
#### 6.2.1 Primary map



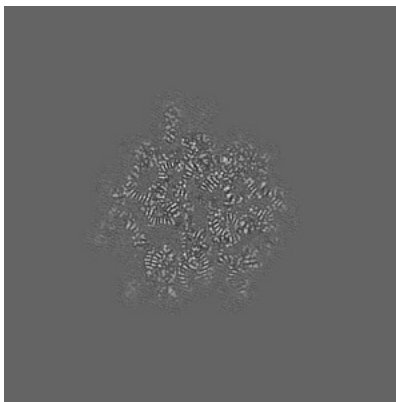
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

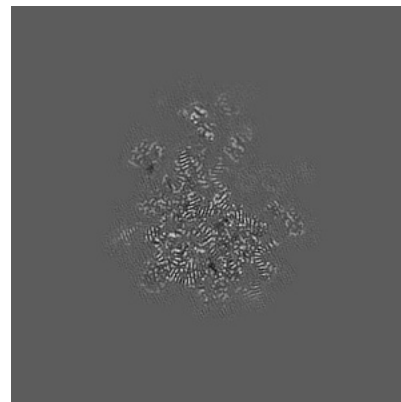
### 6.3.1 Primary map



X Index: 231



Y Index: 182

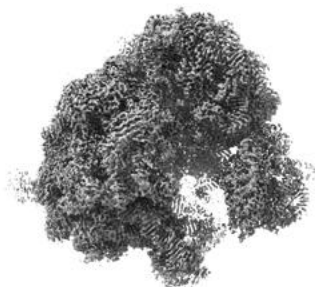


Z Index: 264

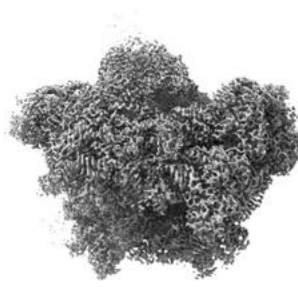
The images above show the largest variance slices of the map in three orthogonal directions.

## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.05. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

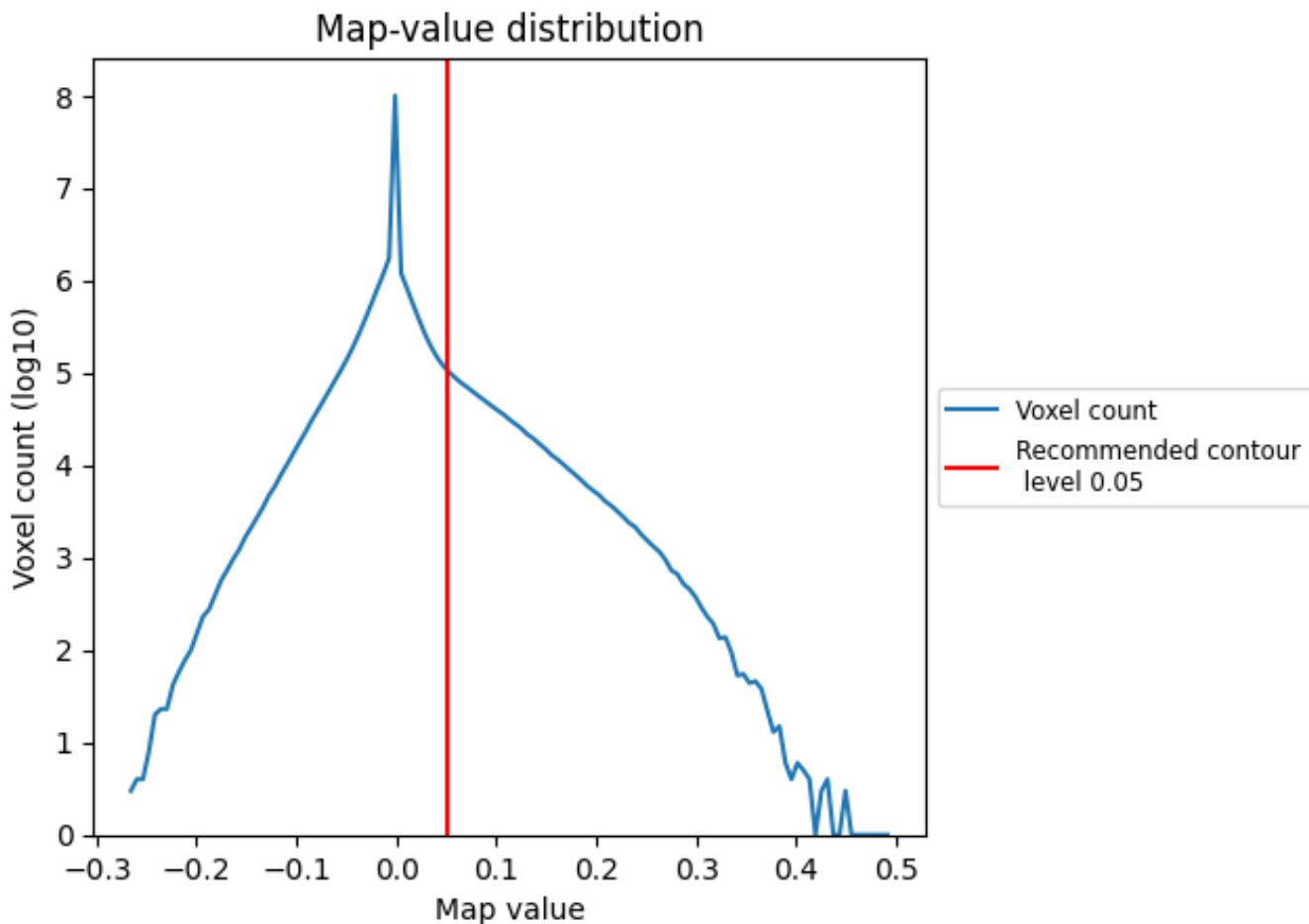
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

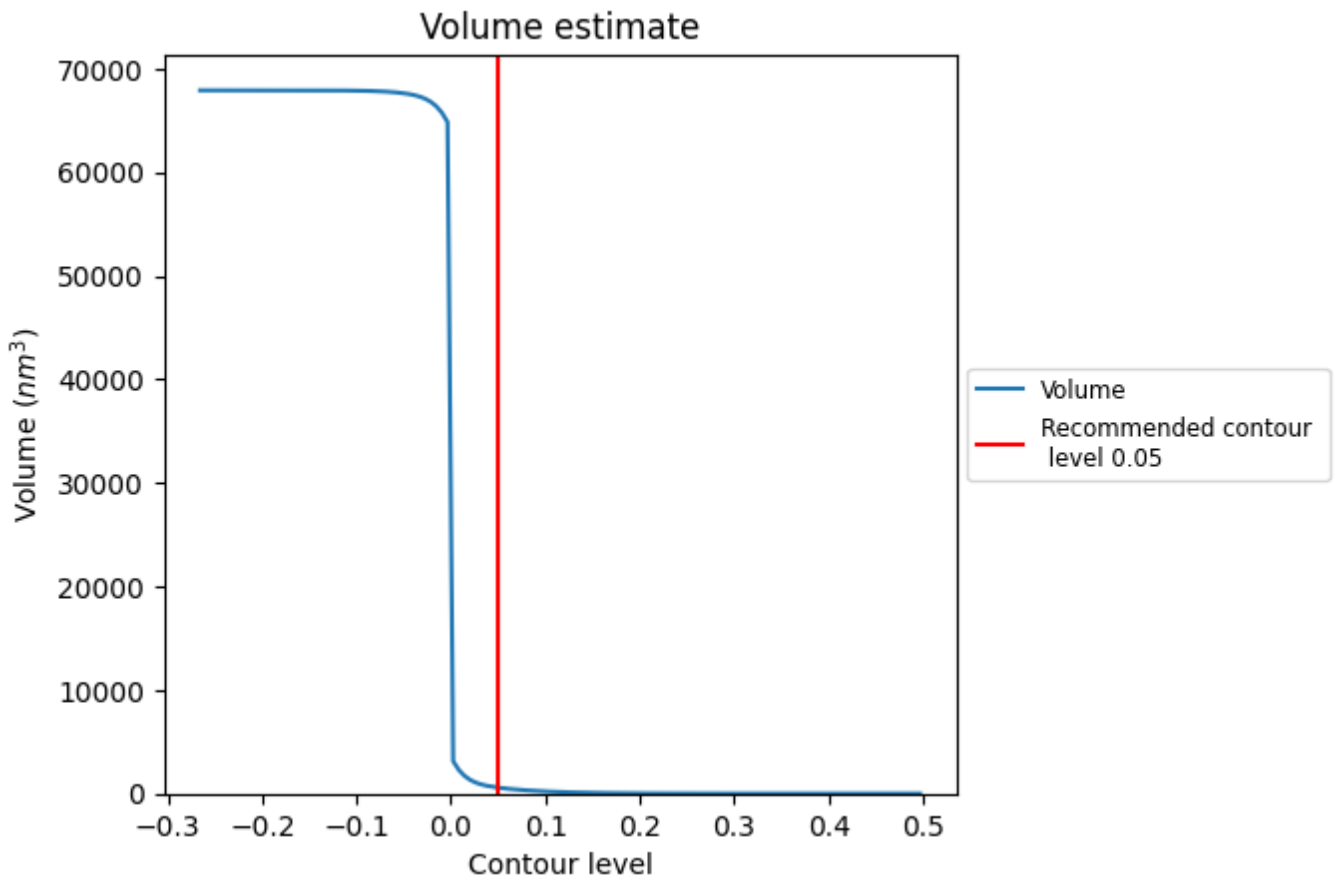
This section contains the results of statistical analysis of the map.

### 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 [\(i\)](#)

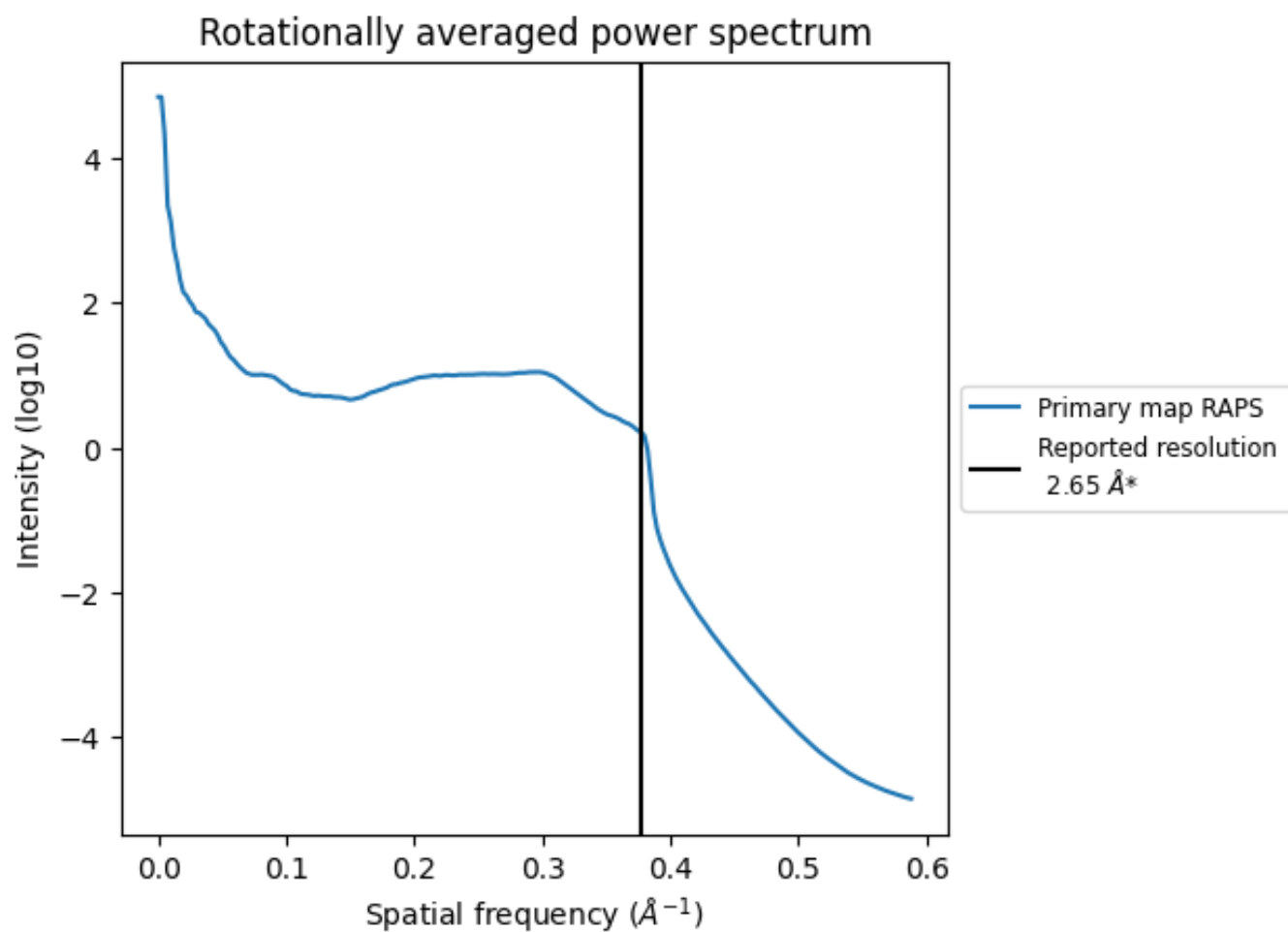


The volume at the recommended contour level is 576 nm<sup>3</sup>; this corresponds to an approximate mass of 520 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.



### 7.3 Rotationally averaged power spectrum [i](#)



\*Reported resolution corresponds to spatial frequency of  $0.377 \text{\AA}^{-1}$

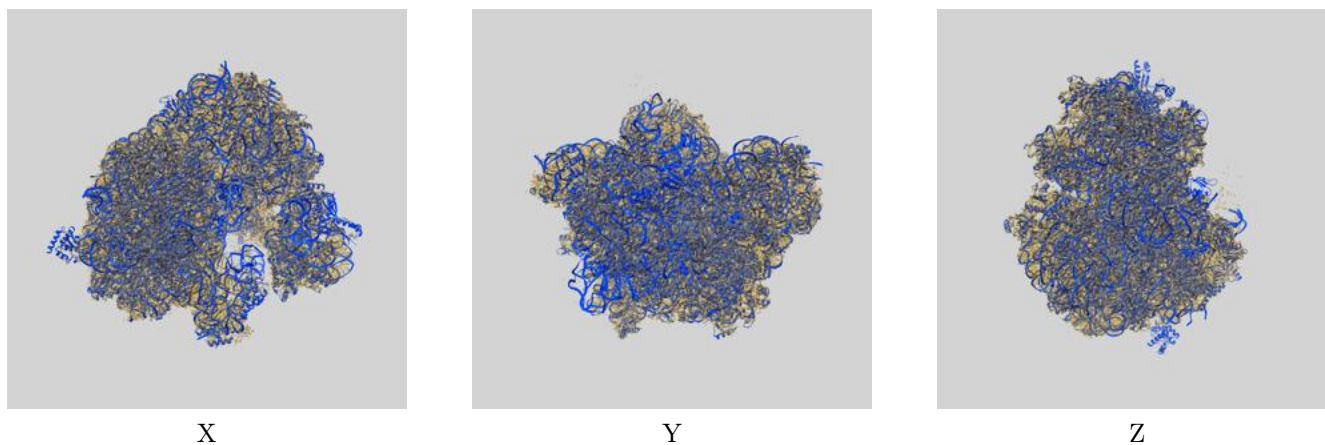
## 8 Fourier-Shell correlation

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

## 9 Map-model fit [i](#)

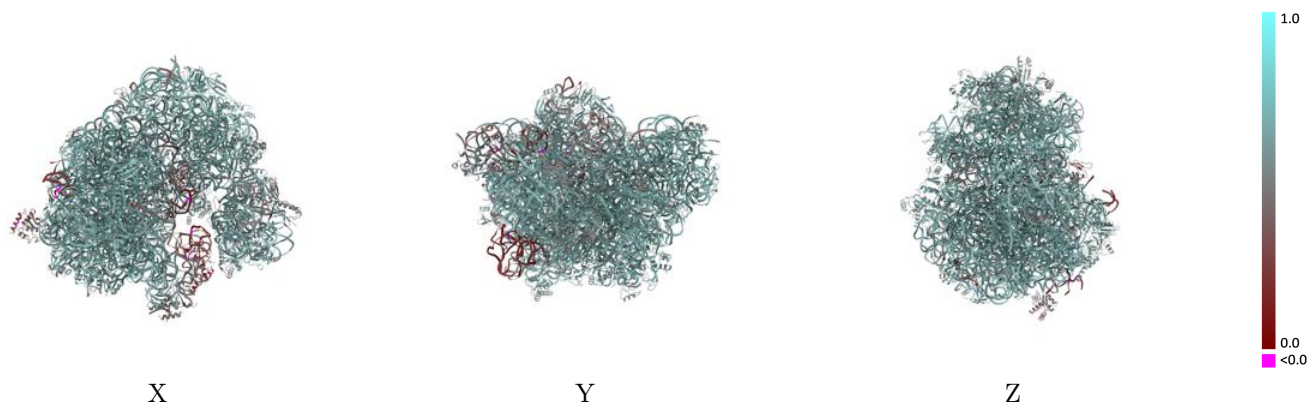
This section contains information regarding the fit between EMDB map EMD-10224 and PDB model 6SKG. Per-residue inclusion information can be found in section 3 on page 19.

### 9.1 Map-model overlay [i](#)



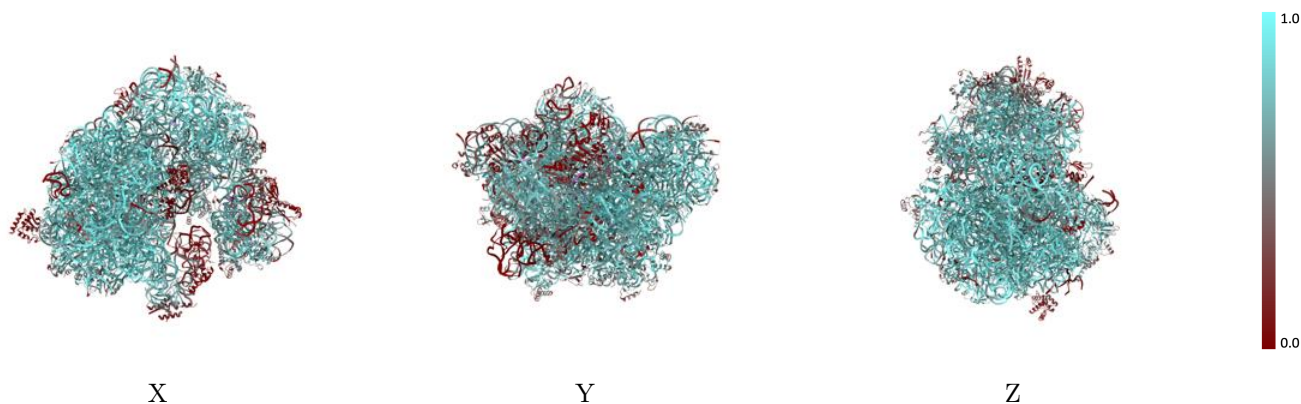
The images above show the 3D surface view of the map at the recommended contour level 0.05 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



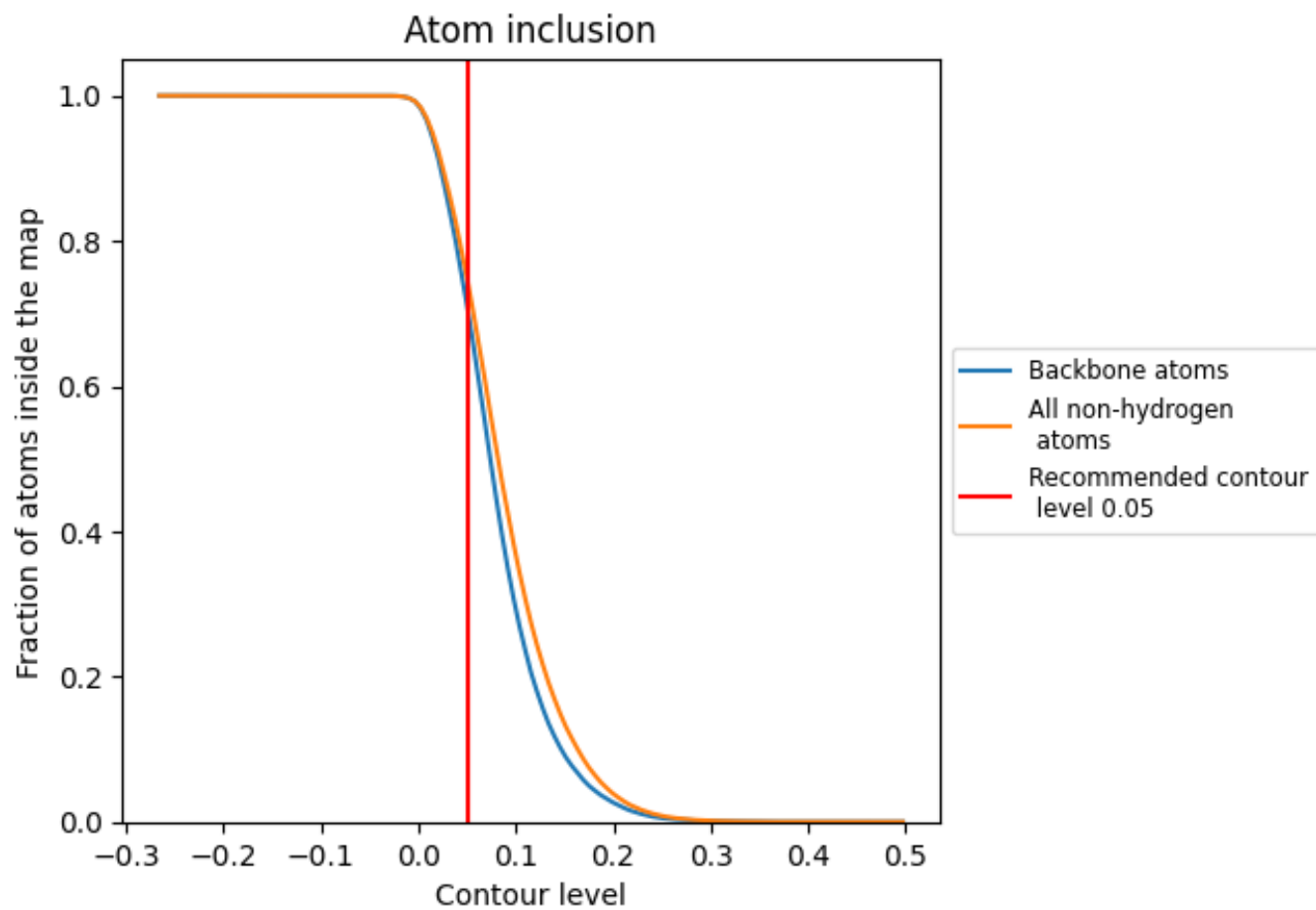
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.05).







































































## 9.4 Atom inclusion [i](#)



At the recommended contour level, 71% of all backbone atoms, 75% 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 (0.05) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7490	 0.6140
Aa	 0.7927	 0.6280
Ab	 0.5924	 0.6100
Ac	 0.2463	 0.5360
Ad	 0.5664	 0.5980
Ae	 0.6810	 0.6190
Af	 0.7274	 0.6350
Ag	 0.7098	 0.6200
Ah	 0.4000	 0.5690
Ai	 0.4940	 0.5990
Aj	 0.7553	 0.6300
Ak	 0.7110	 0.6310
Al	 0.5270	 0.5990
Am	 0.2703	 0.5380
An	 0.6043	 0.6020
Ao	 0.6923	 0.6310
Ap	 0.5219	 0.5790
Aq	 0.6911	 0.6190
Ar	 0.4755	 0.5880
As	 0.7243	 0.6360
At	 0.3019	 0.5440
Au	 0.5405	 0.6000
Av	 0.5909	 0.6110
Aw	 0.6085	 0.5970
Ax	 0.5551	 0.5970
Ay	 0.3013	 0.5600
Az	 0.7255	 0.6200
BA	 0.8478	 0.6220
BB	 0.6378	 0.5600
BC	 0.8460	 0.6450
BD	 0.8391	 0.6530
BE	 0.8164	 0.6360
BF	 0.1501	 0.3600
BG	 0.6964	 0.6090
BH	 0.5727	 0.5750



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Chain	Atom inclusion	Q-score
BI	 0.0568	 0.3850
BJ	 0.7458	 0.6240
BK	 0.8161	 0.6490
BL	 0.7818	 0.6350
BM	 0.5859	 0.5680
BN	 0.6077	 0.5780
BO	 0.6684	 0.6110
BP	 0.8953	 0.6670
BQ	 0.5273	 0.5510
BR	 0.8238	 0.6440
BS	 0.8035	 0.6340
BT	 0.6926	 0.6170
BU	 0.8368	 0.6540
BV	 0.8375	 0.6560
BW	 0.7216	 0.6220
BX	 0.7981	 0.6320
BY	 0.7937	 0.6420
BZ	 0.6173	 0.5890
Ba	 0.8071	 0.6440
Bb	 0.6504	 0.5960
Bc	 0.7730	 0.6340
Bd	 0.8350	 0.6530
Be	 0.8228	 0.6360
Bf	 0.7601	 0.6300
Bg	 0.7888	 0.6300
Bh	 0.9502	 0.6770
Bi	 0.8495	 0.6450
Bj	 0.7640	 0.6420
Bk	 0.8019	 0.6360
Bl	 0.7513	 0.6440
Bm	 0.5271	 0.5600
Bn	 0.5294	 0.6000
Bo	 0.8142	 0.6430
Bp	 0.0126	 0.3690
Bq	 0.0235	 0.4650