



wwPDB EM Validation Summary Report ⓘ

Dec 31, 2024 – 12:33 AM EST

PDB ID : 8FLD
EMDB ID : EMD-29275
Title : Human nuclear pre-60S ribosomal subunit (State L1)
Authors : Vanden Broeck, A.; Klinge, S.
Deposited on : 2022-12-21
Resolution : 2.58 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.dev113
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4.02b-467
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.40

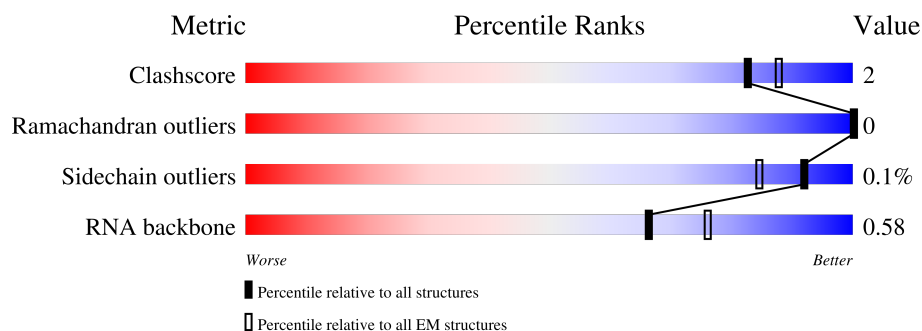
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.58 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



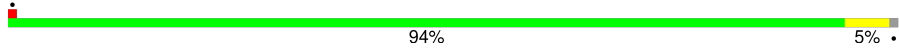

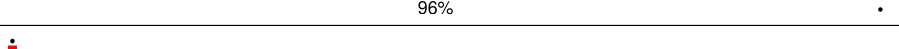
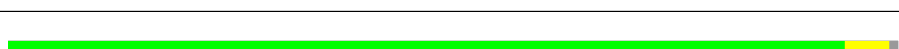


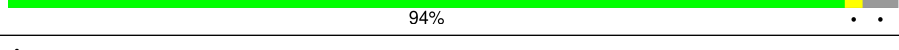
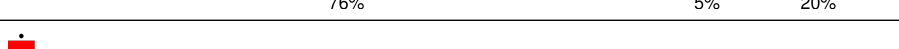



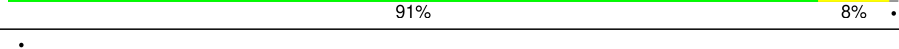

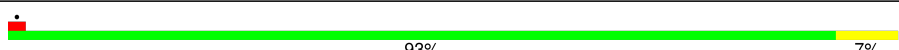


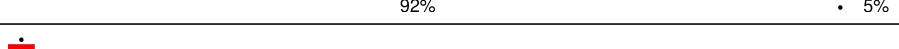



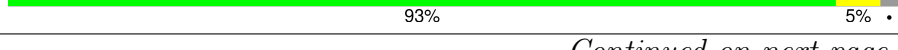



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415
RNA backbone	6643	2191

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$. The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion $< 40\%$). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	BE	214	
2	L1	157	
3	L2	1167	
4	L3	5070	
5	L4	121	
6	L5	178	
7	L6	211	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
8	L7	203	
9	L8	215	
10	L9	204	
11	LA	184	
12	LB	188	
13	LC	176	
14	LD	196	
15	LE	160	
16	LF	128	
17	LG	140	
18	LH	156	
19	LI	145	
20	LJ	136	
21	LK	148	
22	LL	137	
23	LM	159	
24	LN	403	
25	LO	115	
26	LP	125	
27	LQ	135	
28	LR	117	
29	LS	123	
30	LT	110	
31	LU	105	
32	LV	106	

Continued on next page...

Continued from previous page...

Mol	Chain	Length	Quality of chain
33	LW	97	
34	LX	92	
35	LY	70	
36	LZ	51	
37	NK	129	
38	NL	478	
39	NP	134	
40	SA	427	
41	SB	297	
42	SC	288	
43	SD	248	
44	SE	266	
45	SF	257	
46	SG	192	
47	SH	293	
48	SI	255	
49	SK	245	
50	SL	490	
51	SM	588	
52	SQ	239	
53	SR	634	
54	SV	163	

2 Entry composition

There are 56 unique types of molecules in this entry. The entry contains 148569 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 protein called 60S ribosomal protein L10.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	BE	160	Total	C	N	O	S	0	0
			1295	827	242	218	8		

- Molecule 2 is a RNA chain called 5.8S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	L1	154	Total	C	N	O	P	0	0
			3278	1463	581	1080	154		

- Molecule 3 is a RNA chain called ITS2 rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	L2	72	Total	C	N	O	P	0	0
			1535	683	278	502	72		

- Molecule 4 is a RNA chain called 28S rRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	L3	3372	Total	C	N	O	P	0	0
			72381	32259	13258	23492	3372		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
5	L4	120	Total	C	N	O	P	0	0
			2561	1141	456	844	120		

- Molecule 6 is a protein called 60S ribosomal protein L11.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	L5	168	Total	C	N	O	S	0	0
			1349	853	251	239	6		

- Molecule 7 is a protein called 60S ribosomal protein L13.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	L6	210	Total	C	N	O	S	0	0
			1701	1064	352	281	4		

- Molecule 8 is a protein called 60S ribosomal protein L13a.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	L7	201	Total	C	N	O	S	0	0
			1650	1063	321	261	5		

- Molecule 9 is a protein called 60S ribosomal protein L14.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	L8	135	Total	C	N	O	S	0	0
			1111	713	213	178	7		

- Molecule 10 is a protein called 60S ribosomal protein L15.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	L9	203	Total	C	N	O	S	0	0
			1701	1072	359	266	4		

- Molecule 11 is a protein called 60S ribosomal protein L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	LA	153	Total	C	N	O	S	0	0
			1242	776	241	216	9		

- Molecule 12 is a protein called 60S ribosomal protein L18.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	LB	187	Total	C	N	O	S	0	0
			1512	944	314	249	5		

- Molecule 13 is a protein called 60S ribosomal protein L18a.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	LC	176	Total	C	N	O	S	0	0
			1461	930	284	236	11		

- Molecule 14 is a protein called 60S ribosomal protein L19.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	LD	154	Total	C	N	O	S	0	0
			1289	805	277	198	9		

- Molecule 15 is a protein called 60S ribosomal protein L21.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	LE	154	Total	C	N	O	S	0	0
			1264	803	246	210	5		

- Molecule 16 is a protein called 60S ribosomal protein L22.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	LF	103	Total	C	N	O	S	0	0
			842	538	148	154	2		

- Molecule 17 is a protein called 60S ribosomal protein L23.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	LG	139	Total	C	N	O	S	0	0
			1034	648	199	182	5		

- Molecule 18 is a protein called 60S ribosomal protein L23a.

Mol	Chain	Residues	Atoms					AltConf	Trace
18	LH	143	Total	C	N	O	S	0	0
			1156	740	220	195	1		

- Molecule 19 is a protein called 60S ribosomal protein L26.

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

- Molecule 20 is a protein called 60S ribosomal protein L27.

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

- Molecule 21 is a protein called 60S ribosomal protein L27a.

Mol	Chain	Residues	Atoms					AltConf	Trace
21	LK	147	Total	C	N	O	S	0	0
			1162	736	237	186	3		

- Molecule 22 is a protein called 60S ribosomal protein L28.

Mol	Chain	Residues	Atoms					AltConf	Trace
22	LL	125	Total	C	N	O	S	0	0
			1002	622	207	168	5		

- Molecule 23 is a protein called 60S ribosomal protein L29.

Mol	Chain	Residues	Atoms					AltConf	Trace
23	LM	94	Total	C	N	O	S	0	0
			775	483	170	118	4		

- Molecule 24 is a protein called 60S ribosomal protein L3.

Mol	Chain	Residues	Atoms					AltConf	Trace
24	LN	402	Total	C	N	O	S	0	0
			3239	2061	608	556	14		

- Molecule 25 is a protein called 60S ribosomal protein L30.

Mol	Chain	Residues	Atoms					AltConf	Trace
25	LO	95	Total	C	N	O	S	0	0
			738	468	131	133	6		

- Molecule 26 is a protein called 60S ribosomal protein L31.

Mol	Chain	Residues	Atoms					AltConf	Trace
26	LP	106	Total	C	N	O	S	0	0
			879	555	170	152	2		

- Molecule 27 is a protein called 60S ribosomal protein L32.

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

- Molecule 28 is a protein called 60S ribosomal protein L34.

Mol	Chain	Residues	Atoms					AltConf	Trace
28	LR	112	Total	C	N	O	S	0	0
			888	555	183	144	6		

- Molecule 29 is a protein called 60S ribosomal protein L35.

Mol	Chain	Residues	Atoms					AltConf	Trace
29	LS	122	Total	C	N	O	S	0	0
			1015	641	205	168	1		

- Molecule 30 is a protein called 60S ribosomal protein L35a.

Mol	Chain	Residues	Atoms					AltConf	Trace
30	LT	109	Total	C	N	O	S	0	0
			876	555	174	144	3		

- Molecule 31 is a protein called 60S ribosomal protein L36.

Mol	Chain	Residues	Atoms					AltConf	Trace
31	LU	102	Total	C	N	O	S	0	0
			832	521	177	129	5		

- Molecule 32 is a protein called 60S ribosomal protein L36a.

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

- Molecule 33 is a protein called 60S ribosomal protein L37.

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

- Molecule 34 is a protein called 60S ribosomal protein L37a.

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

- Molecule 35 is a protein called 60S ribosomal protein L38.

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

- Molecule 36 is a protein called 60S ribosomal protein L39.

Mol	Chain	Residues	Atoms					AltConf	Trace
36	LZ	50	Total	C	N	O	S	0	0
			444	281	98	64	1		

- Molecule 37 is a protein called Protein LLP homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
37	NK	67	Total	C	N	O	S	0	0
			581	363	128	88	2		

- Molecule 38 is a protein called Ribosome biogenesis protein NOP53.

Mol	Chain	Residues	Atoms					AltConf	Trace
38	NL	323	Total	C	N	O	S	0	0
			2666	1658	531	475	2		

- Molecule 39 is a protein called Zinc finger protein 593.

Mol	Chain	Residues	Atoms					AltConf	Trace
39	NP	108	Total	C	N	O	S	0	0
			876	537	182	153	4		

- Molecule 40 is a protein called 60S ribosomal protein L4.

Mol	Chain	Residues	Atoms					AltConf	Trace
40	SA	358	Total	C	N	O	S	0	0
			2853	1797	570	473	13		

- Molecule 41 is a protein called 60S ribosomal protein L5.

Mol	Chain	Residues	Atoms					AltConf	Trace
41	SB	292	Total	C	N	O	S	0	0
			2376	1499	433	430	14		

- Molecule 42 is a protein called 60S ribosomal protein L6.

Mol	Chain	Residues	Atoms					AltConf	Trace
42	SC	217	Total	C	N	O	S	0	0
			1747	1124	332	287	4		

- Molecule 43 is a protein called 60S ribosomal protein L7.

Mol	Chain	Residues	Atoms					AltConf	Trace
43	SD	225	Total	C	N	O	S	0	0
			1870	1202	358	301	9		

- Molecule 44 is a protein called 60S ribosomal protein L7a.

Mol	Chain	Residues	Atoms					AltConf	Trace
44	SE	231	Total	C	N	O	S	1	0
			1869	1191	361	313	4		

- Molecule 45 is a protein called 60S ribosomal protein L8.

Mol	Chain	Residues	Atoms					AltConf	Trace
45	SF	245	Total	C	N	O	S	0	0
			1876	1177	383	310	6		

- Molecule 46 is a protein called 60S ribosomal protein L9.

Mol	Chain	Residues	Atoms					AltConf	Trace
46	SG	190	Total	C	N	O	S	0	0
			1518	956	284	272	6		

- Molecule 47 is a protein called MKI67 FHA domain-interacting nucleolar phosphoprotein.

Mol	Chain	Residues	Atoms					AltConf	Trace
47	SH	150	Total	C	N	O	S	1	0
			1275	824	227	220	4		

- Molecule 48 is a protein called 60S ribosomal protein L7-like 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
48	SI	234	Total	C	N	O	S	1	0
			1937	1254	363	316	4		

- Molecule 49 is a protein called Eukaryotic translation initiation factor 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
49	SK	226	Total	C	N	O	S	0	0
			1721	1070	296	343	12		

- Molecule 50 is a protein called Ribosomal L1 domain-containing protein 1.

Mol	Chain	Residues	Atoms					AltConf	Trace
50	SL	238	Total	C	N	O	S	0	0
			1917	1227	337	347	6		

- Molecule 51 is a protein called Pescadillo homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
51	SM	399	Total	C	N	O	S	0	0
			3278	2120	576	571	11		

- Molecule 52 is a protein called mRNA turnover protein 4 homolog.

Mol	Chain	Residues	Atoms				AltConf	Trace
52	SQ	123	Total	C	N	O	0	0
			610	364	123	123		

- Molecule 53 is a protein called GTP-binding protein 4.

Mol	Chain	Residues	Atoms					AltConf	Trace
53	SR	243	Total	C	N	O	S	0	0
			2002	1238	377	375	12		

- Molecule 54 is a protein called Probable ribosome biogenesis protein RLP24.

Mol	Chain	Residues	Atoms					AltConf	Trace
54	SV	139	Total	C	N	O	S	0	0
			1184	754	229	191	10		

- Molecule 55 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
55	L1	5	Total	Mg	0
			5	5	
55	L3	72	Total	Mg	0
			72	72	
55	L4	3	Total	Mg	0
			3	3	

Continued on next page...

Continued from previous page...

Mol	Chain	Residues	Atoms		AltConf
55	LG	1	Total 1	Mg 1	0
55	LQ	1	Total 1	Mg 1	0
55	LR	1	Total 1	Mg 1	0
55	LT	1	Total 1	Mg 1	0
55	LW	1	Total 1	Mg 1	0
55	SA	1	Total 1	Mg 1	0
55	SF	1	Total 1	Mg 1	0

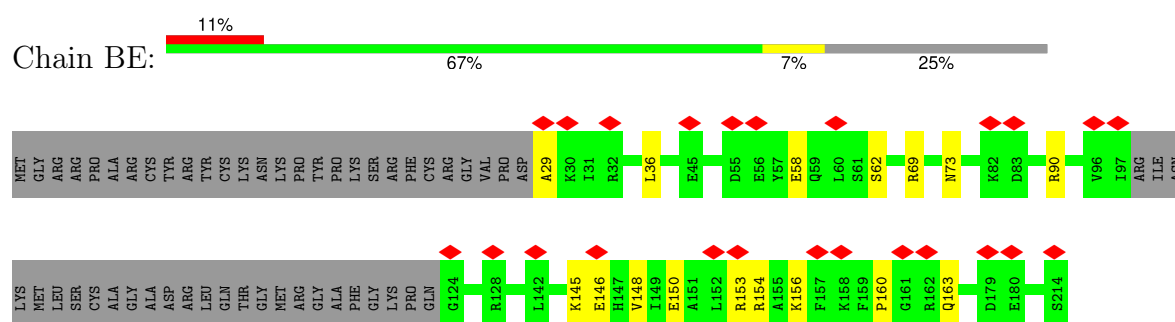
- Molecule 56 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
56	LR	1	Total 1	Zn 1	0
56	LV	1	Total 1	Zn 1	0
56	LW	1	Total 1	Zn 1	0
56	LX	1	Total 1	Zn 1	0
56	NP	1	Total 1	Zn 1	0
56	SV	1	Total 1	Zn 1	0

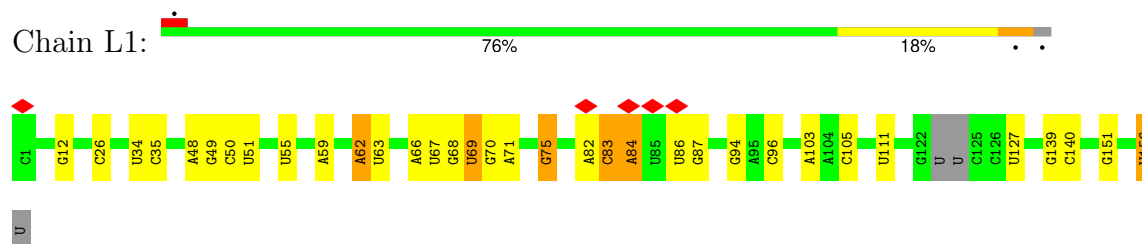
3 Residue-property plots [i](#)

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.

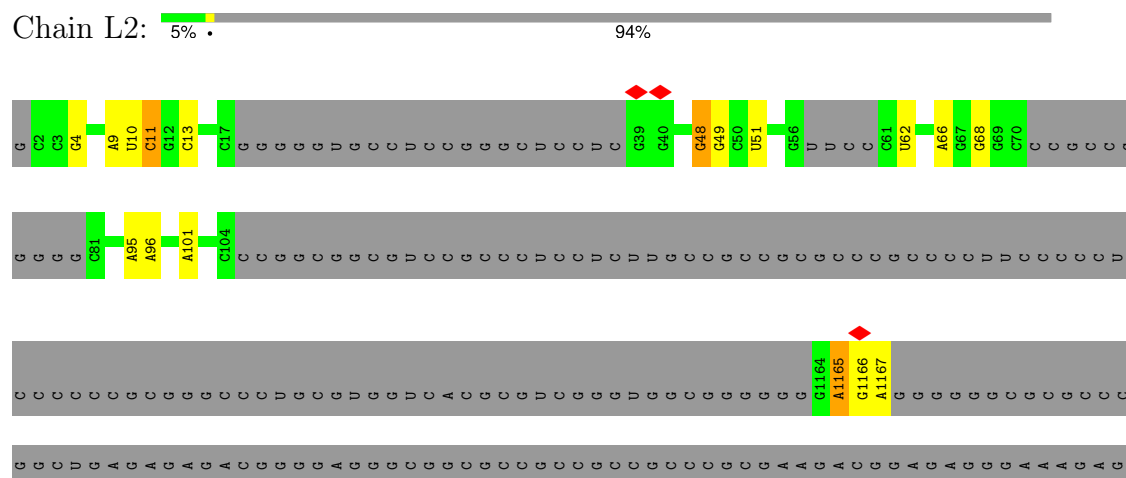
- Molecule 1: 60S ribosomal protein L10

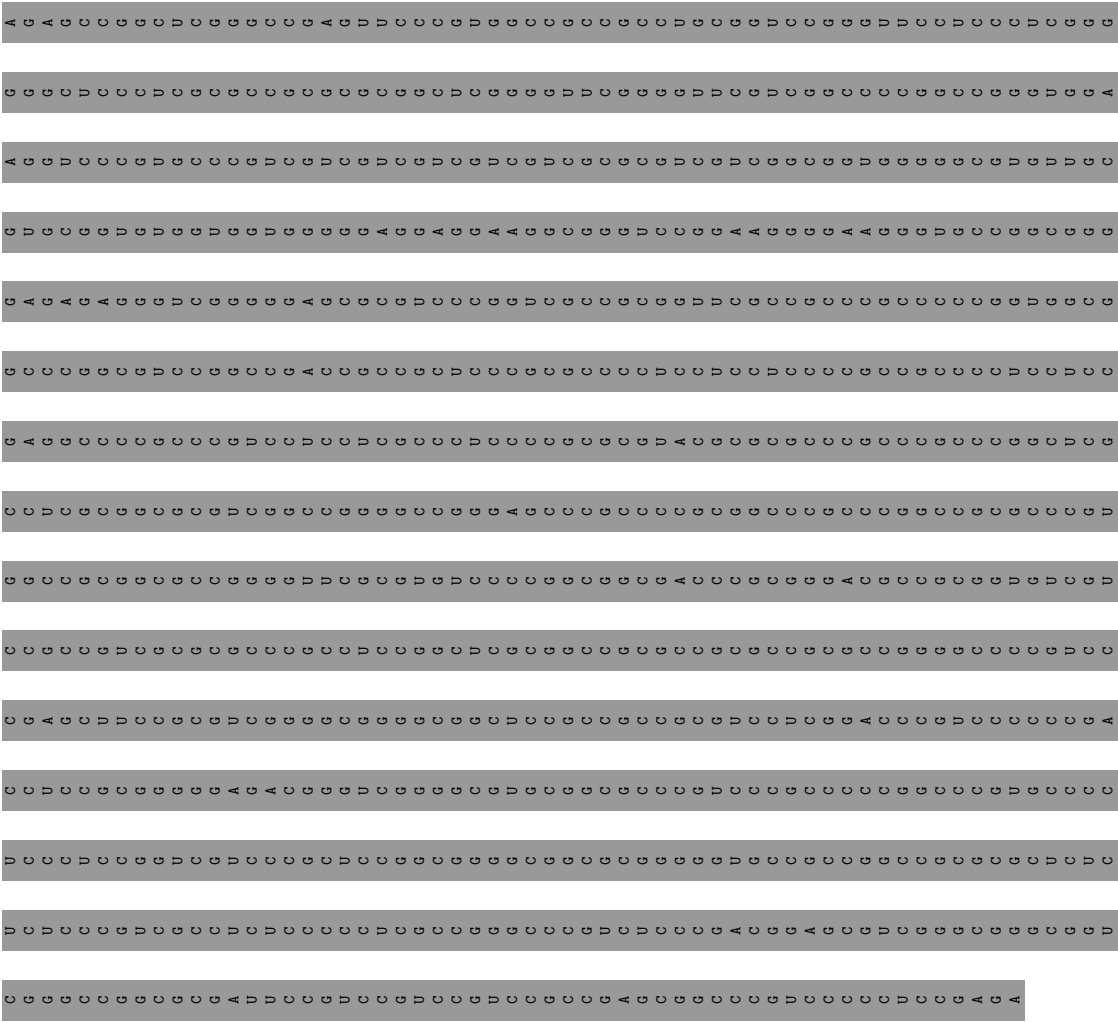


- Molecule 2: 5.8S rRNA

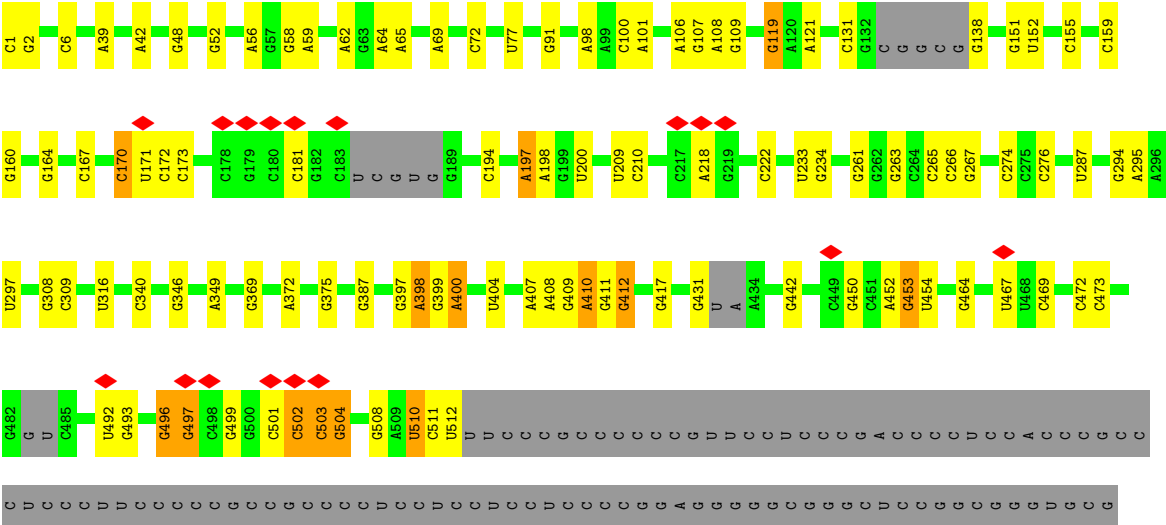


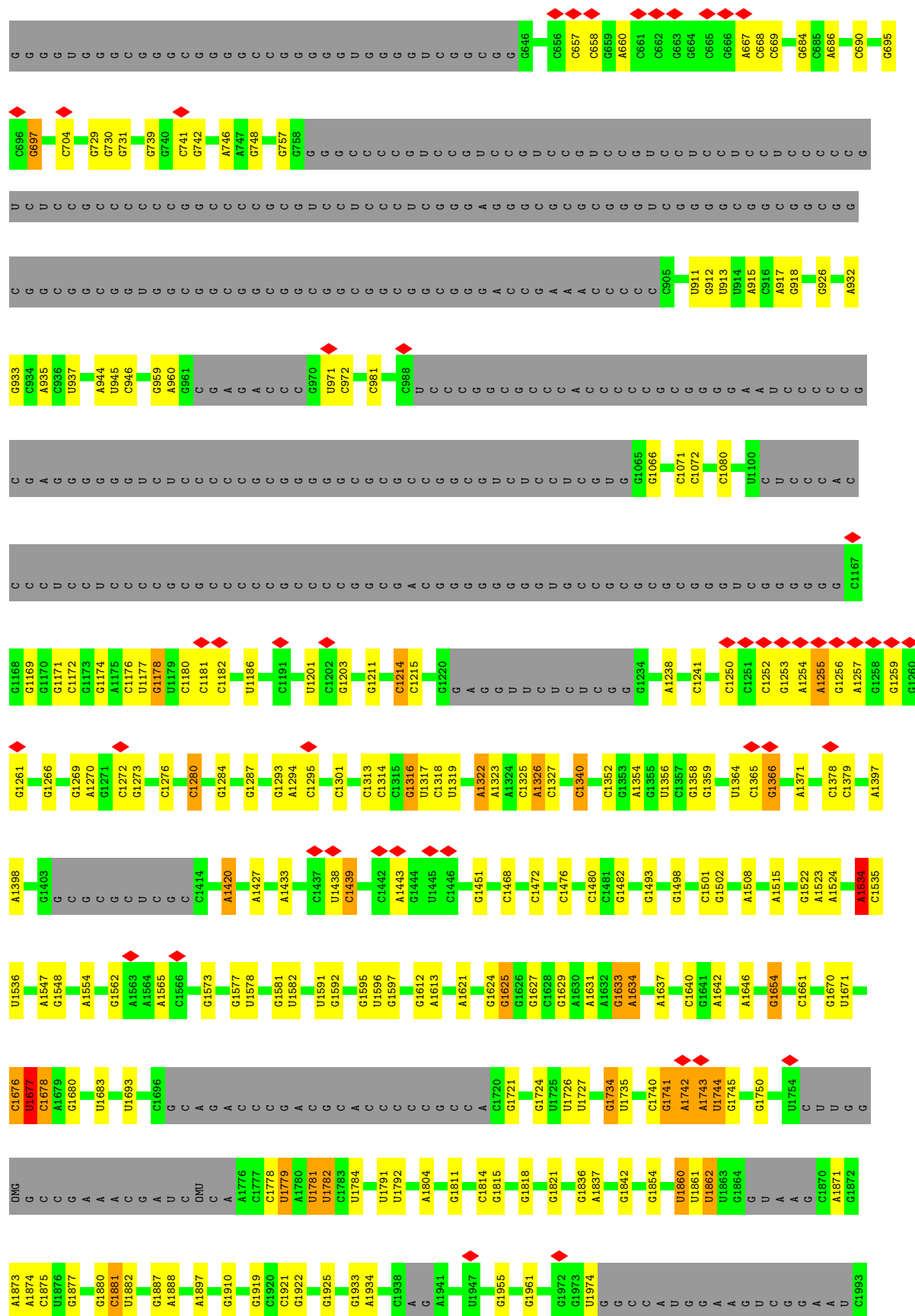
- Molecule 3: ITS2 rRNA



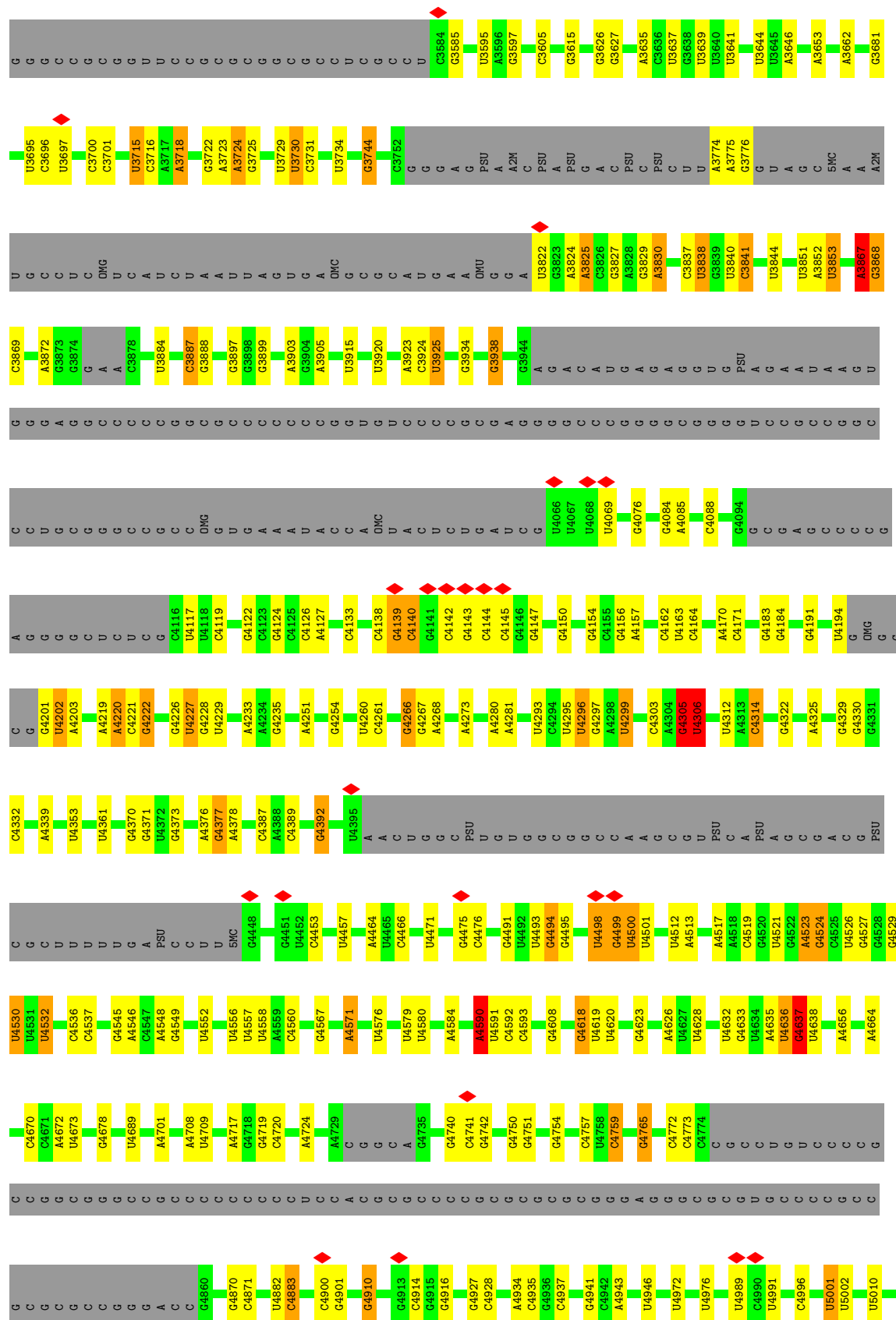


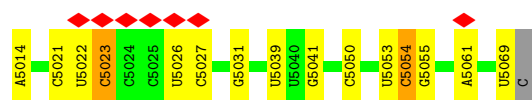
• Molecule 4: 28S rRNA



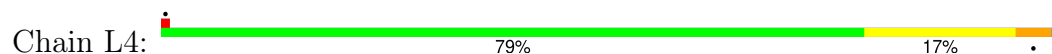




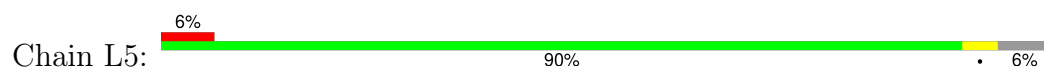




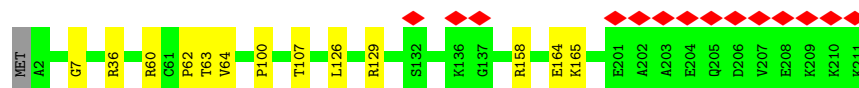
- Molecule 5: 5S rRNA



- Molecule 6: 60S ribosomal protein L11



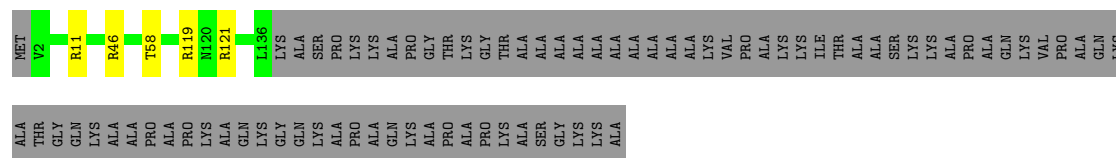
- Molecule 7: 60S ribosomal protein L13



- Molecule 8: 60S ribosomal protein L13a



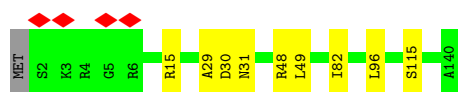
- Molecule 9: 60S ribosomal protein L14



- Molecule 10: 60S ribosomal protein L15



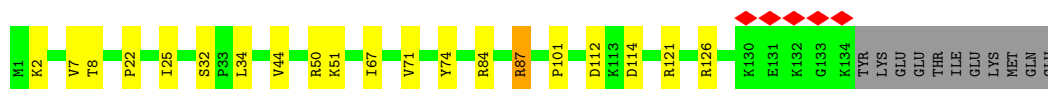
- Chain LG:  93% 6%



- Molecule 18: 60S ribosomal protein L23a



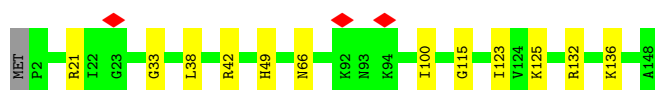
- Molecule 19: 60S ribosomal protein L26



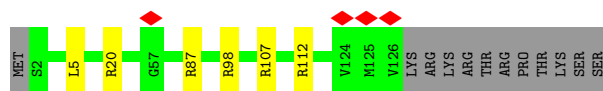
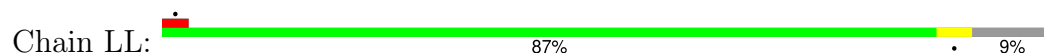
- Molecule 20: 60S ribosomal protein L27



- Molecule 21: 60S ribosomal protein L27a

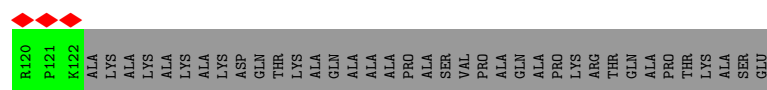


- Molecule 22: 60S ribosomal protein L28

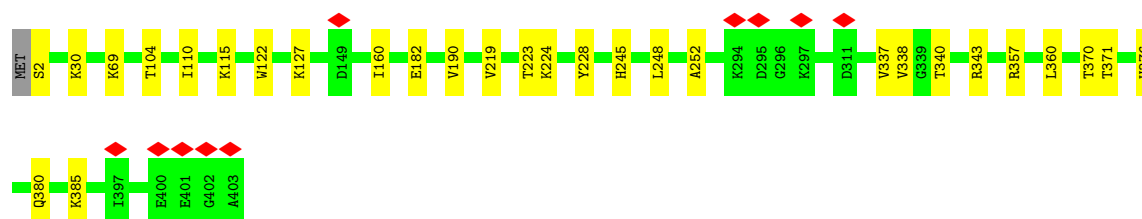


- Molecule 23: 60S ribosomal protein L29

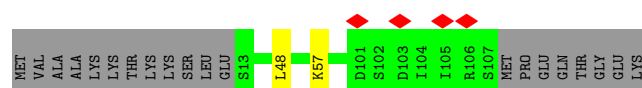
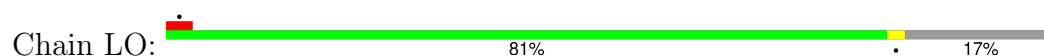




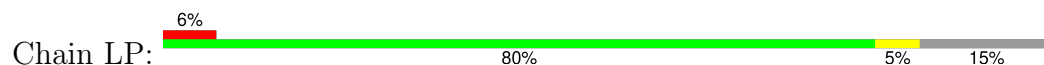
- Molecule 24: 60S ribosomal protein L3



- Molecule 25: 60S ribosomal protein L30



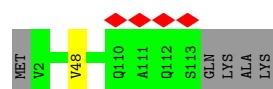
- Molecule 26: 60S ribosomal protein L31



- Molecule 27: 60S ribosomal protein L32

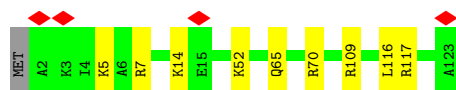


- Molecule 28: 60S ribosomal protein L34



- Molecule 29: 60S ribosomal protein L35





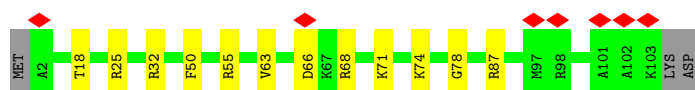
- Molecule 30: 60S ribosomal protein L35a

Chain LT: 95%



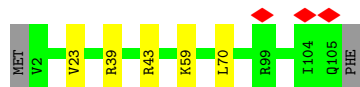
- Molecule 31: 60S ribosomal protein L36

Chain LU: 7% 86% 11%



- Molecule 32: 60S ribosomal protein L36a

Chain LV: 7% 93% 5%



- Molecule 33: 60S ribosomal protein L37

Chain LW: 82% 6% 11%



- Molecule 34: 60S ribosomal protein L37a

Chain LX: 5% 93% 5%




- Molecule 35: 60S ribosomal protein L38

Chain LY: 7% 90% 9%



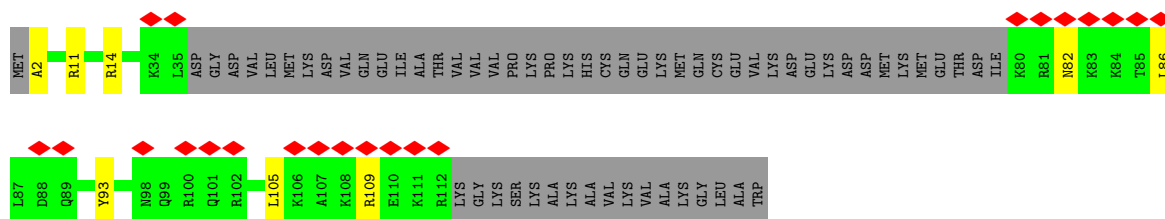
- Molecule 36: 60S ribosomal protein L39

Chain LZ:  88% 10%



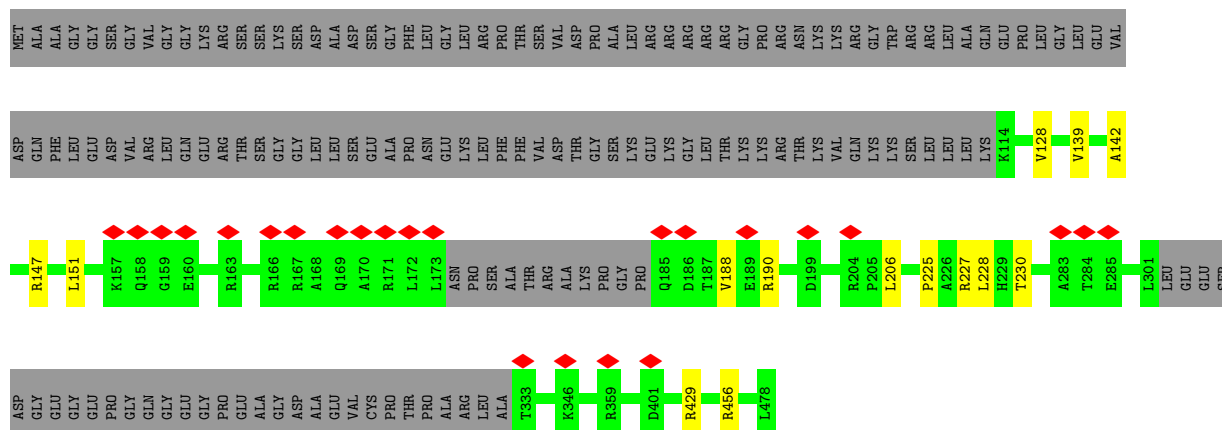
- Molecule 37: Protein LLP homolog

Chain NK:  17% 46% 6% 48%




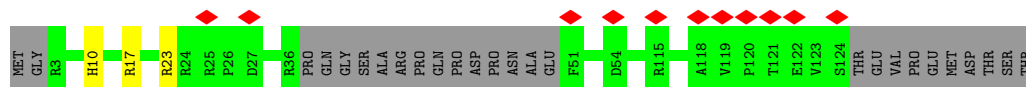
- Molecule 38: Ribosome biogenesis protein NOP53

Chain NL:  5% 65% 32%




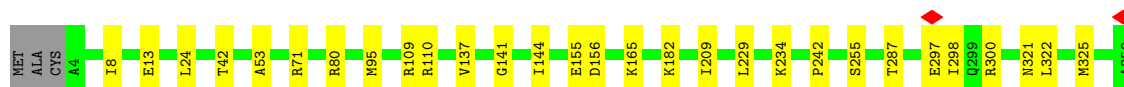
- Molecule 39: Zinc finger protein 593

Chain NP:  8% 78% 19%

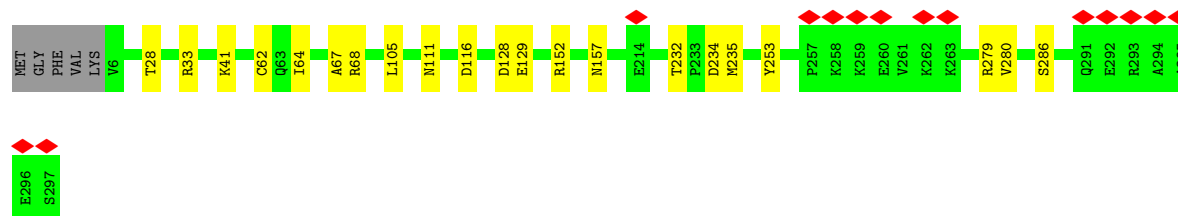


- Molecule 40: 60S ribosomal protein L4

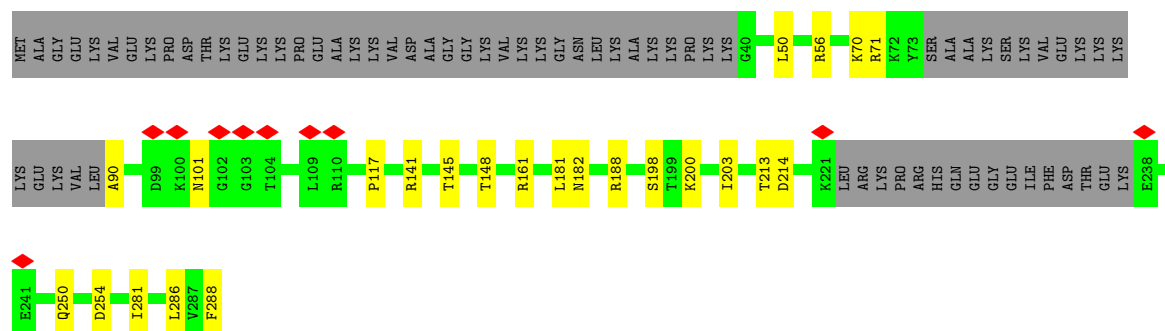
Chain SA:  77% 7% 16%



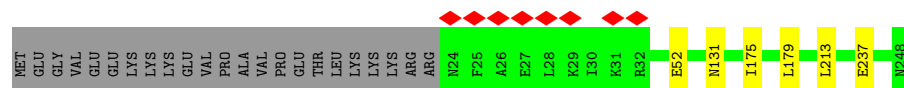
- Molecule 41: 60S ribosomal protein L5



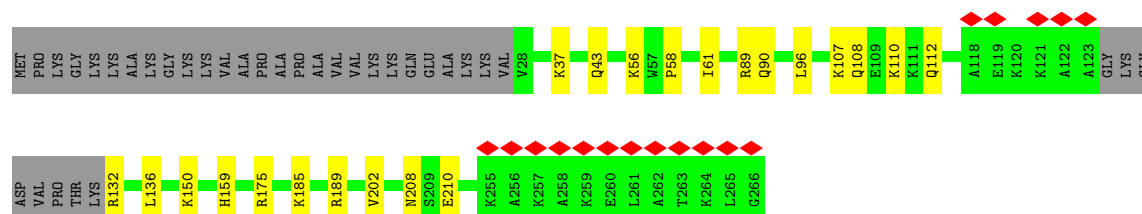
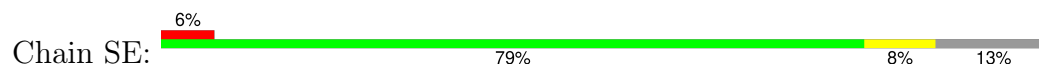
- Molecule 42: 60S ribosomal protein L6



- Molecule 43: 60S ribosomal protein L7



- Molecule 44: 60S ribosomal protein L7a



- | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | | |
|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|------|------|------|------|------|------|------|------|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|
| M1 | A2 | M21 | D44 | T45 | V49 | H50 | R57 | Q79 | L88 | T94 | R95 | R96 | E99 | A103 | V107 | D112 | H118 | T125 | E126 | R139 | Q145 | S155 | M166 | P163 | T186 | L215 | E226 | ALA | GLN | PRO | SER | THR | ILE | ALA | THR | SER | MET | ARG |
|----|----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|------|------|------|------|------|------|------|------|------|------|------|------|------|------|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|-----|

MET	PRO	LYS	SER	LYS	ARG	ASP	LYS	VAL	LEU	THR	LYS	THR	ALA	LYS	LYS	GLY	L20	L30	R31	K32	C33	V34	D35	T36	Y37	F41	I42	F43	S44	V45	A46	M47	M48	R49	L53	K54	D55	I56	R57	N58	A59	W60	K61	H62	S63	R64	M65	F66	N70	V75	A76
-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----	-----

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	88174	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 ($e^-/\text{\AA}^2$)	60	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	64000	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	11.150	Depositor
Minimum map value	-0.181	Depositor
Average map value	0.047	Depositor
Map value standard deviation	0.194	Depositor
Recommended contour level	0.95	Depositor
Map size (\AA)	514.56, 514.56, 514.56	wwPDB
Map dimensions	480, 480, 480	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (\AA)	1.072, 1.072, 1.072	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: A2M, UR3, 6MZ, OMC, OMG, MG, HIC, PSU, 1MA, ZN, OMU

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	BE	0.27	0/1323	0.56	0/1767
2	L1	0.88	0/3589	0.80	0/5589
3	L2	0.63	0/1709	0.87	0/2653
4	L3	0.74	0/78428	0.81	14/122316 (0.0%)
5	L4	1.12	0/2861	0.90	3/4459 (0.1%)
6	L5	0.35	0/1372	0.59	0/1836
7	L6	0.33	0/1732	0.60	0/2315
8	L7	0.39	0/1682	0.57	0/2250
9	L8	0.36	0/1133	0.55	0/1516
10	L9	0.43	0/1746	0.65	0/2338
11	LA	0.35	0/1268	0.55	0/1701
12	LB	0.41	0/1536	0.67	0/2052
13	LC	0.45	0/1501	0.63	0/2013
14	LD	0.31	0/1305	0.62	0/1727
15	LE	0.45	0/1291	0.59	0/1724
16	LF	0.32	0/856	0.53	0/1149
17	LG	0.36	0/1048	0.59	0/1402
18	LH	0.35	0/1175	0.53	0/1572
19	LI	0.35	0/1132	0.59	0/1504
20	LJ	0.41	0/1130	0.57	0/1507
21	LK	0.40	0/1191	0.57	0/1591
22	LL	0.33	0/1017	0.61	0/1364
23	LM	0.33	0/788	0.59	0/1040
24	LN	0.35	0/3294	0.57	0/4406
25	LO	0.34	0/748	0.49	0/1004
26	LP	0.34	0/894	0.60	0/1204
27	LQ	0.36	0/1071	0.57	0/1429
28	LR	0.36	0/898	0.64	0/1197
29	LS	0.35	0/1023	0.59	0/1351
30	LT	0.38	0/895	0.62	0/1198
31	LU	0.31	0/843	0.61	0/1115
32	LV	0.44	0/864	0.63	0/1140

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
33	LW	0.39	0/720	0.68	0/952
34	LX	0.34	0/718	0.56	0/953
35	LY	0.31	0/575	0.53	0/761
36	LZ	0.35	0/454	0.63	0/599
37	NK	0.28	0/587	0.59	0/767
38	NL	0.31	0/2705	0.59	0/3624
39	NP	0.29	0/893	0.62	0/1194
40	SA	0.34	0/2907	0.60	0/3905
41	SB	0.47	0/2421	0.57	0/3241
42	SC	0.31	0/1781	0.56	0/2388
43	SD	0.38	0/1905	0.58	0/2539
44	SE	0.43	0/1903	0.60	0/2559
45	SF	0.37	0/1914	0.62	0/2567
46	SG	0.36	0/1537	0.56	0/2066
47	SH	0.38	0/1309	0.55	0/1756
48	SI	0.36	0/1980	0.56	0/2656
49	SK	0.33	0/1745	0.55	0/2374
50	SL	0.32	0/1949	0.52	0/2621
51	SM	0.42	0/3357	0.54	0/4529
52	SQ	0.22	0/608	0.43	0/845
53	SR	0.28	0/2026	0.56	0/2690
54	SV	0.39	0/1207	0.58	0/1600
All	All	0.62	0/156544	0.73	17/228615 (0.0%)

There are no bond length outliers.

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
5	L4	48	G	N3-C4-N9	6.67	130.00	126.00
4	L3	2486	G	N1-C6-O6	-6.41	116.06	119.90
4	L3	2519	U	O4'-C1'-N1	6.27	113.21	108.20
4	L3	1881	C	C2-N1-C1'	6.12	125.53	118.80
4	L3	2469	C	C2-N1-C1'	6.05	125.46	118.80

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	BE	1295	0	1331	11	0
2	L1	3278	0	1665	21	0
3	L2	1535	0	789	11	0
4	L3	72381	0	36675	297	0
5	L4	2561	0	1295	10	0
6	L5	1349	0	1383	5	0
7	L6	1701	0	1818	15	0
8	L7	1650	0	1794	8	0
9	L8	1111	0	1174	6	0
10	L9	1701	0	1749	7	0
11	LA	1242	0	1269	7	0
12	LB	1512	0	1628	9	0
13	LC	1461	0	1502	4	0
14	LD	1289	0	1429	7	0
15	LE	1264	0	1328	3	0
16	LF	842	0	864	4	0
17	LG	1034	0	1097	8	0
18	LH	1156	0	1268	4	0
19	LI	1115	0	1205	15	0
20	LJ	1107	0	1182	8	0
21	LK	1162	0	1213	11	0
22	LL	1002	0	1068	6	0
23	LM	775	0	841	7	0
24	LN	3239	0	3377	23	0
25	LO	738	0	774	1	0
26	LP	879	0	924	4	0
27	LQ	1053	0	1147	4	0
28	LR	888	0	977	1	0
29	LS	1015	0	1148	9	0
30	LT	876	0	912	4	0
31	LU	832	0	917	10	0
32	LV	851	0	920	5	0
33	LW	705	0	737	5	0
34	LX	708	0	756	6	0
35	LY	569	0	637	5	0
36	LZ	444	0	483	4	0
37	NK	581	0	656	6	0
38	NL	2666	0	2774	11	0
39	NP	876	0	881	3	0
40	SA	2853	0	3028	24	0
41	SB	2376	0	2403	15	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
42	SC	1747	0	1897	17	0
43	SD	1870	0	1996	5	0
44	SE	1869	0	2014	18	0
45	SF	1876	0	1970	12	0
46	SG	1518	0	1601	10	0
47	SH	1275	0	1304	4	0
48	SI	1937	0	2070	13	0
49	SK	1721	0	1695	14	0
50	SL	1917	0	2018	18	0
51	SM	3278	0	3332	14	0
52	SQ	610	0	257	1	0
53	SR	2002	0	2072	9	0
54	SV	1184	0	1248	10	0
55	L1	5	0	0	0	0
55	L3	72	0	0	0	0
55	L4	3	0	0	0	0
55	LG	1	0	0	0	0
55	LQ	1	0	0	0	0
55	LR	1	0	0	0	0
55	LT	1	0	0	0	0
55	LW	1	0	0	0	0
55	SA	1	0	0	0	0
55	SF	1	0	0	0	0
56	LR	1	0	0	0	0
56	LV	1	0	0	0	0
56	LW	1	0	0	0	0
56	LX	1	0	0	0	0
56	NP	1	0	0	0	0
56	SV	1	0	0	0	0
All	All	148569	0	112492	573	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 2.

The worst 5 of 573 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)
2:L1:51:U:OP2	36:LZ:21:ARG:NH2	2.06	0.88
5:L4:40:U:O2	6:L5:75:ARG:NH1	2.06	0.88
4:L3:2520:C:O2	4:L3:2640:G:N2	2.08	0.84
4:L3:2898:G:OP2	14:LD:135:LYS:NZ	2.11	0.83

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
32:LV:23:VAL:HG12	32:LV:70:LEU:HD23	1.59	0.83

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	BE	156/214 (73%)	149 (96%)	7 (4%)	0	100	100
6	L5	166/178 (93%)	164 (99%)	2 (1%)	0	100	100
7	L6	208/211 (99%)	203 (98%)	5 (2%)	0	100	100
8	L7	199/203 (98%)	196 (98%)	3 (2%)	0	100	100
9	L8	133/215 (62%)	130 (98%)	3 (2%)	0	100	100
10	L9	201/204 (98%)	197 (98%)	4 (2%)	0	100	100
11	LA	151/184 (82%)	146 (97%)	5 (3%)	0	100	100
12	LB	185/188 (98%)	180 (97%)	5 (3%)	0	100	100
13	LC	174/176 (99%)	171 (98%)	3 (2%)	0	100	100
14	LD	152/196 (78%)	152 (100%)	0	0	100	100
15	LE	150/160 (94%)	145 (97%)	5 (3%)	0	100	100
16	LF	101/128 (79%)	99 (98%)	2 (2%)	0	100	100
17	LG	137/140 (98%)	132 (96%)	5 (4%)	0	100	100
18	LH	141/156 (90%)	140 (99%)	1 (1%)	0	100	100
19	LI	132/145 (91%)	130 (98%)	2 (2%)	0	100	100
20	LJ	133/136 (98%)	131 (98%)	2 (2%)	0	100	100
21	LK	145/148 (98%)	142 (98%)	3 (2%)	0	100	100
22	LL	123/137 (90%)	120 (98%)	3 (2%)	0	100	100
23	LM	90/159 (57%)	87 (97%)	3 (3%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
24	LN	399/403 (99%)	383 (96%)	16 (4%)	0	100	100
25	LO	93/115 (81%)	93 (100%)	0	0	100	100
26	LP	104/125 (83%)	101 (97%)	3 (3%)	0	100	100
27	LQ	126/135 (93%)	126 (100%)	0	0	100	100
28	LR	110/117 (94%)	110 (100%)	0	0	100	100
29	LS	120/123 (98%)	120 (100%)	0	0	100	100
30	LT	107/110 (97%)	106 (99%)	1 (1%)	0	100	100
31	LU	100/105 (95%)	99 (99%)	1 (1%)	0	100	100
32	LV	102/106 (96%)	98 (96%)	4 (4%)	0	100	100
33	LW	84/97 (87%)	83 (99%)	1 (1%)	0	100	100
34	LX	89/92 (97%)	86 (97%)	3 (3%)	0	100	100
35	LY	67/70 (96%)	67 (100%)	0	0	100	100
36	LZ	48/51 (94%)	47 (98%)	1 (2%)	0	100	100
37	NK	63/129 (49%)	62 (98%)	1 (2%)	0	100	100
38	NL	317/478 (66%)	315 (99%)	2 (1%)	0	100	100
39	NP	104/134 (78%)	102 (98%)	2 (2%)	0	100	100
40	SA	356/427 (83%)	350 (98%)	6 (2%)	0	100	100
41	SB	290/297 (98%)	287 (99%)	3 (1%)	0	100	100
42	SC	211/288 (73%)	202 (96%)	9 (4%)	0	100	100
43	SD	223/248 (90%)	216 (97%)	7 (3%)	0	100	100
44	SE	228/266 (86%)	225 (99%)	3 (1%)	0	100	100
45	SF	243/257 (95%)	232 (96%)	11 (4%)	0	100	100
46	SG	188/192 (98%)	184 (98%)	4 (2%)	0	100	100
47	SH	149/293 (51%)	145 (97%)	4 (3%)	0	100	100
48	SI	231/255 (91%)	228 (99%)	3 (1%)	0	100	100
49	SK	224/245 (91%)	218 (97%)	6 (3%)	0	100	100
50	SL	236/490 (48%)	226 (96%)	10 (4%)	0	100	100
51	SM	393/588 (67%)	387 (98%)	6 (2%)	0	100	100
52	SQ	119/239 (50%)	118 (99%)	1 (1%)	0	100	100
53	SR	237/634 (37%)	235 (99%)	2 (1%)	0	100	100
54	SV	137/163 (84%)	136 (99%)	1 (1%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	8375/10550 (79%)	8201 (98%)	174 (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	
1	BE	137/181 (76%)	137 (100%)	0	100	100
6	L5	142/149 (95%)	142 (100%)	0	100	100
7	L6	176/177 (99%)	176 (100%)	0	100	100
8	L7	173/174 (99%)	173 (100%)	0	100	100
9	L8	115/161 (71%)	115 (100%)	0	100	100
10	L9	171/172 (99%)	171 (100%)	0	100	100
11	LA	134/163 (82%)	134 (100%)	0	100	100
12	LB	164/165 (99%)	164 (100%)	0	100	100
13	LC	157/157 (100%)	157 (100%)	0	100	100
14	LD	138/175 (79%)	138 (100%)	0	100	100
15	LE	136/140 (97%)	136 (100%)	0	100	100
16	LF	93/115 (81%)	93 (100%)	0	100	100
17	LG	106/107 (99%)	106 (100%)	0	100	100
18	LH	124/133 (93%)	124 (100%)	0	100	100
19	LI	124/135 (92%)	123 (99%)	1 (1%)	79	91
20	LJ	117/118 (99%)	117 (100%)	0	100	100
21	LK	120/121 (99%)	120 (100%)	0	100	100
22	LL	109/121 (90%)	109 (100%)	0	100	100
23	LM	80/126 (64%)	78 (98%)	2 (2%)	42	66
24	LN	347/348 (100%)	347 (100%)	0	100	100
25	LO	80/97 (82%)	80 (100%)	0	100	100

Continued on next page...

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
26	LP	97/110 (88%)	97 (100%)	0	100	100
27	LQ	114/121 (94%)	114 (100%)	0	100	100
28	LR	96/100 (96%)	96 (100%)	0	100	100
29	LS	109/110 (99%)	109 (100%)	0	100	100
30	LT	88/89 (99%)	88 (100%)	0	100	100
31	LU	86/89 (97%)	86 (100%)	0	100	100
32	LV	92/94 (98%)	92 (100%)	0	100	100
33	LW	73/80 (91%)	73 (100%)	0	100	100
34	LX	74/75 (99%)	74 (100%)	0	100	100
35	LY	64/65 (98%)	64 (100%)	0	100	100
36	LZ	47/48 (98%)	47 (100%)	0	100	100
37	NK	61/115 (53%)	61 (100%)	0	100	100
38	NL	280/402 (70%)	280 (100%)	0	100	100
39	NP	92/114 (81%)	92 (100%)	0	100	100
40	SA	298/348 (86%)	298 (100%)	0	100	100
41	SB	246/250 (98%)	246 (100%)	0	100	100
42	SC	192/252 (76%)	192 (100%)	0	100	100
43	SD	194/215 (90%)	194 (100%)	0	100	100
44	SE	198/223 (89%)	197 (100%)	1 (0%)	86	95
45	SF	188/199 (94%)	188 (100%)	0	100	100
46	SG	169/171 (99%)	169 (100%)	0	100	100
47	SH	141/274 (52%)	141 (100%)	0	100	100
48	SI	210/228 (92%)	210 (100%)	0	100	100
49	SK	196/213 (92%)	195 (100%)	1 (0%)	86	95
50	SL	221/437 (51%)	221 (100%)	0	100	100
51	SM	354/509 (70%)	354 (100%)	0	100	100
53	SR	218/574 (38%)	217 (100%)	1 (0%)	86	95
54	SV	128/149 (86%)	128 (100%)	0	100	100
All	All	7269/8889 (82%)	7263 (100%)	6 (0%)	92	98

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
44	SE	175	ARG
49	SK	57	ARG
53	SR	488	ARG
23	LM	14	ARG
19	LI	87	ARG

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 15 such sidechains are listed below:

Mol	Chain	Res	Type
41	SB	244	HIS
48	SI	225	HIS
42	SC	101	ASN
51	SM	170	HIS
44	SE	112	GLN

5.3.3 RNA [i](#)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
2	L1	152/157 (96%)	17 (11%)	0
3	L2	67/1167 (5%)	9 (13%)	0
4	L3	3339/5070 (65%)	441 (13%)	6 (0%)
5	L4	119/121 (98%)	13 (10%)	1 (0%)
All	All	3677/6515 (56%)	480 (13%)	7 (0%)

5 of 480 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
2	L1	34	U
2	L1	35	C
2	L1	49	G
2	L1	59	A
2	L1	62	A

5 of 7 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	L3	1625	OMG
4	L3	1633	G
5	L4	109	U
4	L3	2095	A
4	L3	503	C

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

110 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
4	PSU	L3	1677	4	18,21,22	1.07	2 (11%)	21,30,33	1.94	4 (19%)
4	6MZ	L3	4220	4	17,25,26	1.28	3 (17%)	15,36,39	2.35	4 (26%)
4	PSU	L3	3695	4	18,21,22	1.12	2 (11%)	21,30,33	1.98	5 (23%)
4	OMC	L3	3869	4	19,22,23	0.70	0	25,31,34	0.76	0
4	PSU	L3	1860	4	18,21,22	1.01	2 (11%)	21,30,33	1.92	4 (19%)
4	PSU	L3	4293	4	18,21,22	1.05	2 (11%)	21,30,33	1.76	4 (19%)
4	OMC	L3	4456	4	19,22,23	0.66	0	25,31,34	0.60	0
4	OMC	L3	1340	4	19,22,23	0.75	1 (5%)	25,31,34	0.72	0
4	PSU	L3	1781	4	18,21,22	1.10	1 (5%)	21,30,33	1.99	5 (23%)
4	PSU	L3	3884	4	18,21,22	1.11	2 (11%)	21,30,33	2.02	5 (23%)
4	OMG	L3	4228	4	19,26,27	1.48	3 (15%)	21,38,41	0.99	1 (4%)
4	PSU	L3	3639	4	18,21,22	1.06	2 (11%)	21,30,33	1.89	4 (19%)
4	A2M	L3	4523	4	18,25,26	1.30	3 (16%)	20,36,39	1.27	1 (5%)
4	OMC	L3	2824	4	19,22,23	0.65	0	25,31,34	0.66	0
4	OMG	L3	4494	4	19,26,27	1.35	3 (15%)	21,38,41	0.81	1 (4%)
4	OMC	L3	2351	55,4	19,22,23	0.73	1 (5%)	25,31,34	0.85	1 (4%)
4	A2M	L3	2401	4	18,25,26	1.32	3 (16%)	20,36,39	1.41	1 (5%)
4	PSU	L3	2508	4	18,21,22	1.03	2 (11%)	21,30,33	2.04	5 (23%)
4	OMG	L3	3899	4	19,26,27	1.32	3 (15%)	21,38,41	0.91	1 (4%)
4	PSU	L3	1536	4	18,21,22	1.04	2 (11%)	21,30,33	2.00	5 (23%)
4	A2M	L3	2787	4	18,25,26	1.31	3 (16%)	20,36,39	1.51	2 (10%)
4	PSU	L3	1582	4	18,21,22	1.03	2 (11%)	21,30,33	2.09	5 (23%)
4	A2M	L3	2363	55,4	18,25,26	1.38	3 (16%)	20,36,39	1.36	1 (5%)
4	OMC	L3	2804	4	19,22,23	0.69	0	25,31,34	0.65	0
4	PSU	L3	4493	4	18,21,22	1.05	1 (5%)	21,30,33	2.02	2 (9%)
4	PSU	L3	4972	4	18,21,22	1.03	1 (5%)	21,30,33	1.90	4 (19%)
4	A2M	L3	1524	4	18,25,26	1.33	3 (16%)	20,36,39	1.66	3 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	OMG	L3	1316	4	19,26,27	1.27	3 (15%)	21,38,41	0.86	1 (4%)
4	A2M	L3	3830	4	18,25,26	1.34	3 (16%)	20,36,39	1.42	2 (10%)
4	OMG	L3	1625	4	19,26,27	1.30	3 (15%)	21,38,41	0.85	1 (4%)
4	OMG	L3	2424	4	19,26,27	1.31	3 (15%)	21,38,41	0.76	1 (4%)
4	OMC	L3	3887	4	19,22,23	0.64	0	25,31,34	0.61	0
4	PSU	L3	1779	4	18,21,22	1.09	1 (5%)	21,30,33	2.01	6 (28%)
4	PSU	L3	3730	4	18,21,22	1.08	1 (5%)	21,30,33	1.98	5 (23%)
4	PSU	L3	2632	4	18,21,22	1.06	1 (5%)	21,30,33	1.92	3 (14%)
4	OMU	L3	2415	4	19,22,23	1.95	6 (31%)	25,31,34	1.97	5 (20%)
4	A2M	L3	4571	4	18,25,26	1.34	3 (16%)	20,36,39	1.26	1 (5%)
4	1MA	L3	1322	4	17,25,26	0.92	2 (11%)	17,37,40	1.12	2 (11%)
4	PSU	L3	4689	4	18,21,22	1.06	2 (11%)	21,30,33	1.93	3 (14%)
4	PSU	L3	1792	4	18,21,22	1.04	2 (11%)	21,30,33	2.03	4 (19%)
4	A2M	L3	3867	4	18,25,26	1.32	3 (16%)	20,36,39	1.63	2 (10%)
4	PSU	L3	4471	4	18,21,22	1.11	2 (11%)	21,30,33	1.95	5 (23%)
4	A2M	L3	3718	4	18,25,26	1.36	3 (16%)	20,36,39	1.23	1 (5%)
4	PSU	L3	1744	4	18,21,22	1.09	1 (5%)	21,30,33	1.99	5 (23%)
4	OMC	L3	2365	4	19,22,23	0.68	0	25,31,34	0.59	0
4	PSU	L3	4500	4	18,21,22	1.11	1 (5%)	21,30,33	1.97	5 (23%)
2	PSU	L1	69	2	18,21,22	1.09	2 (11%)	21,30,33	1.95	5 (23%)
4	OMU	L3	4498	4	19,22,23	2.10	7 (36%)	25,31,34	1.80	5 (20%)
4	OMU	L3	4227	4	19,22,23	1.90	5 (26%)	25,31,34	1.87	5 (20%)
4	A2M	L3	4590	4	18,25,26	1.37	3 (16%)	20,36,39	1.57	3 (15%)
4	PSU	L3	3715	4	18,21,22	1.08	2 (11%)	21,30,33	2.01	6 (28%)
4	PSU	L3	4361	4	18,21,22	1.12	2 (11%)	21,30,33	2.06	5 (23%)
2	PSU	L1	55	2	18,21,22	1.05	2 (11%)	21,30,33	2.10	5 (23%)
4	PSU	L3	1862	4	18,21,22	1.05	1 (5%)	21,30,33	2.00	5 (23%)
4	OMG	L3	3627	4	19,26,27	1.26	3 (15%)	21,38,41	0.85	1 (4%)
4	OMC	L3	3701	4	19,22,23	0.59	0	25,31,34	0.65	0
4	OMG	L3	1522	4	19,26,27	1.32	3 (15%)	21,38,41	0.88	1 (4%)
24	HIC	LN	245	24	8,11,12	1.60	2 (25%)	5,14,16	0.86	0
4	OMG	L3	4637	4	19,26,27	1.27	3 (15%)	21,38,41	0.90	1 (4%)
4	PSU	L3	4457	4	18,21,22	1.11	2 (11%)	21,30,33	2.03	5 (23%)
4	PSU	L3	3851	4	18,21,22	1.09	1 (5%)	21,30,33	1.95	4 (19%)
4	PSU	L3	5010	4	18,21,22	1.06	2 (11%)	21,30,33	1.91	4 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PSU	L3	4521	4	18,21,22	1.04	2 (11%)	21,30,33	1.94	4 (19%)
4	PSU	L3	3920	55,4	18,21,22	1.11	2 (11%)	21,30,33	1.93	4 (19%)
4	PSU	L3	4576	4	18,21,22	1.10	2 (11%)	21,30,33	2.05	6 (28%)
4	PSU	L3	3853	4	18,21,22	1.01	2 (11%)	21,30,33	1.89	5 (23%)
4	PSU	L3	3822	4	18,21,22	1.18	1 (5%)	21,30,33	1.96	5 (23%)
4	OMC	L3	2861	4	19,22,23	0.62	0	25,31,34	0.63	0
4	OMG	L3	4623	4	19,26,27	1.28	3 (15%)	21,38,41	0.91	1 (4%)
4	A2M	L3	1326	4	18,25,26	1.38	3 (16%)	20,36,39	1.71	4 (20%)
4	OMG	L3	4370	4	19,26,27	1.37	3 (15%)	21,38,41	1.07	2 (9%)
4	A2M	L3	1534	55,4	18,25,26	1.33	3 (16%)	20,36,39	1.53	4 (20%)
4	OMG	L3	3744	4	19,26,27	1.21	3 (15%)	21,38,41	0.89	1 (4%)
4	PSU	L3	4353	4	18,21,22	1.09	2 (11%)	21,30,33	2.07	6 (28%)
4	PSU	L3	4628	4	18,21,22	1.00	2 (11%)	21,30,33	2.01	6 (28%)
4	PSU	L3	1683	4	18,21,22	1.22	3 (16%)	21,30,33	2.17	5 (23%)
4	PSU	L3	2839	4	18,21,22	1.06	2 (11%)	21,30,33	1.91	4 (19%)
4	A2M	L3	3724	4	18,25,26	1.29	3 (16%)	20,36,39	1.34	1 (5%)
4	OMC	L3	3841	4	19,22,23	0.65	0	25,31,34	0.75	1 (4%)
4	A2M	L3	1871	4	18,25,26	1.31	2 (11%)	20,36,39	1.49	2 (10%)
4	PSU	L3	1782	4	18,21,22	1.09	1 (5%)	21,30,33	1.98	5 (23%)
4	PSU	L3	4312	4	18,21,22	1.07	2 (11%)	21,30,33	2.11	5 (23%)
4	PSU	L3	4673	4	18,21,22	1.10	2 (11%)	21,30,33	1.99	5 (23%)
4	OMG	L3	4392	4	19,26,27	1.31	3 (15%)	21,38,41	0.85	1 (4%)
4	OMG	L3	4499	4	19,26,27	1.11	2 (10%)	21,38,41	0.84	1 (4%)
4	OMC	L3	2422	55,4	19,22,23	0.64	0	25,31,34	0.80	1 (4%)
4	UR3	L3	4530	4	19,22,23	1.22	3 (15%)	26,32,35	1.49	2 (7%)
2	OMG	L1	75	2	19,26,27	1.22	3 (15%)	21,38,41	0.86	1 (4%)
4	OMG	L3	2876	4	19,26,27	1.26	3 (15%)	21,38,41	0.81	1 (4%)
4	OMG	L3	4618	4	19,26,27	1.29	3 (15%)	21,38,41	0.92	1 (4%)
4	OMU	L3	2837	4	19,22,23	1.97	6 (31%)	25,31,34	1.92	5 (20%)
4	OMG	L3	2364	4	19,26,27	1.27	3 (15%)	21,38,41	0.92	1 (4%)
4	OMU	L3	3925	4	19,22,23	1.89	6 (31%)	25,31,34	1.96	5 (20%)
4	A2M	L3	398	4	18,25,26	1.35	3 (16%)	20,36,39	1.41	2 (10%)
4	OMU	L3	4306	4	19,22,23	1.90	5 (26%)	25,31,34	2.08	4 (16%)
4	PSU	L3	3637	4	18,21,22	1.01	1 (5%)	21,30,33	1.98	5 (23%)
4	A2M	L3	400	4	18,25,26	1.34	3 (16%)	20,36,39	1.41	2 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	PSU	L3	4579	4	18,21,22	1.01	2 (11%)	21,30,33	1.83	4 (19%)
4	PSU	L3	3734	4	18,21,22	1.08	1 (5%)	21,30,33	1.99	5 (23%)
4	A2M	L3	3825	4	18,25,26	1.34	3 (16%)	20,36,39	1.26	1 (5%)
4	OMC	L3	4536	4	19,22,23	0.65	0	25,31,34	0.68	0
4	OMU	L3	4620	4	19,22,23	1.93	5 (26%)	25,31,34	1.74	4 (16%)
4	PSU	L3	4296	4	18,21,22	1.09	2 (11%)	21,30,33	2.13	6 (28%)
4	PSU	L3	3844	4	18,21,22	1.12	2 (11%)	21,30,33	2.03	5 (23%)
4	PSU	L3	4636	4	18,21,22	1.05	1 (5%)	21,30,33	2.14	6 (28%)
4	A2M	L3	2815	4	18,25,26	1.30	3 (16%)	20,36,39	1.29	2 (10%)
4	PSU	L3	4532	4	18,21,22	1.11	2 (11%)	21,30,33	2.11	5 (23%)
4	PSU	L3	4552	4	18,21,22	1.03	2 (11%)	21,30,33	2.01	5 (23%)
4	PSU	L3	4299	4	18,21,22	1.04	2 (11%)	21,30,33	2.11	3 (14%)
4	PSU	L3	5001	4	18,21,22	1.13	3 (16%)	21,30,33	2.09	5 (23%)

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
4	PSU	L3	1677	4	-	3/7/25/26	0/2/2/2
4	6MZ	L3	4220	4	-	2/5/27/28	0/3/3/3
4	PSU	L3	3695	4	-	0/7/25/26	0/2/2/2
4	OMC	L3	3869	4	-	0/9/27/28	0/2/2/2
4	PSU	L3	1860	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	4293	4	-	1/7/25/26	0/2/2/2
4	OMC	L3	4456	4	-	0/9/27/28	0/2/2/2
4	OMC	L3	1340	4	-	1/9/27/28	0/2/2/2
4	PSU	L3	1781	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	3884	4	-	1/7/25/26	0/2/2/2
4	OMG	L3	4228	4	-	0/5/27/28	0/3/3/3
4	PSU	L3	3639	4	-	0/7/25/26	0/2/2/2
4	A2M	L3	4523	4	-	2/5/27/28	0/3/3/3
4	OMC	L3	2824	4	-	0/9/27/28	0/2/2/2
4	OMG	L3	4494	4	-	0/5/27/28	0/3/3/3
4	OMC	L3	2351	55,4	-	4/9/27/28	0/2/2/2
4	A2M	L3	2401	4	-	2/5/27/28	0/3/3/3
4	PSU	L3	2508	4	-	0/7/25/26	0/2/2/2
4	OMG	L3	3899	4	-	0/5/27/28	0/3/3/3

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PSU	L3	1536	4	-	0/7/25/26	0/2/2/2
4	A2M	L3	2787	4	-	2/5/27/28	0/3/3/3
4	PSU	L3	1582	4	-	0/7/25/26	0/2/2/2
4	A2M	L3	2363	55,4	-	1/5/27/28	0/3/3/3
4	OMC	L3	2804	4	-	1/9/27/28	0/2/2/2
4	PSU	L3	4493	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	4972	4	-	0/7/25/26	0/2/2/2
4	A2M	L3	1524	4	-	0/5/27/28	0/3/3/3
4	OMG	L3	1316	4	-	1/5/27/28	0/3/3/3
4	A2M	L3	3830	4	-	1/5/27/28	0/3/3/3
4	OMG	L3	1625	4	-	2/5/27/28	0/3/3/3
4	OMG	L3	2424	4	-	2/5/27/28	0/3/3/3
4	OMC	L3	3887	4	-	2/9/27/28	0/2/2/2
4	PSU	L3	1779	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	3730	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	2632	4	-	0/7/25/26	0/2/2/2
4	OMU	L3	2415	4	-	0/9/27/28	0/2/2/2
4	A2M	L3	4571	4	-	1/5/27/28	0/3/3/3
4	1MA	L3	1322	4	-	2/3/25/26	0/3/3/3
4	PSU	L3	4689	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	1792	4	-	0/7/25/26	0/2/2/2
4	A2M	L3	3867	4	-	3/5/27/28	0/3/3/3
4	PSU	L3	4471	4	-	0/7/25/26	0/2/2/2
4	A2M	L3	3718	4	-	1/5/27/28	0/3/3/3
4	PSU	L3	1744	4	-	0/7/25/26	0/2/2/2
4	OMC	L3	2365	4	-	0/9/27/28	0/2/2/2
4	PSU	L3	4500	4	-	0/7/25/26	0/2/2/2
2	PSU	L1	69	2	-	0/7/25/26	0/2/2/2
4	OMU	L3	4498	4	-	0/9/27/28	0/2/2/2
4	OMU	L3	4227	4	-	0/9/27/28	0/2/2/2
4	A2M	L3	4590	4	-	3/5/27/28	0/3/3/3
4	PSU	L3	3715	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	4361	4	-	0/7/25/26	0/2/2/2
2	PSU	L1	55	2	-	0/7/25/26	0/2/2/2
4	PSU	L3	1862	4	-	2/7/25/26	0/2/2/2
4	OMG	L3	3627	4	-	0/5/27/28	0/3/3/3
4	OMC	L3	3701	4	-	6/9/27/28	0/2/2/2
4	OMG	L3	1522	4	-	0/5/27/28	0/3/3/3
24	HIC	LN	245	24	-	0/5/6/8	0/1/1/1
4	OMG	L3	4637	4	-	1/5/27/28	0/3/3/3

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PSU	L3	4457	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	3851	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	5010	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	4521	4	-	2/7/25/26	0/2/2/2
4	PSU	L3	3920	55,4	-	0/7/25/26	0/2/2/2
4	PSU	L3	4576	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	3853	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	3822	4	-	0/7/25/26	0/2/2/2
4	OMC	L3	2861	4	-	1/9/27/28	0/2/2/2
4	OMG	L3	4623	4	-	0/5/27/28	0/3/3/3
4	A2M	L3	1326	4	-	3/5/27/28	0/3/3/3
4	OMG	L3	4370	4	-	0/5/27/28	0/3/3/3
4	A2M	L3	1534	55,4	-	1/5/27/28	0/3/3/3
4	OMG	L3	3744	4	-	1/5/27/28	0/3/3/3
4	PSU	L3	4353	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	4628	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	1683	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	2839	4	-	3/7/25/26	0/2/2/2
4	A2M	L3	3724	4	-	1/5/27/28	0/3/3/3
4	OMC	L3	3841	4	-	1/9/27/28	0/2/2/2
4	A2M	L3	1871	4	-	1/5/27/28	0/3/3/3
4	PSU	L3	1782	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	4312	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	4673	4	-	0/7/25/26	0/2/2/2
4	OMG	L3	4392	4	-	1/5/27/28	0/3/3/3
4	OMG	L3	4499	4	-	1/5/27/28	0/3/3/3
4	OMC	L3	2422	55,4	-	2/9/27/28	0/2/2/2
4	UR3	L3	4530	4	-	0/7/25/26	0/2/2/2
2	OMG	L1	75	2	-	1/5/27/28	0/3/3/3
4	OMG	L3	2876	4	-	0/5/27/28	0/3/3/3
4	OMG	L3	4618	4	-	4/5/27/28	0/3/3/3
4	OMU	L3	2837	4	-	1/9/27/28	0/2/2/2
4	OMG	L3	2364	4	-	2/5/27/28	0/3/3/3
4	OMU	L3	3925	4	-	1/9/27/28	0/2/2/2
4	A2M	L3	398	4	-	1/5/27/28	0/3/3/3
4	OMU	L3	4306	4	-	4/9/27/28	0/2/2/2
4	PSU	L3	3637	4	-	0/7/25/26	0/2/2/2
4	A2M	L3	400	4	-	1/5/27/28	0/3/3/3
4	PSU	L3	4579	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	3734	4	-	0/7/25/26	0/2/2/2
4	A2M	L3	3825	4	-	1/5/27/28	0/3/3/3

Continued on next page...

Continued from previous page...

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	OMC	L3	4536	4	-	0/9/27/28	0/2/2/2
4	OMU	L3	4620	4	-	0/9/27/28	0/2/2/2
4	PSU	L3	4296	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	3844	4	-	1/7/25/26	0/2/2/2
4	PSU	L3	4636	4	-	1/7/25/26	0/2/2/2
4	A2M	L3	2815	4	-	1/5/27/28	0/3/3/3
4	PSU	L3	4532	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	4552	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	4299	4	-	0/7/25/26	0/2/2/2
4	PSU	L3	5001	4	-	0/7/25/26	0/2/2/2

The worst 5 of 245 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	L3	4498	OMU	C6-N1	4.86	1.49	1.38
4	L3	4620	OMU	C6-N1	4.51	1.48	1.38
4	L3	2415	OMU	C6-N1	4.48	1.48	1.38
4	L3	2837	OMU	C6-N1	4.46	1.48	1.38
4	L3	4498	OMU	C2-N1	4.32	1.45	1.38

The worst 5 of 334 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	L3	4220	6MZ	C2-N1-C6	6.54	121.67	116.60
4	L3	4306	OMU	C4-N3-C2	-5.94	119.24	126.61
4	L3	3925	OMU	C4-N3-C2	-5.92	119.26	126.61
4	L3	2837	OMU	C4-N3-C2	-5.73	119.50	126.61
4	L3	4227	OMU	C4-N3-C2	-5.67	119.57	126.61

There are no chirality outliers.

5 of 84 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	L1	75	OMG	C1'-C2'-O2'-CM2
4	L3	398	A2M	C1'-C2'-O2'-CM'
4	L3	400	A2M	C1'-C2'-O2'-CM'
4	L3	1316	OMG	C1'-C2'-O2'-CM2
4	L3	1326	A2M	C1'-C2'-O2'-CM'

There are no ring outliers.

57 monomers are involved in 85 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L3	1677	PSU	3	0
4	L3	4220	6MZ	1	0
4	L3	3869	OMC	1	0
4	L3	1860	PSU	1	0
4	L3	1340	OMC	1	0
4	L3	1781	PSU	1	0
4	L3	4494	OMG	1	0
4	L3	2351	OMC	1	0
4	L3	2787	A2M	1	0
4	L3	2363	A2M	3	0
4	L3	2804	OMC	1	0
4	L3	1316	OMG	1	0
4	L3	3830	A2M	1	0
4	L3	2424	OMG	1	0
4	L3	3887	OMC	2	0
4	L3	1779	PSU	1	0
4	L3	3730	PSU	2	0
4	L3	2632	PSU	1	0
4	L3	2415	OMU	1	0
4	L3	4571	A2M	1	0
4	L3	3867	A2M	3	0
4	L3	3718	A2M	1	0
4	L3	1744	PSU	4	0
4	L3	4500	PSU	2	0
2	L1	69	PSU	2	0
4	L3	4227	OMU	3	0
4	L3	4590	A2M	1	0
4	L3	3715	PSU	1	0
4	L3	4637	OMG	2	0
4	L3	3853	PSU	2	0
4	L3	2861	OMC	1	0
4	L3	1326	A2M	5	0
4	L3	1534	A2M	2	0
4	L3	3744	OMG	1	0
4	L3	2839	PSU	1	0
4	L3	3724	A2M	2	0
4	L3	3841	OMC	1	0
4	L3	1782	PSU	1	0
4	L3	4392	OMG	1	0
4	L3	4499	OMG	1	0
4	L3	2422	OMC	1	0
4	L3	4530	UR3	2	0
2	L1	75	OMG	2	0

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	L3	2876	OMG	1	0
4	L3	4618	OMG	2	0
4	L3	2837	OMU	1	0
4	L3	2364	OMG	2	0
4	L3	3925	OMU	2	0
4	L3	398	A2M	2	0
4	L3	4306	OMU	1	0
4	L3	400	A2M	1	0
4	L3	3825	A2M	1	0
4	L3	4536	OMC	1	0
4	L3	4296	PSU	1	0
4	L3	2815	A2M	2	0
4	L3	4299	PSU	1	0
4	L3	5001	PSU	1	0

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 93 ligands modelled in this entry, 93 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.

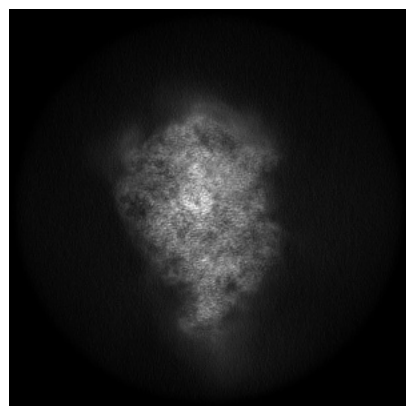
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-29275. These allow visual inspection of the internal detail of the map and identification of artifacts.

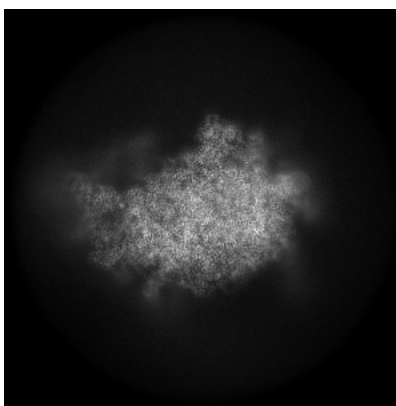
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

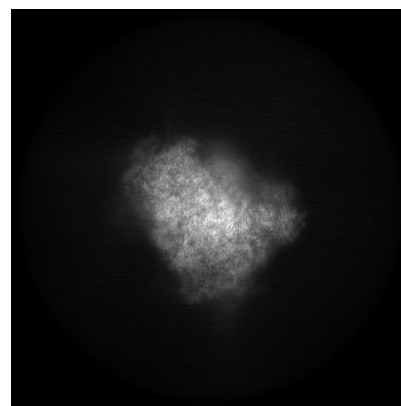
6.1.1 Primary map



X

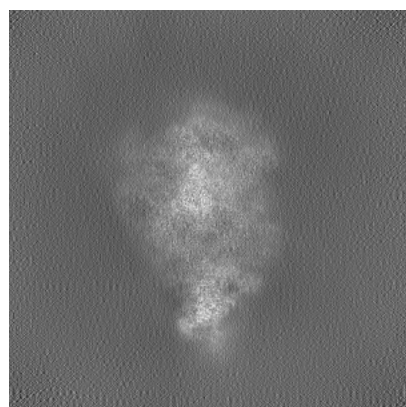


Y

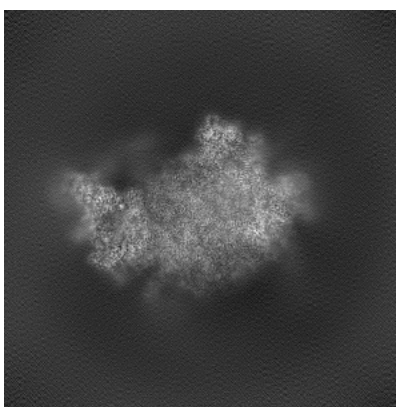


Z

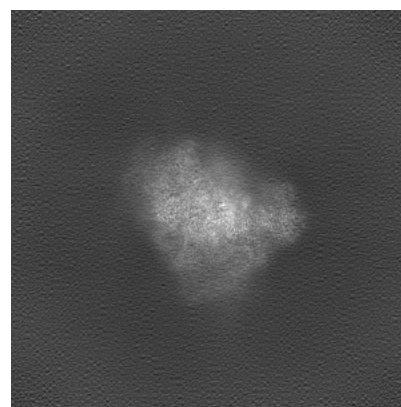
6.1.2 Raw map



X



Y

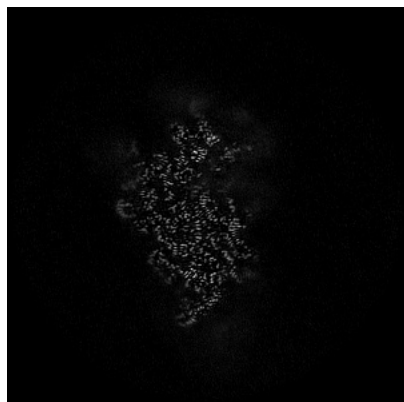


Z

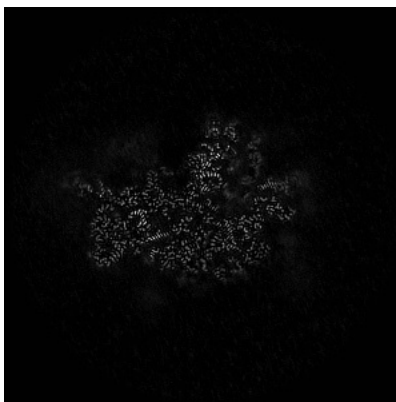
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

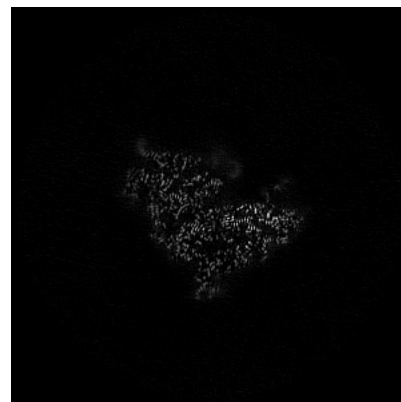
6.2.1 Primary map



X Index: 240

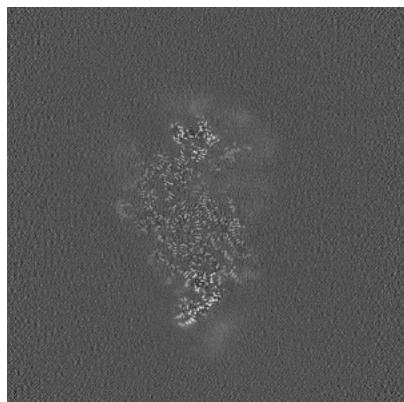


Y Index: 240

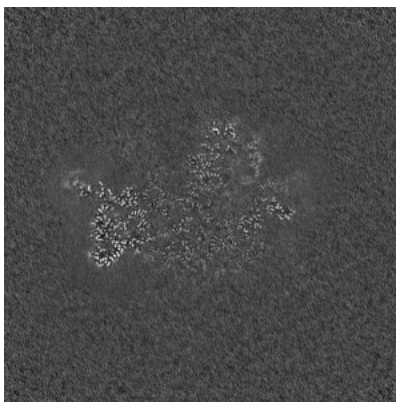


Z Index: 240

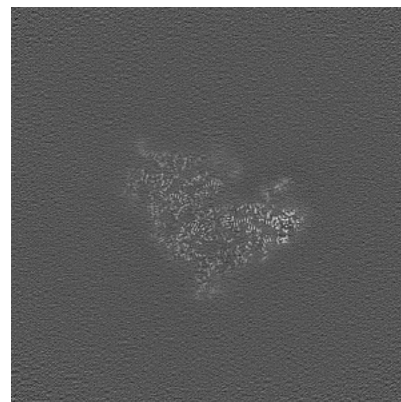
6.2.2 Raw map



X Index: 240



Y Index: 240

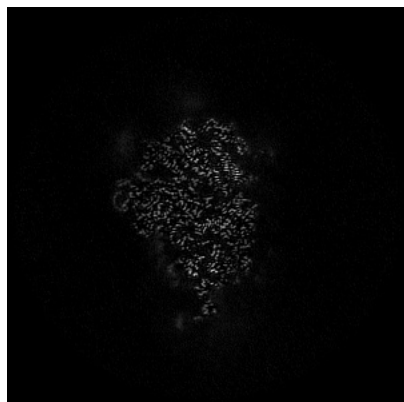


Z Index: 240

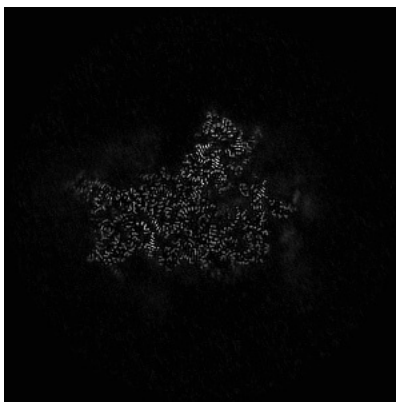
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

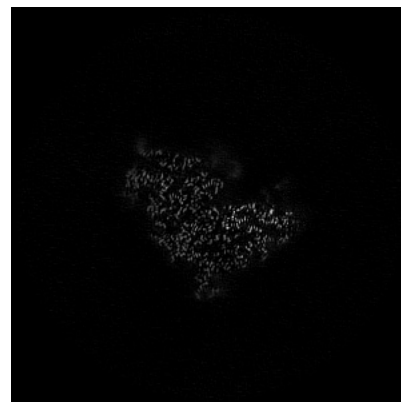
6.3.1 Primary map



X Index: 226

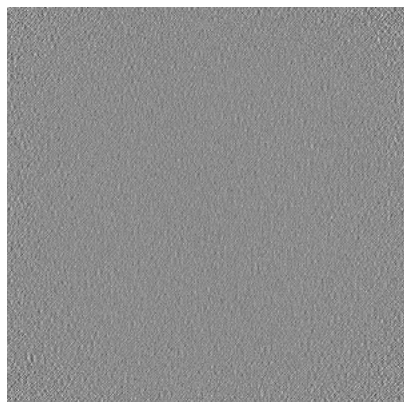


Y Index: 233

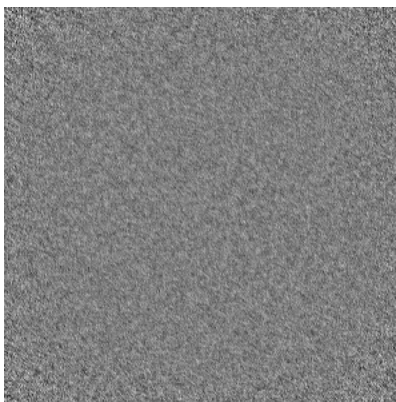


Z Index: 239

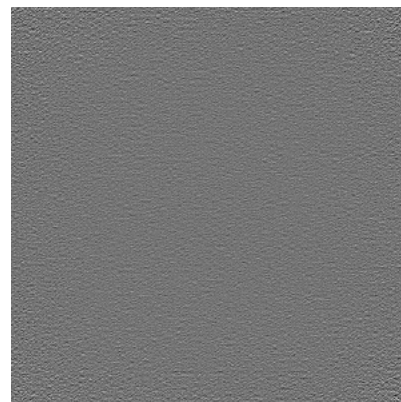
6.3.2 Raw map



X Index: 0



Y Index: 0

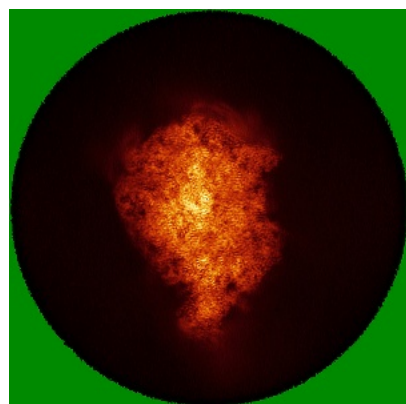


Z Index: 0

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

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

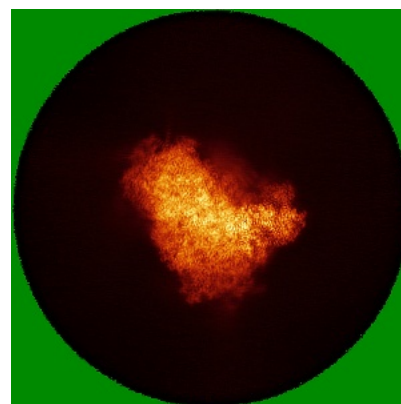
6.4.1 Primary map



X

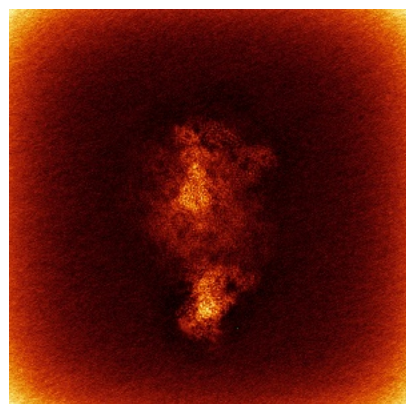


Y

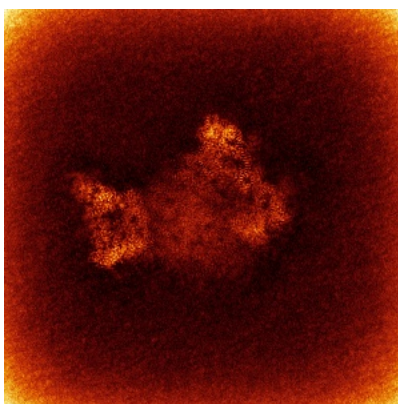


Z

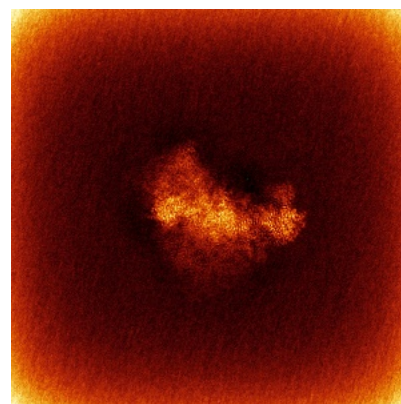
6.4.2 Raw map



X



Y

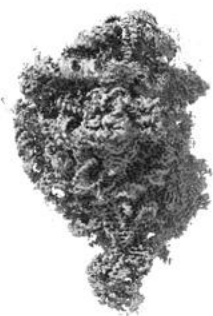


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



X



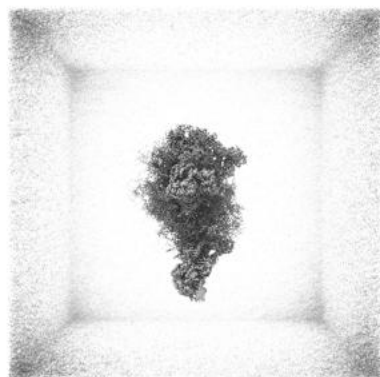
Y



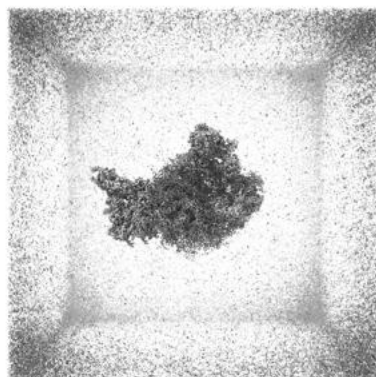
Z

The images above show the 3D surface view of the map at the recommended contour level 0.95. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

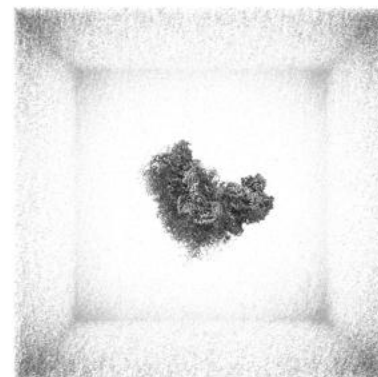
6.5.2 Raw map



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.

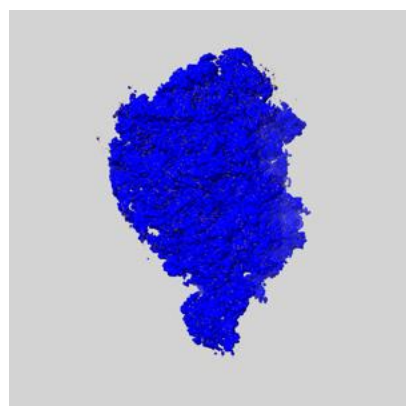
6.6 Mask visualisation [i](#)

This section shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

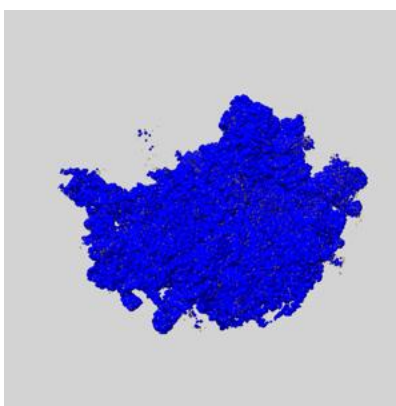
A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

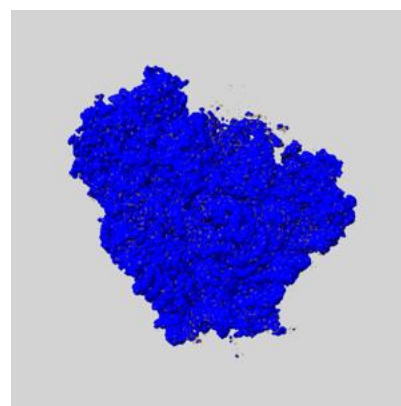
6.6.1 emd_29275_msk_1.map [i](#)



X



Y

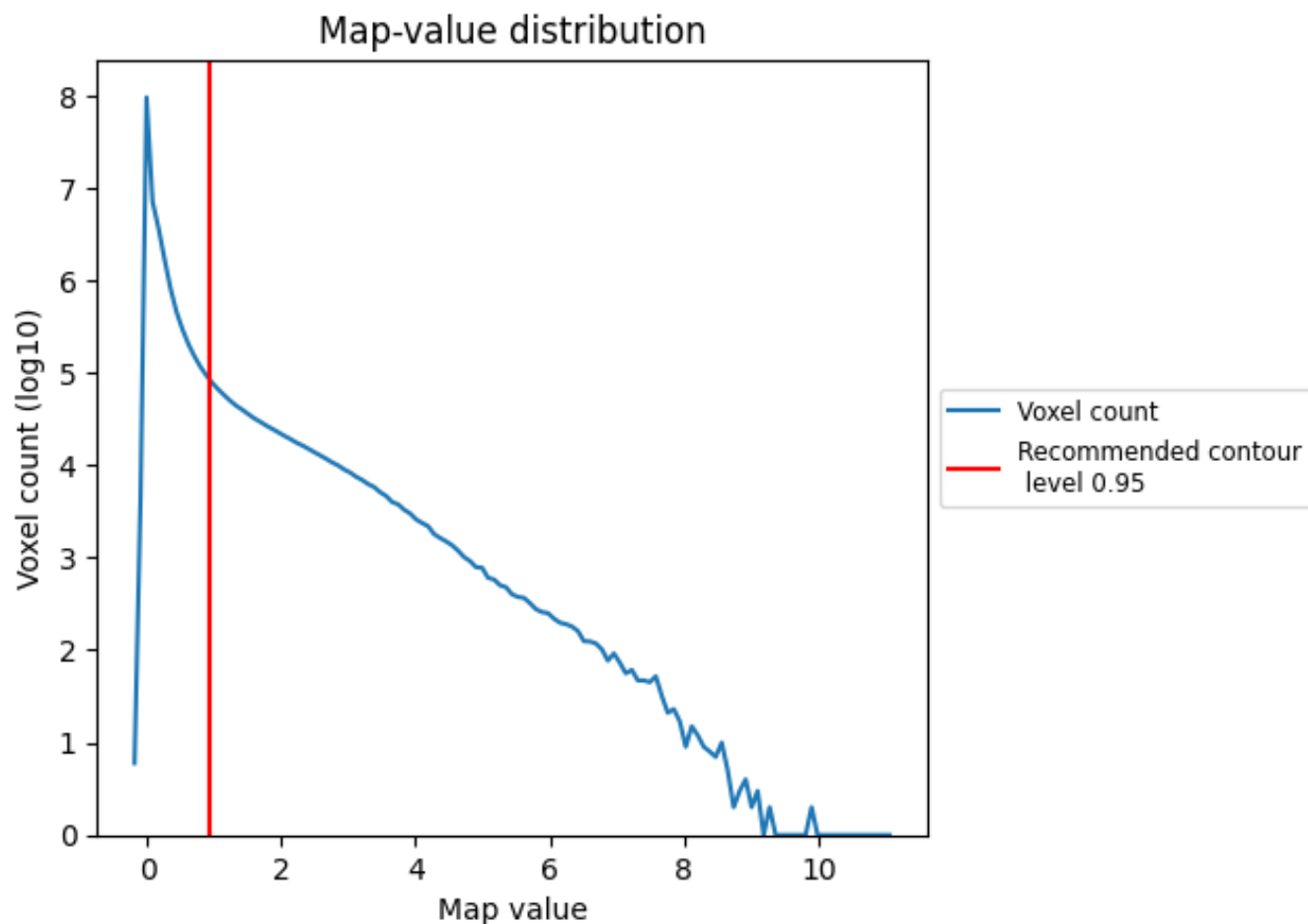


Z

7 Map analysis [i](#)

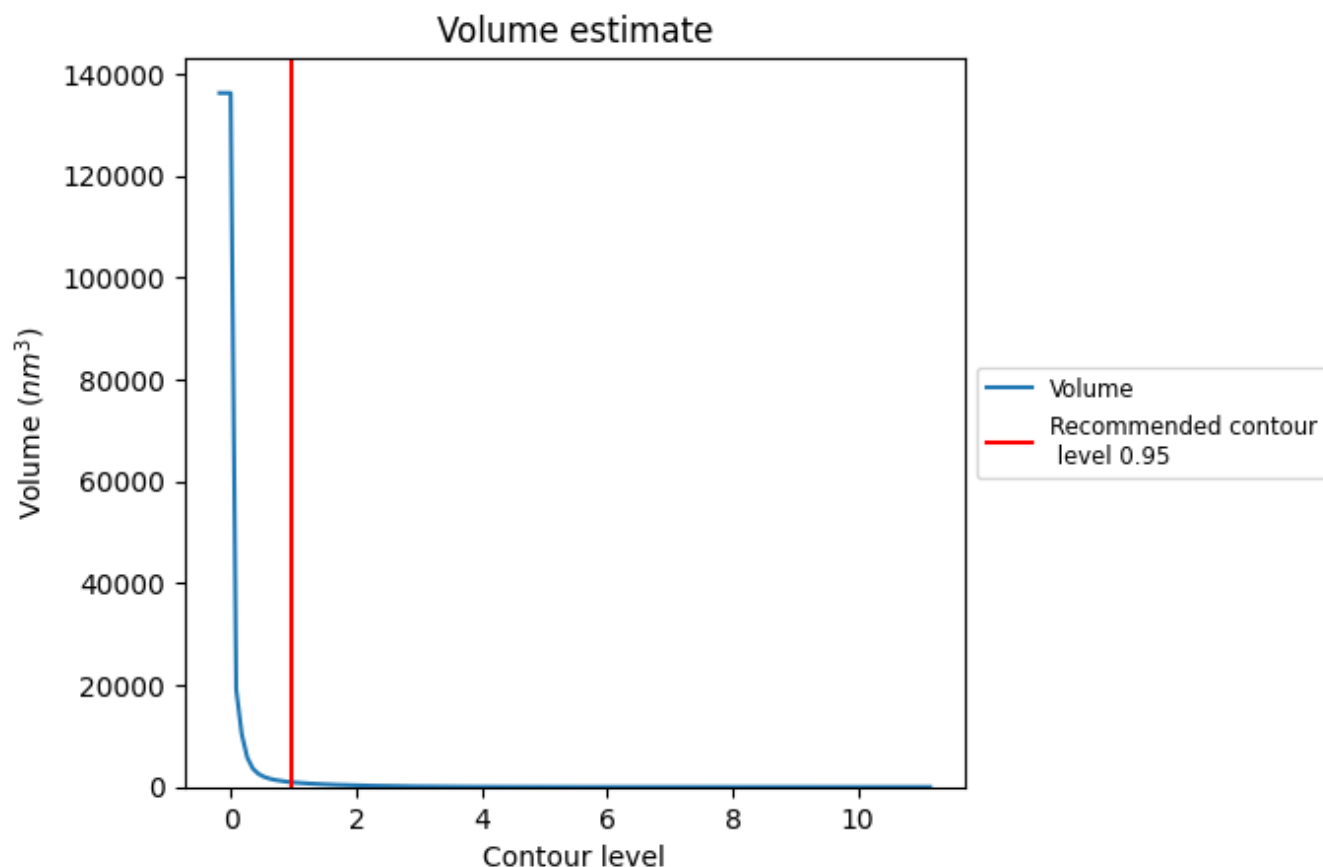
This section contains the results of statistical analysis of the map.

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.

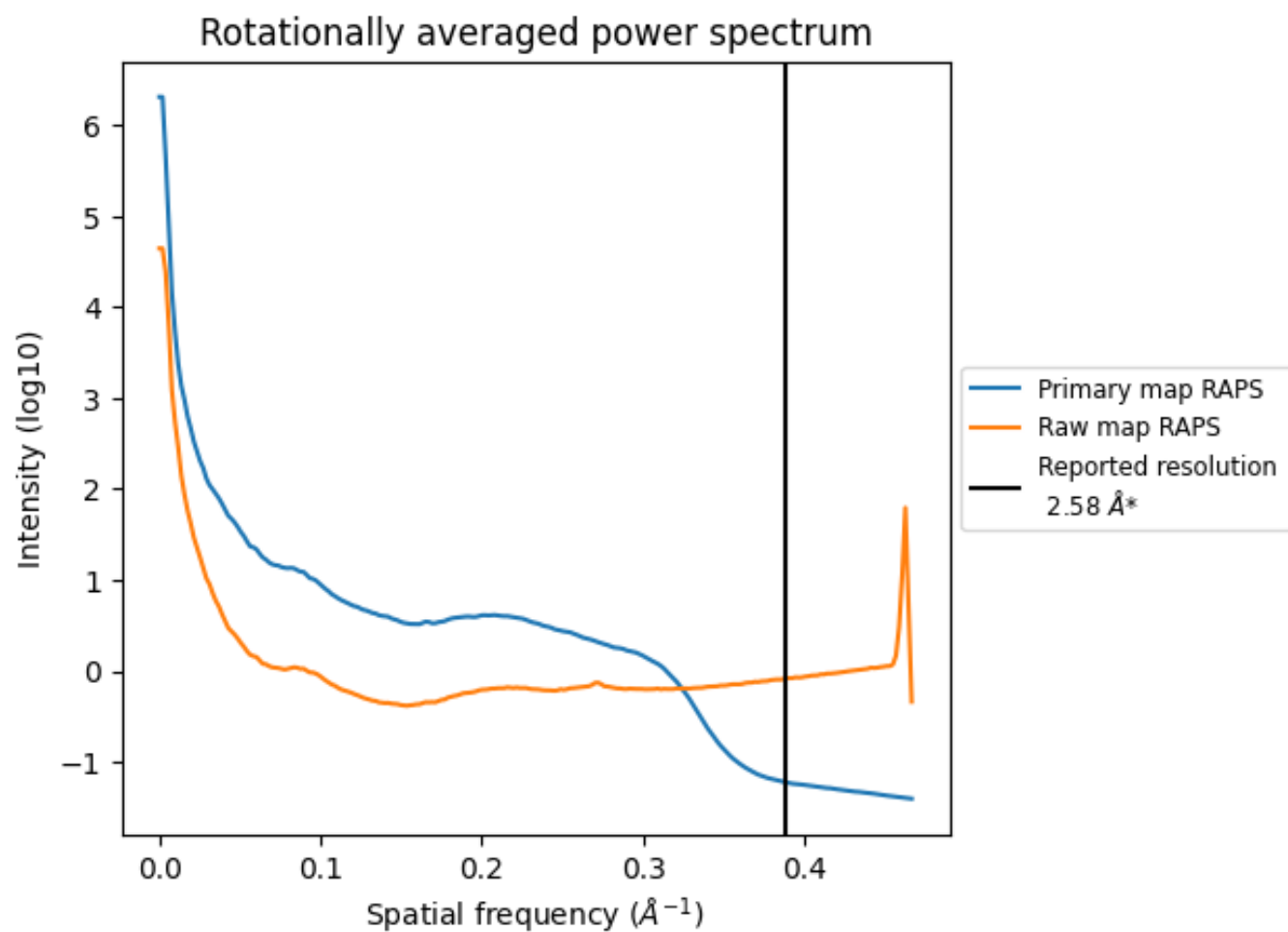
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 953 nm^3 ; this corresponds to an approximate mass of 861 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 ⓘ

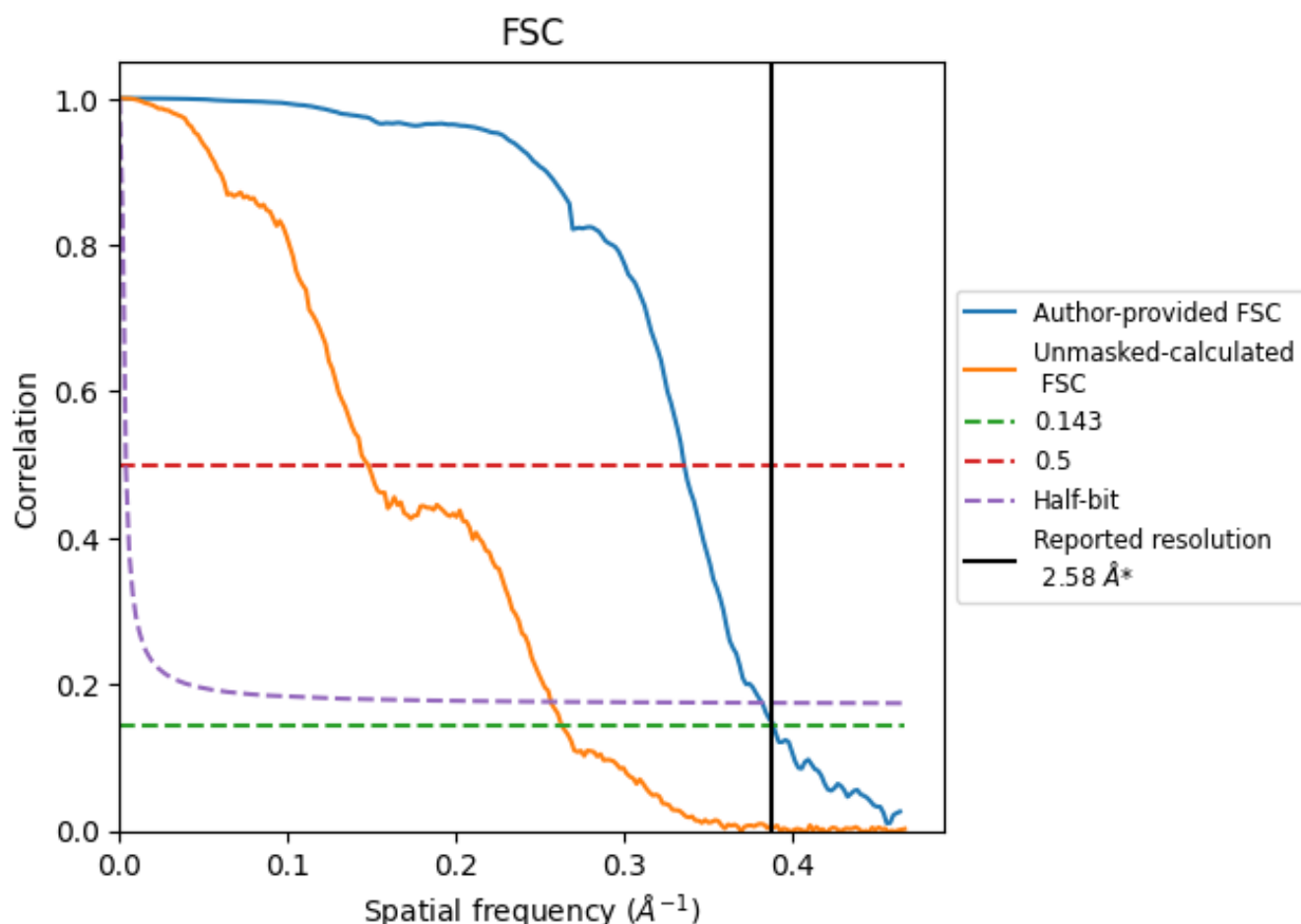


*Reported resolution corresponds to spatial frequency of 0.388 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.388 Å⁻¹

8.2 Resolution estimates [i](#)

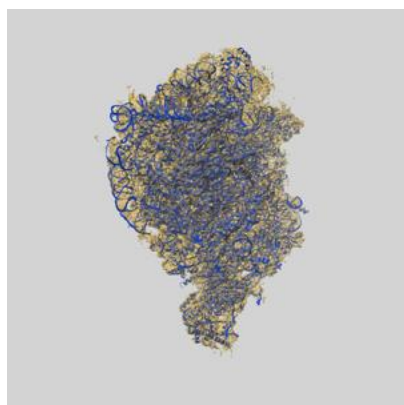
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.58	-	-
Author-provided FSC curve	2.58	2.98	2.62
Unmasked-calculated*	3.81	6.78	3.90

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.81 differs from the reported value 2.58 by more than 10 %

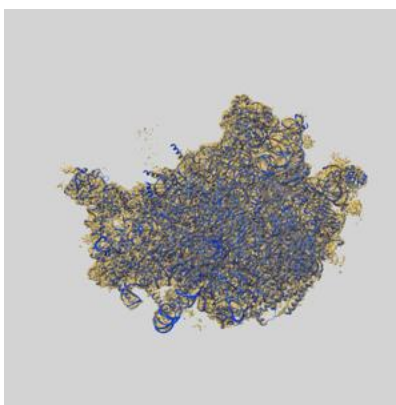
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-29275 and PDB model 8FLD. Per-residue inclusion information can be found in [section 3](#) on [page 14](#).

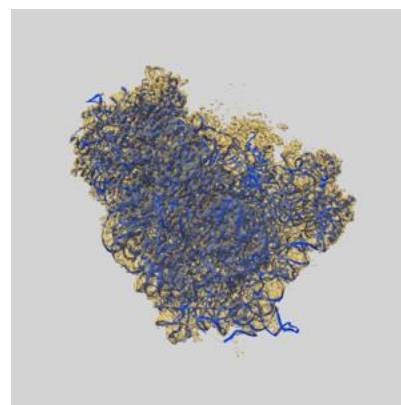
9.1 Map-model overlay [i](#)



X



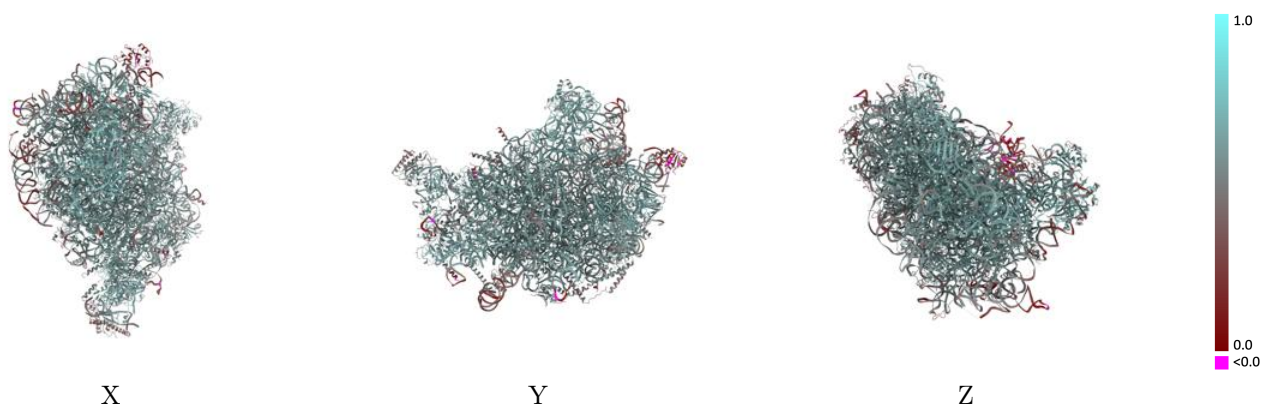
Y



Z

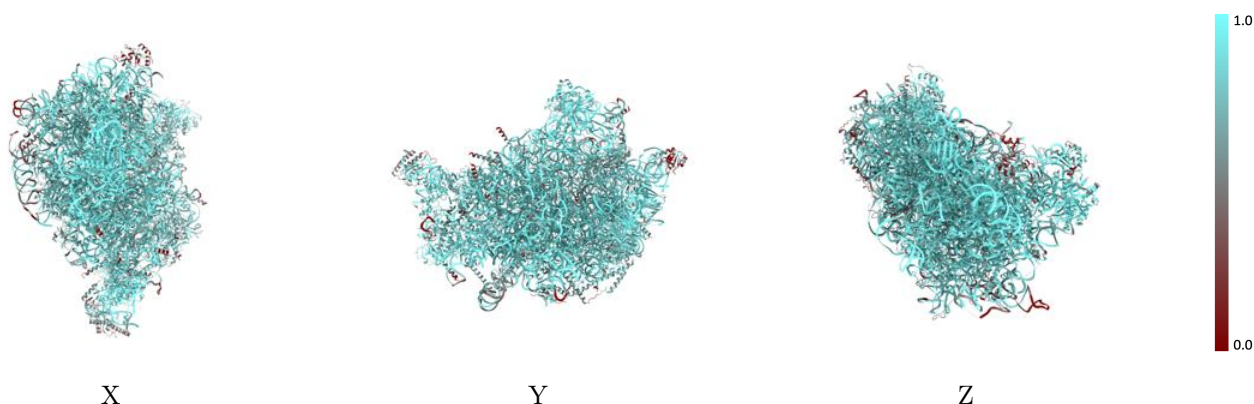
The images above show the 3D surface view of the map at the recommended contour level 0.95 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



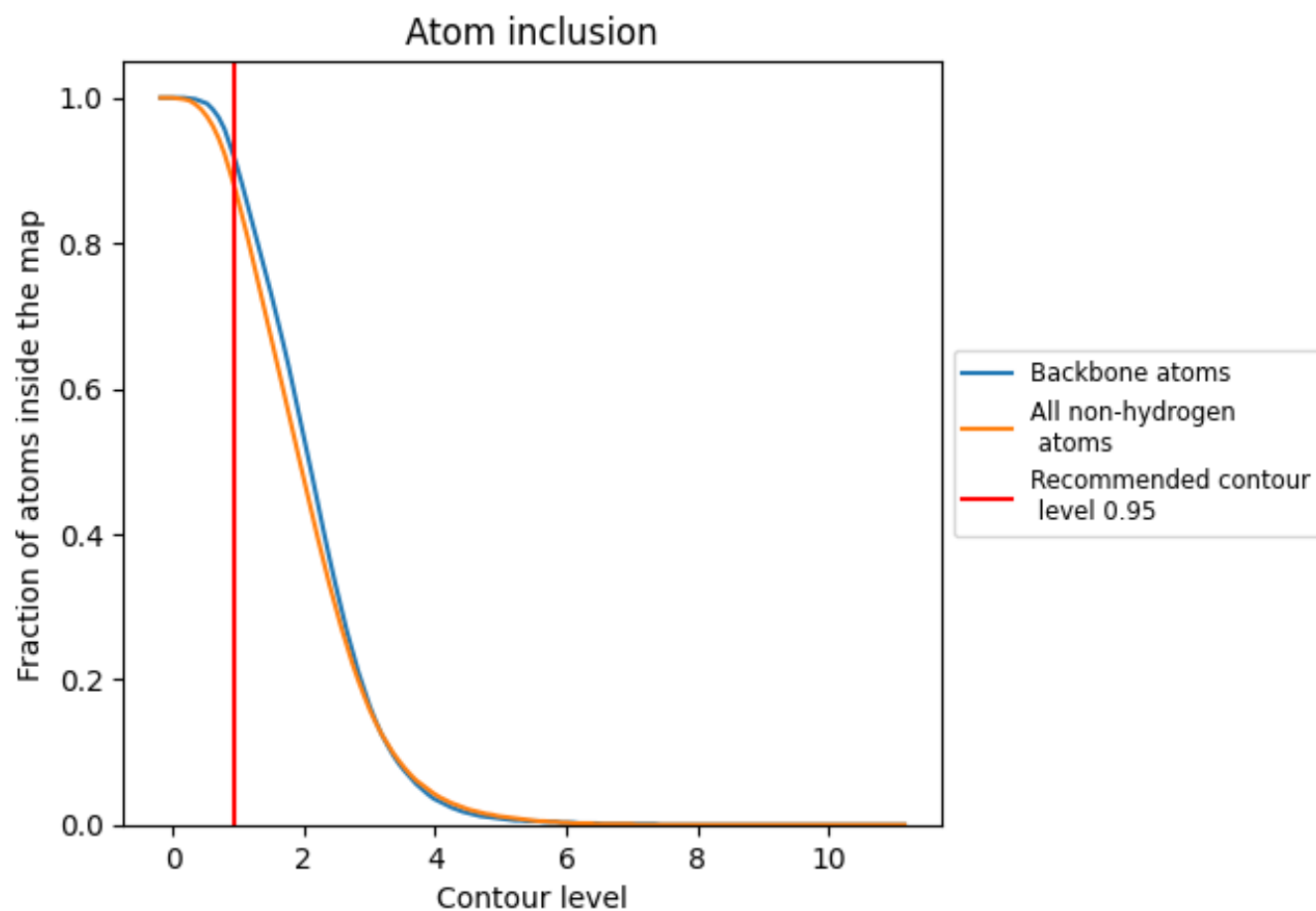
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.95).







































































9.4 Atom inclusion [i](#)



At the recommended contour level, 92% 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 (0.95) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8760	 0.5720
BE	 0.6410	 0.4570
L1	 0.9420	 0.6080
L2	 0.8700	 0.5540
L3	 0.9110	 0.5640
L4	 0.9910	 0.6370
L5	 0.8250	 0.5800
L6	 0.8080	 0.5640
L7	 0.9000	 0.6050
L8	 0.9130	 0.6070
L9	 0.9420	 0.6350
LA	 0.8630	 0.5860
LB	 0.8960	 0.6080
LC	 0.9620	 0.6470
LD	 0.8230	 0.5630
LE	 0.8980	 0.6140
LF	 0.7520	 0.5420
LG	 0.8620	 0.5980
LH	 0.8620	 0.6040
LI	 0.8170	 0.5720
LJ	 0.8760	 0.6020
LK	 0.9000	 0.6080
LL	 0.8610	 0.5790
LM	 0.7480	 0.5520
LN	 0.8560	 0.5850
LO	 0.7540	 0.5590
LP	 0.7990	 0.5580
LQ	 0.8660	 0.5940
LR	 0.8520	 0.5920
LS	 0.8210	 0.5810
LT	 0.8940	 0.6050
LU	 0.7710	 0.5530
LV	 0.8930	 0.6130
LW	 0.9170	 0.6140
LX	 0.7840	 0.5750



Continued on next page...

Continued from previous page...

Chain	Atom inclusion	Q-score
LY	 0.7270	 0.5350
LZ	 0.9080	 0.6090
NK	 0.6080	 0.5020
NL	 0.7570	 0.5540
NP	 0.7390	 0.5400
SA	 0.8730	 0.5920
SB	 0.8850	 0.6090
SC	 0.7680	 0.5410
SD	 0.8640	 0.5890
SE	 0.8690	 0.5980
SF	 0.8910	 0.6130
SG	 0.8860	 0.6030
SH	 0.7460	 0.5310
SI	 0.8450	 0.5860
SK	 0.8450	 0.5760
SL	 0.6710	 0.4970
SM	 0.9030	 0.6260
SQ	 0.3740	 0.2280
SR	 0.7000	 0.5120
SV	 0.7920	 0.5420