



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 01:46 PM UTC

PDB ID : 5A8L / pdb_00005a8l
EMDB ID : EMD-3099
Title : Human eRF1 and the hCMV nascent peptide in the translation termination complex
Authors : Matheisl, S.; Berninghausen, O.; Becker, T.; Beckmann, R.
Deposited on : 2015-07-16
Resolution : 3.80 Å(reported)
Based on initial model : 3E1Y

This is a Full wwPDB EM Validation Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

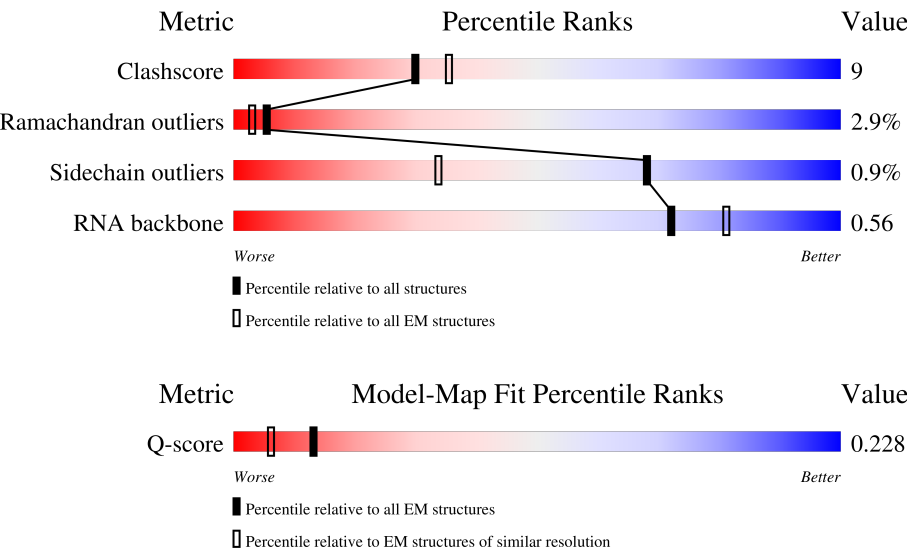
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance i

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

The reported resolution of this entry is 3.80 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	10198 (3.30 - 4.30)

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

Mol	Chain	Length	Quality of chain
1	A	5025	95%
2	B	1869	96%
3	D	184	5% 9% 89%

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Mol	Chain	Length	Quality of chain
4	G	165	79% 62% 19% 16%
5	H	427	97%
6	P	18	6% 72% 22% 6%
7	Q	431	83% 72% 27%
8	R	9	33% 56% 33% 11%
9	Z	22	77% 14% 50% 27% 9%

2 Entry composition

There are 9 unique types of molecules in this entry. The entry contains 12264 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 28S RIBOSOMAL RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	253	Total	C	N	O	P	0	0
			5401	2413	971	1764	253		

- Molecule 2 is a RNA chain called HUMAN 18S RIBOSOMAL RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	67	Total	C	N	O	P	0	0
			1423	636	252	468	67		

- Molecule 3 is a protein called 60S RIBOSOMAL PROTEIN L17.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	D	21	Total	C	N	O	S	0	0
			179	110	40	27	2		

- Molecule 4 is a protein called 60S RIBOSOMAL PROTEIN L12.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	G	138	Total	C	N	O	S	0	0
			1046	654	196	193	3		

- Molecule 5 is a protein called 60S RIBOSOMAL PROTEIN L4.

Mol	Chain	Residues	Atoms				AltConf	Trace
5	H	12	Total	C	N	O	0	0
			88	49	24	15		

- Molecule 6 is a RNA chain called P-SITE TRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	P	18	Total	C	N	O	P	0	0
			382	172	71	122	17		

- Molecule 7 is a protein called EUKARYOTIC RELEASE FACTOR ERF1.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	Q	431	Total	C	N	O	S	0	0
			3410	2168	575	655	12		

- Molecule 8 is a RNA chain called MRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	R	9	Total	C	N	O	P	0	0
			190	86	35	60	9		

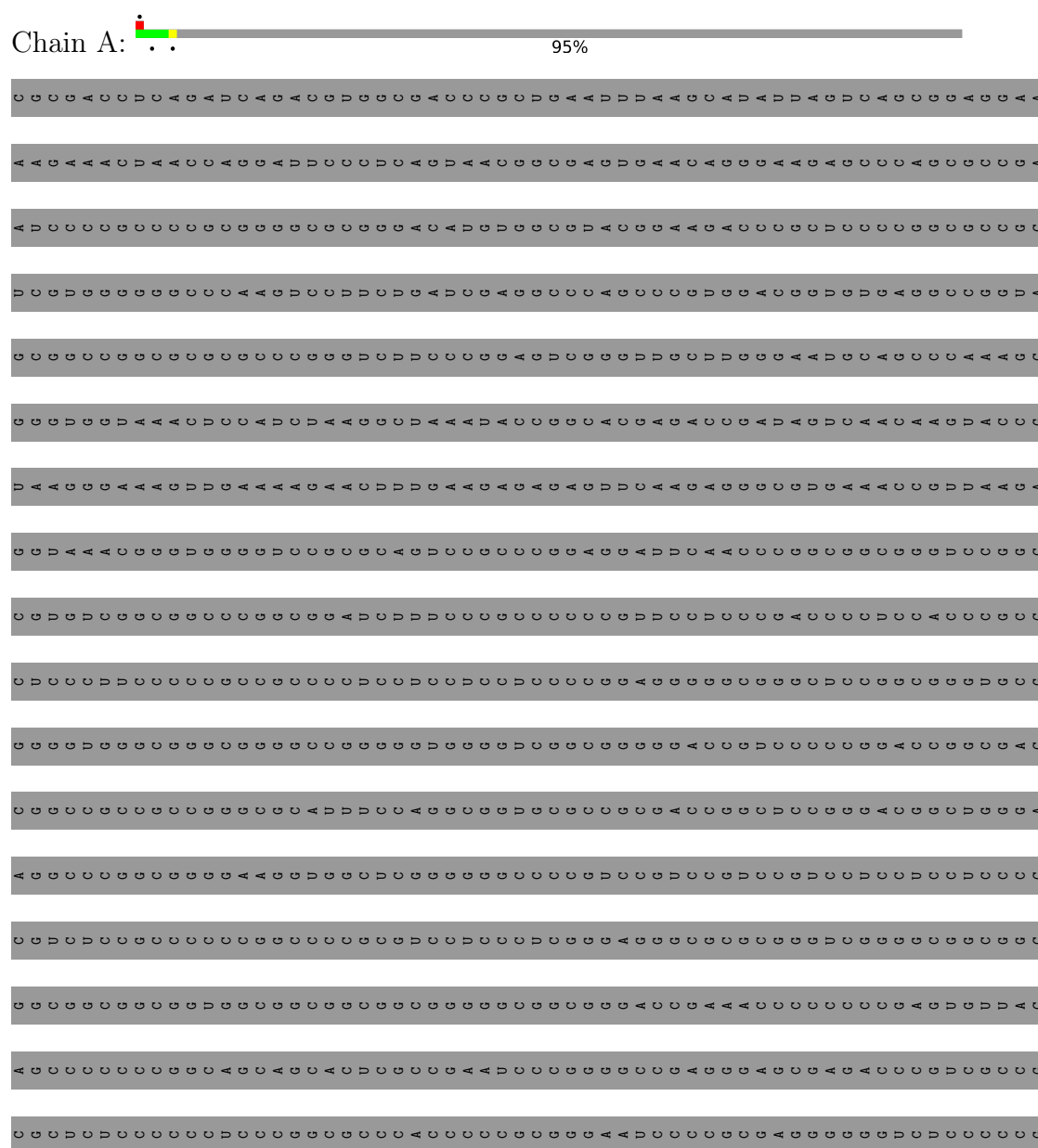
- Molecule 9 is a protein called NASCENT CHAIN.

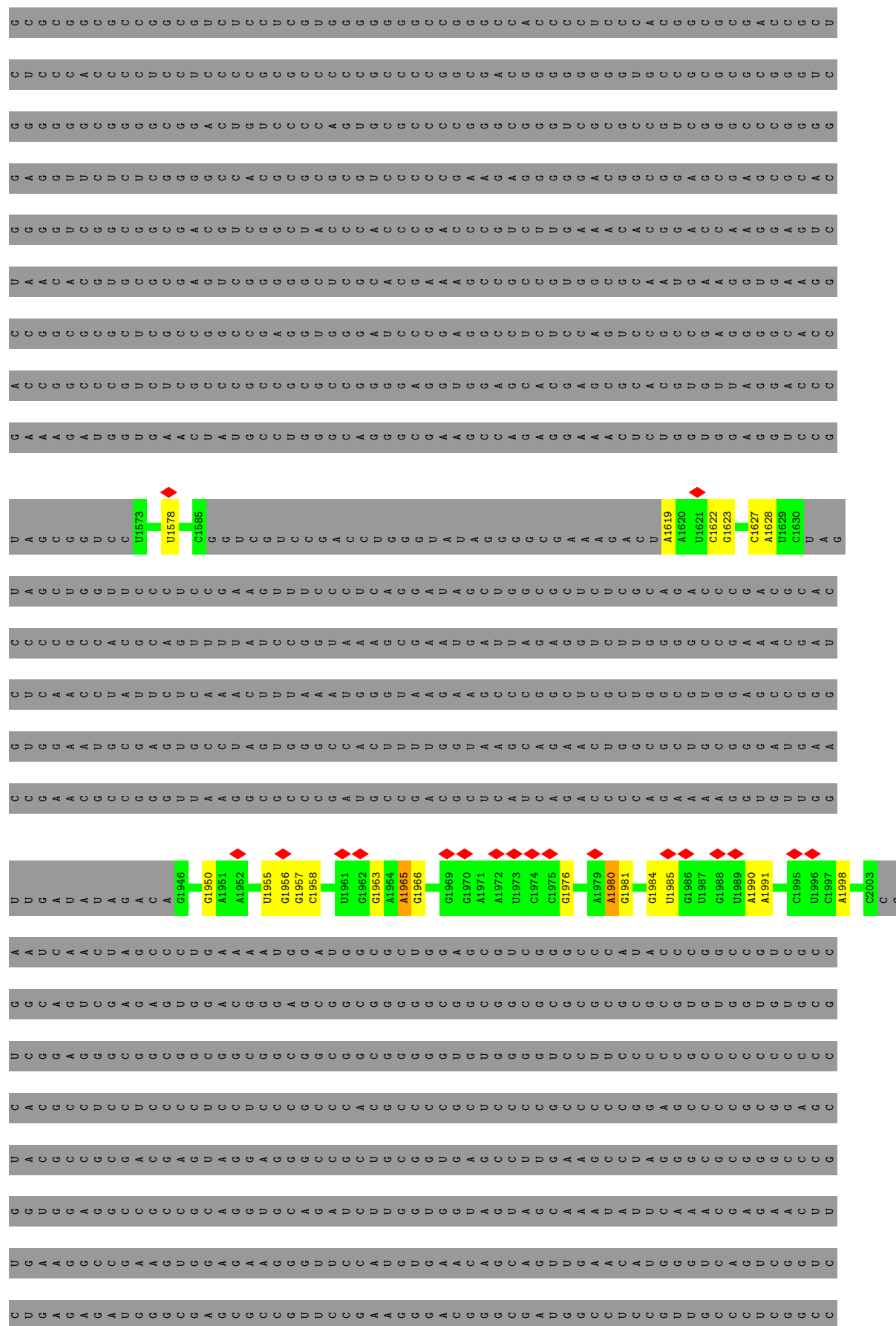
Mol	Chain	Residues	Atoms					AltConf	Trace
9	Z	22	Total	C	N	O	S	0	0
			145	92	25	27	1		

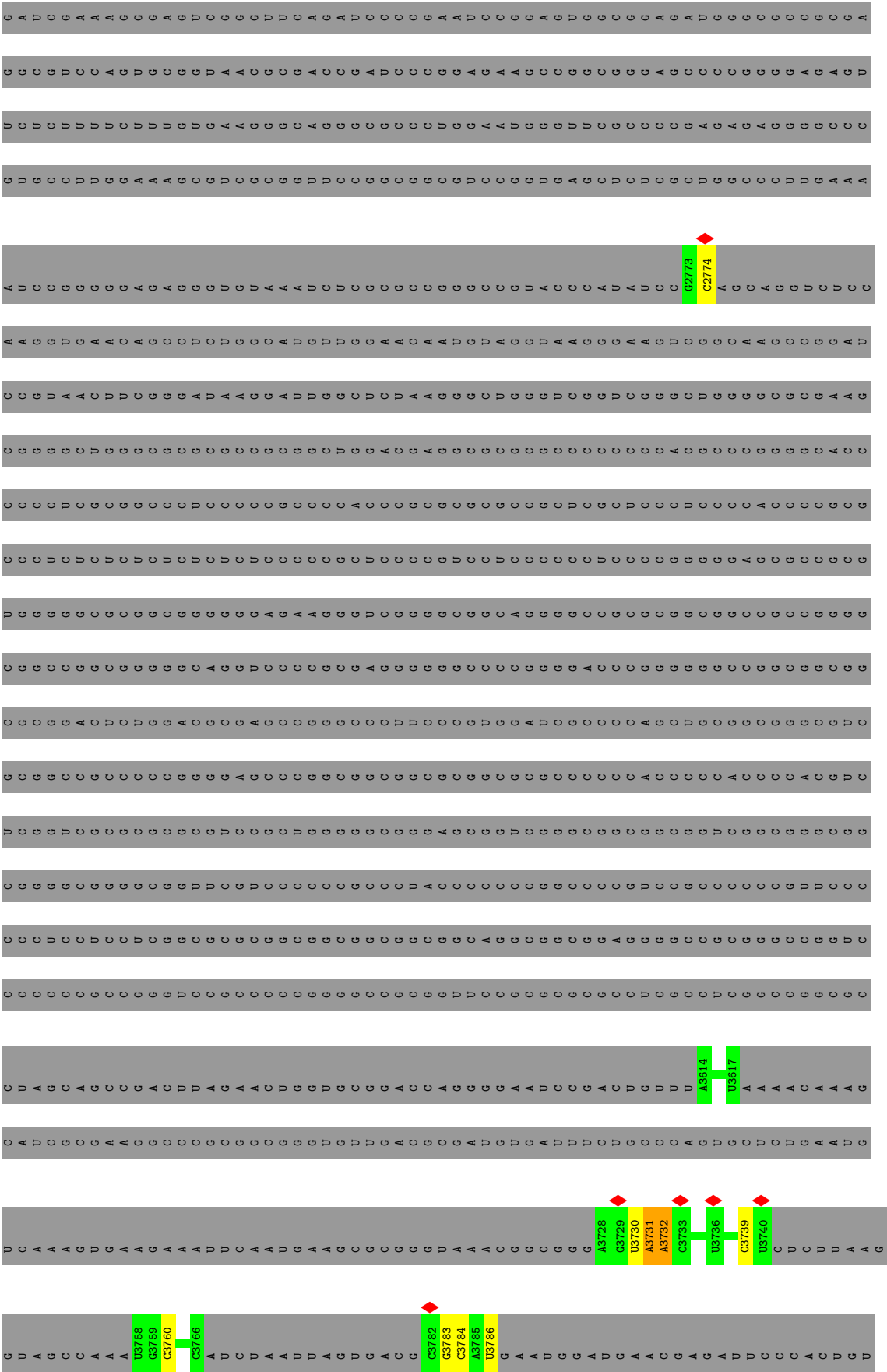
3 Residue-property plots

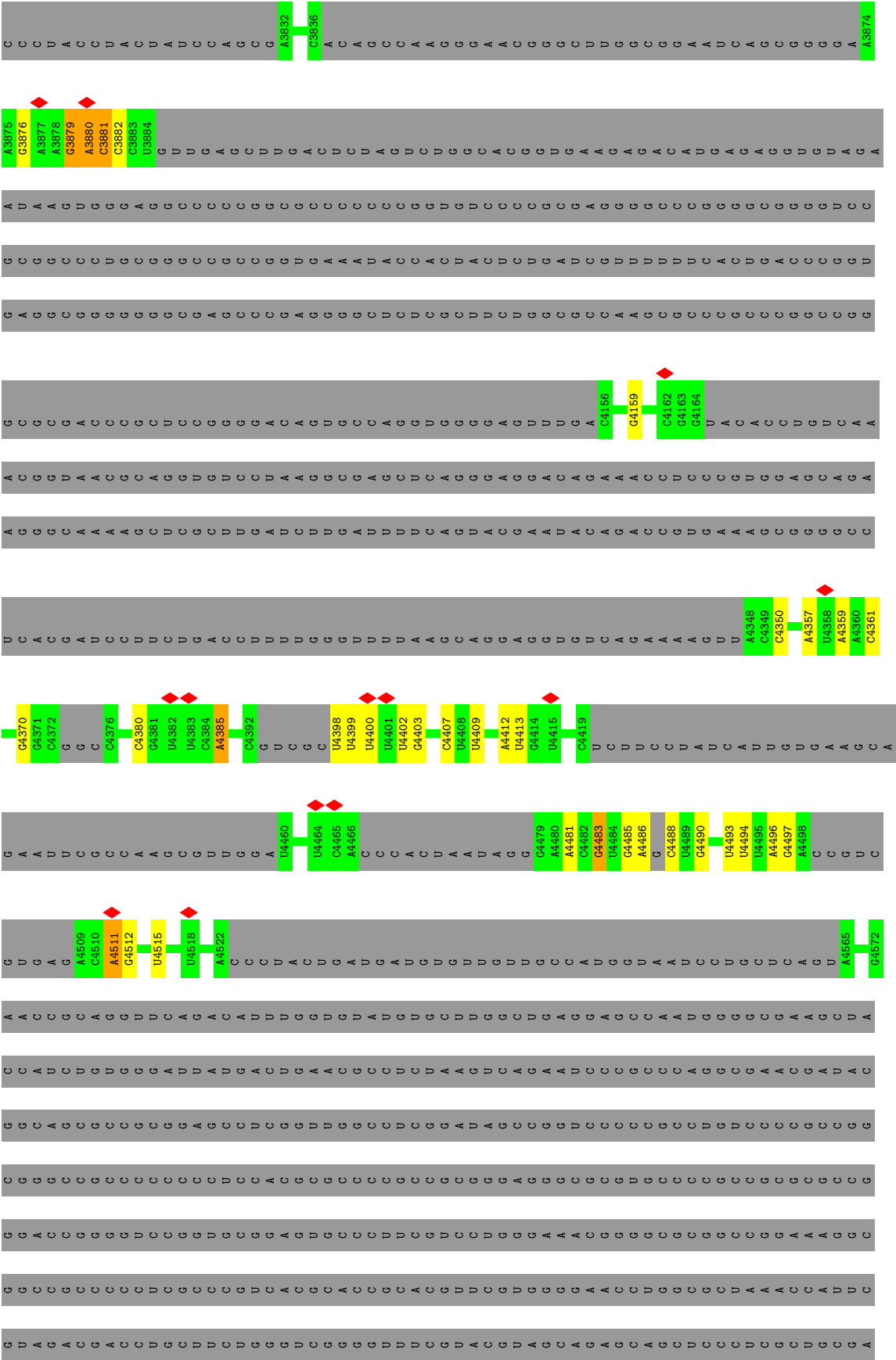
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

• Molecule 1: 28S RIBOSOMAL RNA



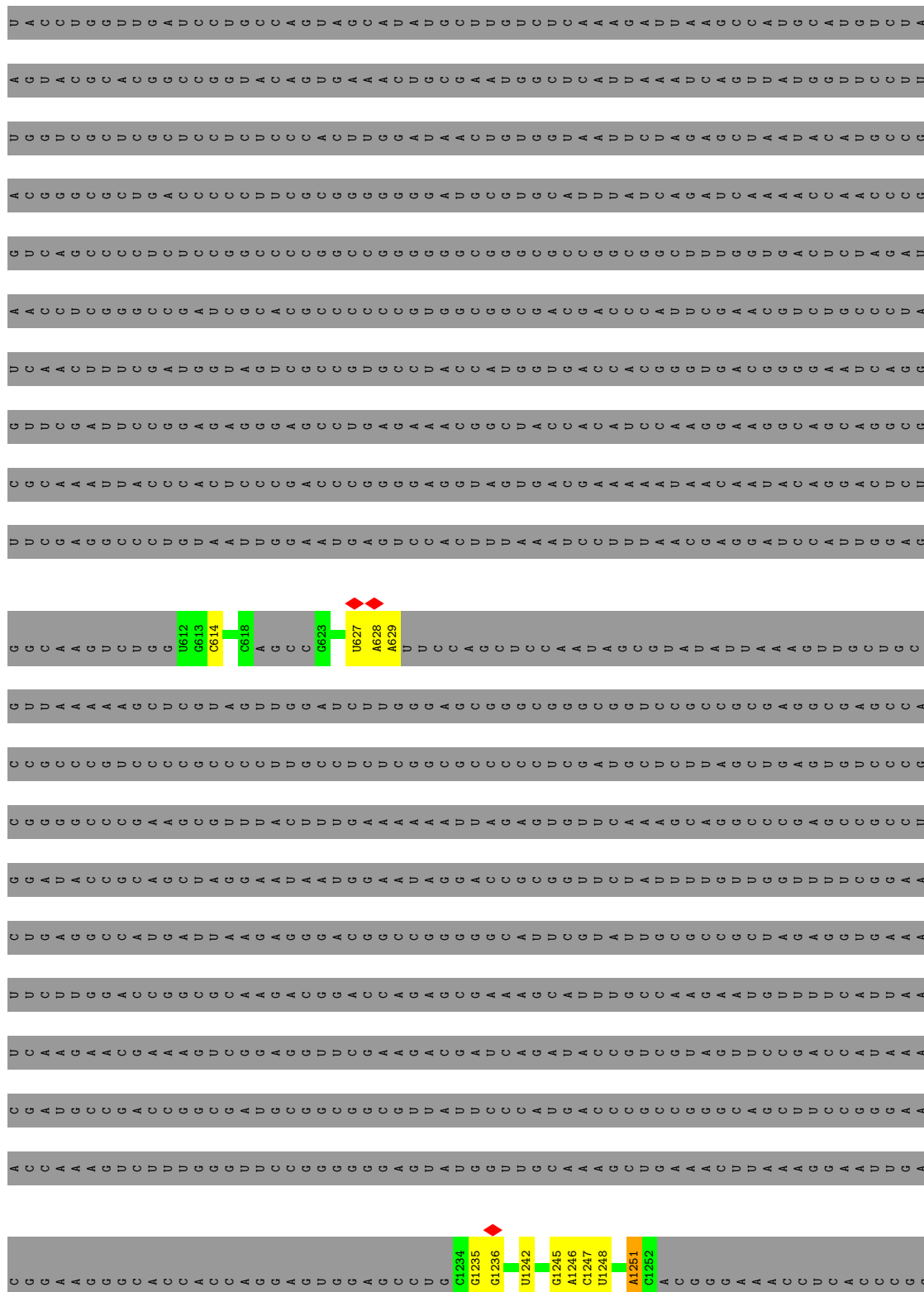


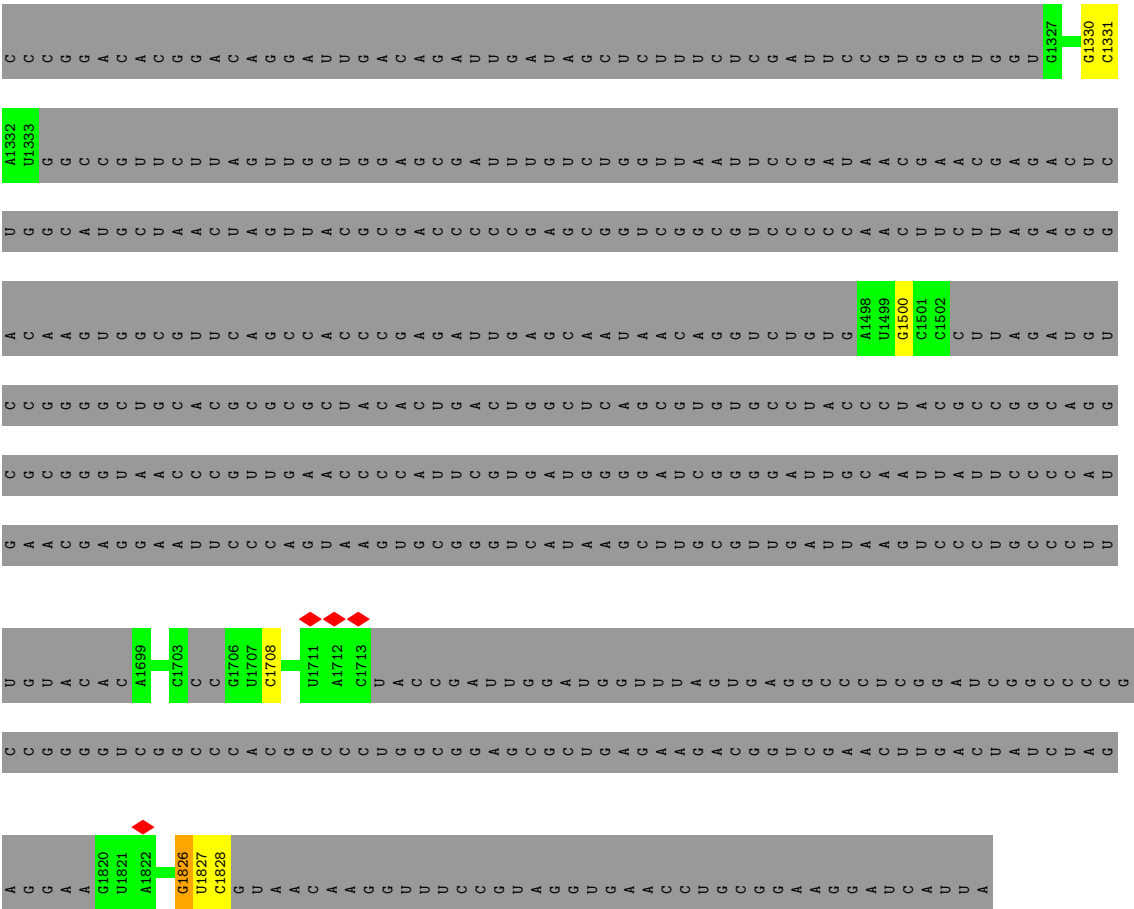




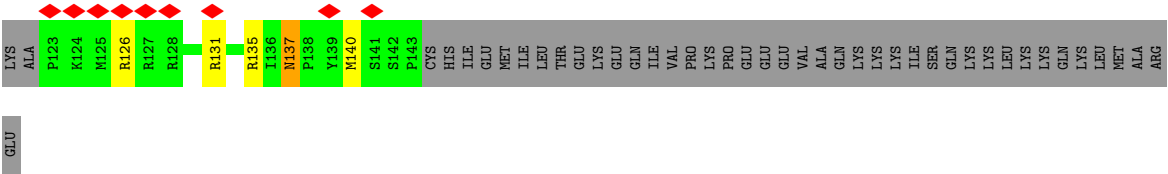
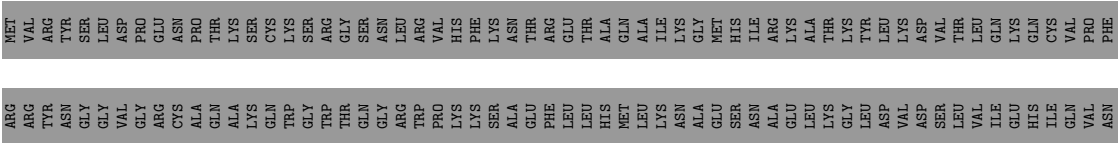
- Molecule 2: HUMAN 18S RIBOSOMAL RNA

Chain B: 96%

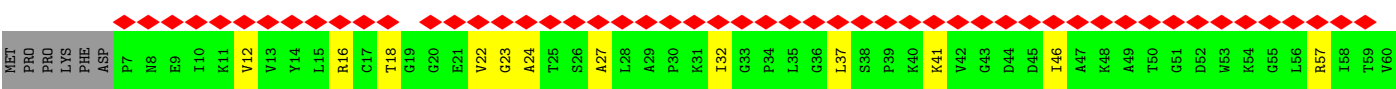
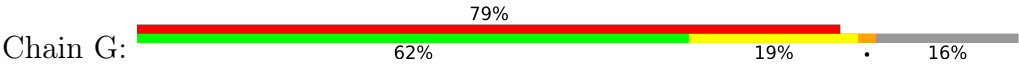




● Molecule 3: 60S RIBOSOMAL PROTEIN L17

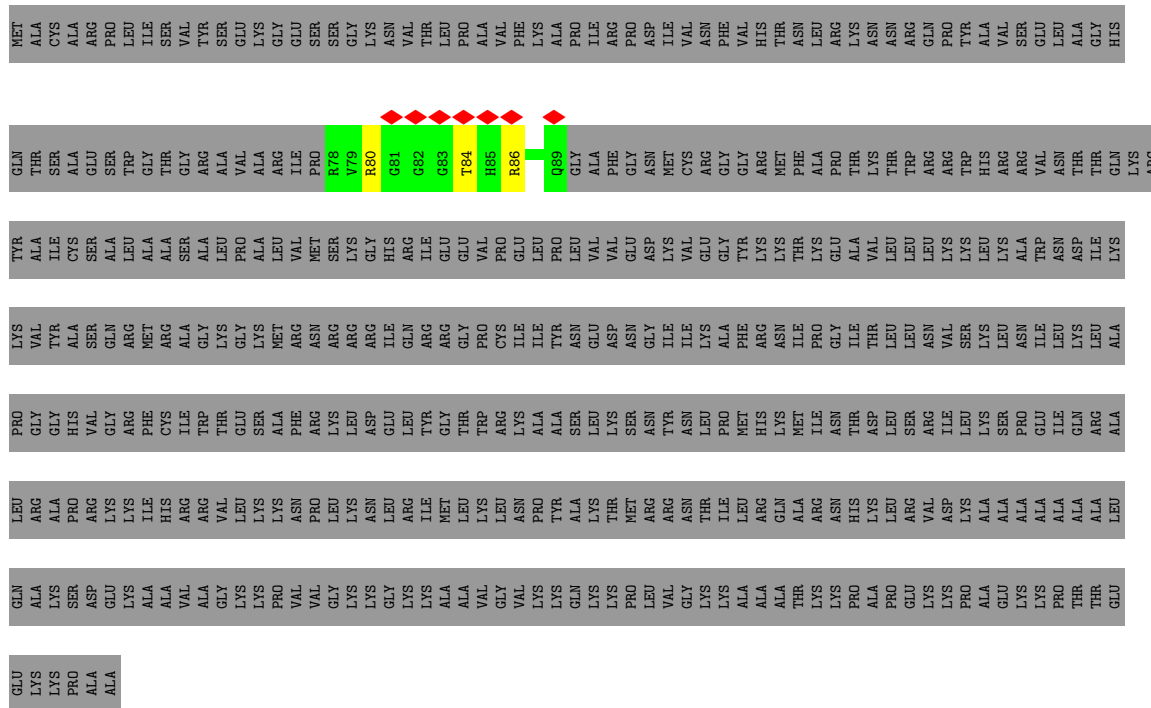


● Molecule 4: 60S RIBOSOMAL PROTEIN L12



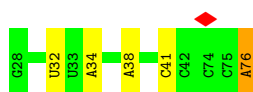
- Molecule 5: 60S RIBOSOMAL PROTEIN L4

Chain H: 97%

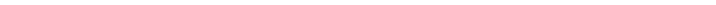


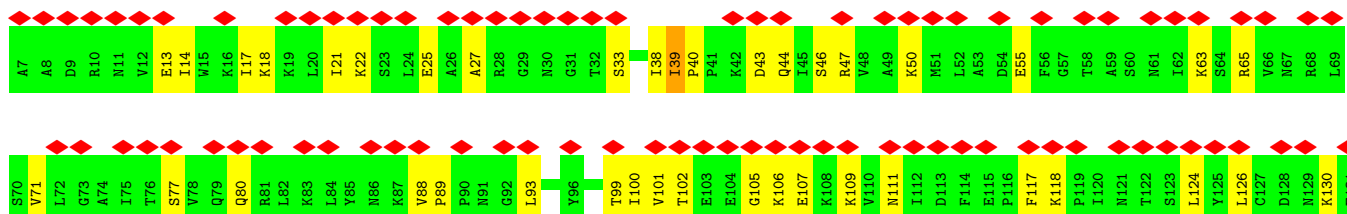
- Molecule 6: P-SITE tRNA

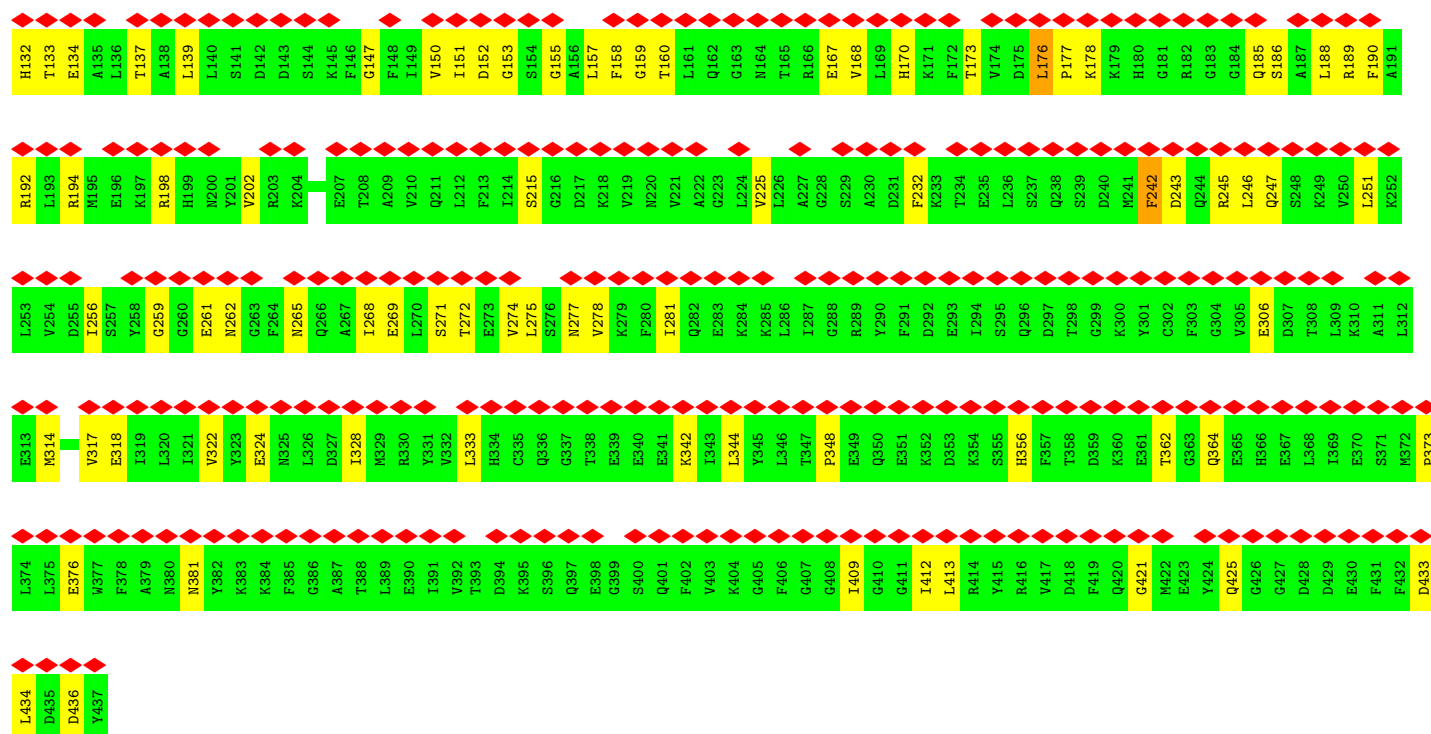
Chain P: 6% 72% 22% 6%



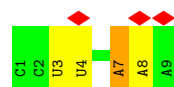
- Molecule 7: EUKARYOTIC RELEASE FACTOR ERF1

Chain Q:  83% 72% 27%

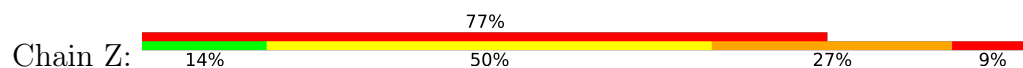




- Molecule 8: MRNA



- Molecule 9: NASCENT CHAIN



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	33165	Depositor
Resolution determination method	Not provided	
CTF correction method	DEFOCUS GROUPS	Depositor
Microscope	FEI TECNAI F20	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	Not provided	
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2700	Depositor
Magnification	Not provided	
Image detector	FEI FALCON II (4k x 4k)	Depositor
Maximum map value	0.003	Depositor
Minimum map value	-0.002	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.000	Depositor
Recommended contour level	0.0008	Depositor
Map size ($\text{\AA}$)	446.04, 446.04, 446.04	wwPDB
Map dimensions	420, 420, 420	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.062, 1.062, 1.062	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# $ Z > 5$	RMSZ	# $ Z > 5$
1	A	0.05	0/6024	0.11	0/9344
2	B	0.04	0/1582	0.08	0/2445
3	D	0.60	0/184	0.92	2/244 (0.8%)
4	G	0.54	1/1058 (0.1%)	0.86	2/1424 (0.1%)
5	H	0.26	0/88	0.56	0/114
6	P	0.04	0/426	0.07	0/660
7	Q	0.40	2/3467 (0.1%)	0.76	8/4661 (0.2%)
8	R	0.06	0/212	0.12	0/327
9	Z	0.91	1/147 (0.7%)	1.42	3/195 (1.5%)
All	All	0.28	4/13188 (0.0%)	0.48	15/19414 (0.1%)

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	Q	40	PRO	CA-C	8.49	1.56	1.51
7	Q	89	PRO	CA-C	7.44	1.55	1.51
4	G	88	PRO	CA-C	7.28	1.58	1.52
9	Z	21	PRO	CA-C	6.05	1.57	1.52

All (15) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
7	Q	88	VAL	CA-C-N	9.81	126.82	119.66
7	Q	88	VAL	C-N-CA	9.81	126.82	119.66
7	Q	39	ILE	CA-C-N	9.41	126.53	119.66
7	Q	39	ILE	C-N-CA	9.41	126.53	119.66
4	G	74	VAL	CA-C-N	7.62	127.63	119.78
4	G	74	VAL	C-N-CA	7.62	127.63	119.78
7	Q	176	LEU	CA-C-N	6.44	126.36	119.28
7	Q	176	LEU	C-N-CA	6.44	126.36	119.28
7	Q	40	PRO	CA-C-N	5.94	125.89	119.78
7	Q	40	PRO	C-N-CA	5.94	125.89	119.78
9	Z	15	LEU	N-CA-C	-5.86	104.23	111.33

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
9	Z	16	THR	N-CA-C	-5.51	105.27	111.28
3	D	137	ASN	CA-C-N	5.11	125.04	119.78
3	D	137	ASN	C-N-CA	5.11	125.04	119.78
9	Z	20	ILE	CB-CA-C	-5.06	107.37	114.00

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	A	5401	0	2742	48	0
2	B	1423	0	730	9	0
3	D	179	0	184	4	0
4	G	1046	0	1116	18	0
5	H	88	0	86	3	0
6	P	382	0	197	3	0
7	Q	3410	0	3419	85	0
8	R	190	0	98	3	0
9	Z	145	0	145	52	0
All	All	12264	0	8717	184	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 9.

All (184) 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)
9:Z:15:LEU:CA	9:Z:18:LYS:HE3	1.62	1.29
9:Z:8:ALA:O	9:Z:12:SER:HB3	1.22	1.26
9:Z:15:LEU:O	9:Z:18:LYS:HD2	1.23	1.26
1:A:3880:A:H2'	9:Z:19:TYR:CE2	1.82	1.14
9:Z:9:LYS:HD2	9:Z:11:LEU:HB2	1.15	1.14
9:Z:15:LEU:HA	9:Z:18:LYS:CE	1.81	1.11
1:A:3880:A:H2'	9:Z:19:TYR:CZ	1.89	1.08

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:Z:15:LEU:O	9:Z:18:LYS:CD	2.04	1.04
9:Z:9:LYS:HD3	9:Z:11:LEU:H	1.20	1.03
1:A:4511:A:C2	7:Q:190:PHE:CE1	2.50	1.00
9:Z:15:LEU:CB	9:Z:18:LYS:HE3	1.90	1.00
9:Z:15:LEU:HA	9:Z:18:LYS:HE3	1.32	0.98
9:Z:8:ALA:O	9:Z:12:SER:CB	2.13	0.95
1:A:4511:A:C4	7:Q:190:PHE:CZ	2.58	0.91
9:Z:9:LYS:CD	9:Z:11:LEU:H	1.84	0.89
9:Z:9:LYS:HD2	9:Z:11:LEU:CB	2.01	0.88
9:Z:13:SER:O	9:Z:16:THR:OG1	1.90	0.88
1:A:4511:A:C4	7:Q:190:PHE:CE2	2.62	0.87
9:Z:15:LEU:CA	9:Z:18:LYS:CE	2.43	0.86
9:Z:10:LYS:HA	9:Z:10:LYS:HE2	1.58	0.84
9:Z:15:LEU:HB3	9:Z:18:LYS:HE3	1.60	0.84
9:Z:16:THR:O	9:Z:20:ILE:HG12	1.81	0.79
9:Z:15:LEU:C	9:Z:18:LYS:HE3	2.09	0.77
1:A:3880:A:C2'	9:Z:19:TYR:CZ	2.67	0.77
7:Q:102:THR:N	7:Q:106:LYS:O	2.15	0.77
7:Q:101:VAL:HA	7:Q:107:GLU:HA	1.67	0.77
1:A:4511:A:C5	7:Q:190:PHE:CZ	2.76	0.72
1:A:4511:A:C2	7:Q:190:PHE:CD1	2.77	0.72
9:Z:9:LYS:CD	9:Z:11:LEU:HB2	2.09	0.72
1:A:4380:C:O2'	7:Q:245:ARG:NH2	2.23	0.72
9:Z:10:LYS:HA	9:Z:10:LYS:CE	2.20	0.70
3:D:126:ARG:HA	3:D:140:MET:HG2	1.73	0.69
7:Q:102:THR:O	7:Q:106:LYS:HB3	1.91	0.69
7:Q:176:LEU:HD12	7:Q:177:PRO:HD2	1.74	0.69
1:A:3760:C:O2'	7:Q:178:LYS:NZ	2.26	0.68
7:Q:251:LEU:HG	7:Q:278:VAL:HG11	1.75	0.67
5:H:84:THR:HG22	5:H:86:ARG:H	1.60	0.66
7:Q:159:GLY:HA2	7:Q:168:VAL:HA	1.78	0.65
1:A:4380:C:N4	1:A:4385:A:O2'	2.30	0.65
7:Q:317:VAL:HG12	7:Q:413:LEU:HD23	1.79	0.65
9:Z:12:SER:HA	9:Z:15:LEU:CD1	2.27	0.64
1:A:3731:A:OP2	7:Q:47:ARG:NH1	2.31	0.63
2:B:628:A:N6	2:B:1500:G:O2'	2.32	0.63
9:Z:15:LEU:HB3	9:Z:18:LYS:CE	2.29	0.62
7:Q:362:THR:HG23	7:Q:364:GLN:H	1.63	0.62
1:A:4511:A:C2	7:Q:190:PHE:CZ	2.88	0.62
1:A:1628:A:OP2	5:H:80:ARG:NH2	2.30	0.61
1:A:4511:A:N1	7:Q:190:PHE:CE1	2.68	0.61

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
7:Q:39:ILE:HD12	7:Q:93:LEU:HD23	1.81	0.60
7:Q:225:VAL:HG22	7:Q:251:LEU:HB2	1.83	0.60
7:Q:256:ILE:HD12	7:Q:259:GLY:HA2	1.83	0.60
1:A:1950:G:H21	1:A:1980:A:H8	1.49	0.60
7:Q:33:SER:HA	7:Q:71:VAL:HG22	1.84	0.59
7:Q:271:SER:HA	7:Q:274:VAL:HG22	1.85	0.59
9:Z:8:ALA:O	9:Z:9:LYS:HG3	2.02	0.59
7:Q:322:VAL:HG22	7:Q:409:ILE:HG13	1.84	0.58
4:G:103:ASN:ND2	4:G:141:CYS:SG	2.77	0.58
7:Q:261:GLU:O	7:Q:265:ASN:ND2	2.37	0.58
9:Z:14:LEU:O	9:Z:17:CYS:N	2.37	0.58
6:P:34:A:H61	8:R:3:U:H3	1.50	0.57
9:Z:15:LEU:O	9:Z:18:LYS:CE	2.52	0.57
1:A:1627:C:OP1	5:H:80:ARG:NH1	2.37	0.57
7:Q:318:GLU:N	7:Q:412:ILE:O	2.36	0.57
7:Q:101:VAL:CG2	7:Q:105:GLY:C	2.78	0.57
1:A:3732:A:N6	7:Q:55:GLU:OE2	2.36	0.57
1:A:3882:C:O2'	1:A:4159:G:N2	2.37	0.57
1:A:1619:A:OP1	1:A:1622:C:N4	2.38	0.57
2:B:1331:C:N4	7:Q:63:LYS:O	2.38	0.56
1:A:4511:A:N3	7:Q:190:PHE:CZ	2.73	0.56
1:A:1963:G:N2	1:A:1990:A:O2'	2.39	0.56
1:A:4511:A:N3	7:Q:190:PHE:CE2	2.73	0.56
2:B:1235:G:OP1	2:B:1246:A:N6	2.39	0.56
9:Z:14:LEU:O	9:Z:18:LYS:HG3	2.06	0.56
7:Q:151:ILE:HG21	7:Q:232:PHE:HB2	1.88	0.55
7:Q:269:GLU:O	7:Q:272:THR:OG1	2.15	0.55
7:Q:158:PHE:HB2	7:Q:170:HIS:HB3	1.88	0.55
4:G:125:LEU:HA	4:G:129:ILE:HD13	1.89	0.55
7:Q:101:VAL:CG2	7:Q:105:GLY:O	2.54	0.55
9:Z:15:LEU:HA	9:Z:18:LYS:CD	2.37	0.55
7:Q:101:VAL:HG23	7:Q:105:GLY:O	2.07	0.54
9:Z:15:LEU:C	9:Z:18:LYS:CD	2.80	0.54
7:Q:262:ASN:HA	7:Q:265:ASN:HD22	1.73	0.54
9:Z:9:LYS:CD	9:Z:11:LEU:N	2.64	0.54
7:Q:243:ASP:HB3	7:Q:246:LEU:HB2	1.90	0.53
9:Z:10:LYS:CE	9:Z:10:LYS:CA	2.86	0.53
7:Q:44:GLN:HE21	7:Q:47:ARG:HB2	1.72	0.53
9:Z:8:ALA:O	9:Z:9:LYS:CG	2.57	0.53
9:Z:12:SER:O	9:Z:16:THR:HG23	2.08	0.53
2:B:1331:C:H41	7:Q:63:LYS:HB3	1.74	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:Z:15:LEU:C	9:Z:18:LYS:CE	2.80	0.52
1:A:3880:A:H2'	9:Z:19:TYR:CD2	2.40	0.52
7:Q:147:GLY:HA2	7:Q:160:THR:HG22	1.91	0.52
1:A:1958:C:OP1	4:G:57:ARG:NH2	2.35	0.52
1:A:1965:A:N7	1:A:1991:A:O2'	2.43	0.52
4:G:24:ALA:HA	7:Q:434:LEU:HD13	1.92	0.52
7:Q:342:LYS:HE3	7:Q:344:LEU:HD23	1.91	0.52
1:A:4409:U:O2'	1:A:4413:U:OP1	2.28	0.52
1:A:3880:A:C8	9:Z:19:TYR:CD2	2.98	0.51
4:G:108:GLU:HA	4:G:112:ILE:HD12	1.92	0.51
1:A:4490:G:OP2	1:A:4490:G:N2	2.29	0.51
9:Z:15:LEU:HA	9:Z:18:LYS:HE2	1.83	0.51
6:P:32:U:H3	6:P:38:A:H61	1.57	0.51
7:Q:27:ALA:HB1	7:Q:100:ILE:HG21	1.93	0.51
2:B:1245:G:O3'	7:Q:65:ARG:NH2	2.44	0.51
7:Q:118:LYS:HD2	7:Q:139:LEU:HA	1.93	0.50
4:G:32:ILE:HD12	4:G:37:LEU:HG	1.93	0.50
7:Q:373:PRO:HG2	7:Q:376:GLU:HB3	1.93	0.50
8:R:7:A:H2'	8:R:7:A:N3	2.26	0.49
4:G:64:ILE:HG23	4:G:68:GLN:H	1.76	0.49
9:Z:12:SER:HA	9:Z:15:LEU:HD11	1.94	0.49
7:Q:38:ILE:HD12	7:Q:124:LEU:HB3	1.95	0.48
7:Q:77:SER:HB3	7:Q:111:ASN:HD21	1.79	0.48
1:A:4511:A:C6	7:Q:190:PHE:CZ	3.00	0.48
7:Q:328:ILE:HG13	7:Q:348:PRO:HD3	1.95	0.48
7:Q:155:GLY:HA3	7:Q:173:THR:HA	1.96	0.48
7:Q:186:SER:O	7:Q:189:ARG:HG2	2.14	0.47
2:B:1236:G:O2'	7:Q:80:GLN:NE2	2.44	0.47
7:Q:160:THR:N	7:Q:167:GLU:O	2.46	0.47
4:G:77:ALA:O	4:G:79:ALA:N	2.48	0.47
9:Z:15:LEU:HA	9:Z:18:LYS:CG	2.44	0.47
1:A:1578:U:O2'	3:D:135:ARG:NH2	2.48	0.47
7:Q:13:GLU:O	7:Q:17:ILE:HG12	2.15	0.47
7:Q:277:ASN:O	7:Q:281:ILE:HD12	2.14	0.47
1:A:1578:U:H2'	3:D:131:ARG:HH21	1.80	0.47
7:Q:17:ILE:O	7:Q:21:ILE:HG12	2.14	0.46
7:Q:421:GLY:O	7:Q:425:GLN:HB2	2.15	0.46
1:A:3880:A:O2'	9:Z:19:TYR:CE1	2.56	0.46
9:Z:14:LEU:HD12	9:Z:17:CYS:HB3	1.96	0.46
7:Q:46:SER:O	7:Q:50:LYS:HG2	2.15	0.46
1:A:3881:C:O4'	9:Z:19:TYR:HE1	1.98	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:G:12:VAL:HG11	4:G:61:LYS:HB3	1.98	0.46
7:Q:101:VAL:HG23	7:Q:106:LYS:C	2.40	0.46
7:Q:152:ASP:OD1	7:Q:153:GLY:N	2.46	0.46
1:A:3876:G:H1'	9:Z:11:LEU:HD13	1.99	0.45
4:G:24:ALA:H	4:G:27:ALA:HB2	1.81	0.45
7:Q:150:VAL:HB	7:Q:157:LEU:HB3	1.99	0.45
7:Q:14:ILE:HG23	7:Q:117:PHE:CE1	2.51	0.45
7:Q:268:ILE:O	7:Q:271:SER:OG	2.25	0.45
6:P:76:A:O2'	7:Q:185:GLN:HG2	2.17	0.45
1:A:1981:G:H1'	1:A:1998:A:H61	1.82	0.45
1:A:4485:G:O2'	1:A:4488:C:OP2	2.30	0.45
4:G:72:GLU:HG3	4:G:74:VAL:HG22	1.99	0.45
7:Q:134:GLU:HA	7:Q:137:THR:HG23	1.99	0.45
7:Q:275:LEU:HA	7:Q:278:VAL:HG22	1.98	0.45
1:A:4511:A:N9	7:Q:190:PHE:CE2	2.84	0.44
4:G:16:ARG:HH12	4:G:23:GLY:HA2	1.83	0.44
7:Q:333:LEU:HD21	7:Q:356:HIS:HD2	1.82	0.44
1:A:3760:C:N4	1:A:3784:C:OP2	2.50	0.44
7:Q:99:THR:HA	7:Q:109:LYS:HA	1.99	0.44
1:A:4370:G:H1	1:A:4398:U:H3	1.65	0.44
4:G:76:SER:OG	4:G:77:ALA:N	2.51	0.44
4:G:41:LYS:HB2	4:G:71:ILE:H	1.82	0.43
2:B:1246:A:N3	2:B:1251:A:O2'	2.49	0.43
7:Q:188:LEU:O	7:Q:192:ARG:HG2	2.17	0.43
1:A:4399:U:H4'	1:A:4400:U:H5'	1.99	0.43
3:D:131:ARG:HG3	3:D:137:ASN:OD1	2.18	0.43
7:Q:157:LEU:HD21	7:Q:168:VAL:HG22	2.00	0.43
1:A:3879:G:O3'	9:Z:19:TYR:OH	2.35	0.43
4:G:74:VAL:HA	4:G:75:PRO:HD2	1.85	0.42
9:Z:9:LYS:HG3	9:Z:12:SER:H	1.83	0.42
7:Q:198:ARG:O	7:Q:202:VAL:HG23	2.19	0.42
2:B:1708:C:H42	2:B:1826:G:H1	1.68	0.42
7:Q:18:LYS:O	7:Q:22:LYS:HG2	2.20	0.42
7:Q:194:ARG:HG2	7:Q:198:ARG:HE	1.83	0.42
9:Z:9:LYS:CG	9:Z:10:LYS:N	2.82	0.42
1:A:1963:G:H5'	7:Q:436:ASP:HA	2.01	0.42
9:Z:12:SER:HA	9:Z:15:LEU:HD12	2.01	0.42
1:A:4402:U:H2'	1:A:4403:G:C8	2.54	0.42
7:Q:242:PHE:HE2	7:Q:247:GLN:HA	1.85	0.41
1:A:4481:A:OP1	1:A:4483:G:O2'	2.27	0.41
7:Q:43:ASP:CG	7:Q:47:ARG:HH21	2.27	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
9:Z:20:ILE:HB	9:Z:21:PRO:HD3	2.03	0.41
9:Z:14:LEU:C	9:Z:17:CYS:H	2.29	0.41
1:A:3731:A:O2'	2:B:1826:G:O2'	2.22	0.41
7:Q:251:LEU:HD23	7:Q:274:VAL:HG23	2.01	0.41
7:Q:25:GLU:OE1	7:Q:133:THR:OG1	2.24	0.41
1:A:4497:G:H1	1:A:4515:U:H3	1.69	0.40
4:G:22:VAL:HA	4:G:46:ILE:HG21	2.02	0.40
4:G:85:LEU:HD22	4:G:100:HIS:CE1	2.55	0.40
7:Q:314:MET:HB3	7:Q:433:ASP:OD1	2.21	0.40
7:Q:63:LYS:HB2	8:R:7:A:H5''	2.04	0.40
7:Q:126:LEU:HD23	7:Q:132:HIS:CD2	2.56	0.40
1:A:3880:A:H2'	9:Z:19:TYR:CE1	2.50	0.40
4:G:18:THR:HG21	7:Q:434:LEU:HD11	2.03	0.40

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	D	19/184 (10%)	18 (95%)	1 (5%)	0	100	100
4	G	136/165 (82%)	103 (76%)	27 (20%)	6 (4%)	2	19
5	H	10/427 (2%)	10 (100%)	0	0	100	100
7	Q	429/431 (100%)	390 (91%)	33 (8%)	6 (1%)	9	37
9	Z	20/22 (91%)	12 (60%)	2 (10%)	6 (30%)	0	0
All	All	614/1229 (50%)	533 (87%)	63 (10%)	18 (3%)	5	25

All (18) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
4	G	98	ILE

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Mol	Chain	Res	Type
4	G	109	ILE
7	Q	242	PHE
9	Z	3	PRO
9	Z	4	LEU
9	Z	6	LEU
9	Z	21	PRO
4	G	88	PRO
4	G	78	SER
7	Q	130	LYS
7	Q	215	SER
7	Q	306	GLU
7	Q	324	GLU
9	Z	7	SER
4	G	64	ILE
7	Q	381	ASN
9	Z	5	VAL
4	G	102	GLY

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	
3	D	19/163 (12%)	19 (100%)	0	100	100
4	G	114/137 (83%)	114 (100%)	0	100	100
5	H	8/348 (2%)	8 (100%)	0	100	100
7	Q	371/371 (100%)	371 (100%)	0	100	100
9	Z	15/21 (71%)	10 (67%)	5 (33%)	0	1
All	All	527/1040 (51%)	522 (99%)	5 (1%)	68	74

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
9	Z	9	LYS
9	Z	10	LYS

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Mol	Chain	Res	Type
9	Z	13	SER
9	Z	15	LEU
9	Z	18	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
4	G	103	ASN
7	Q	44	GLN
7	Q	132	HIS
7	Q	211	GLN
7	Q	265	ASN
7	Q	282	GLN
7	Q	336	GLN
7	Q	350	GLN
7	Q	356	HIS
7	Q	366	HIS

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
1	A	234/5025 (4%)	33 (14%)	1 (0%)
2	B	59/1869 (3%)	11 (18%)	0
6	P	16/18 (88%)	2 (12%)	0
8	R	8/9 (88%)	3 (37%)	0
All	All	317/6921 (4%)	49 (15%)	1 (0%)

All (49) RNA backbone outliers are listed below:

Mol	Chain	Res	Type
1	A	1623	G
1	A	1955	U
1	A	1956	G
1	A	1957	G
1	A	1965	A
1	A	1966	G
1	A	1976	G
1	A	1980	A
1	A	1984	G
1	A	1985	U

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Mol	Chain	Res	Type
1	A	2774	C
1	A	3731	A
1	A	3732	A
1	A	3739	C
1	A	3783	G
1	A	3786	U
1	A	3879	G
1	A	3880	A
1	A	3881	C
1	A	4350	C
1	A	4357	A
1	A	4359	A
1	A	4361	C
1	A	4385	A
1	A	4407	C
1	A	4412	A
1	A	4483	G
1	A	4486	A
1	A	4493	U
1	A	4494	U
1	A	4496	A
1	A	4511	A
1	A	4512	G
2	B	614	C
2	B	627	U
2	B	629	A
2	B	1242	U
2	B	1247	C
2	B	1248	U
2	B	1251	A
2	B	1330	G
2	B	1826	G
2	B	1827	U
2	B	1828	C
6	P	41	C
6	P	76	A
8	R	4	U
8	R	7	A
8	R	8	A

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
1	A	3730	U

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

There are no ligands in this entry.

5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

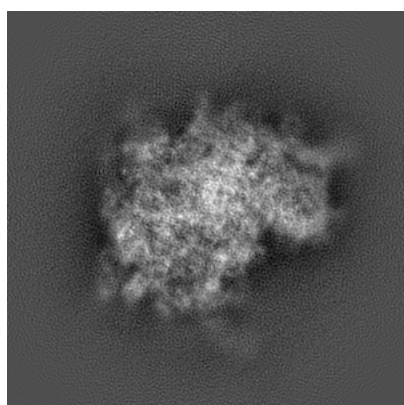
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-3099. These allow visual inspection of the internal detail of the map and identification of artifacts.

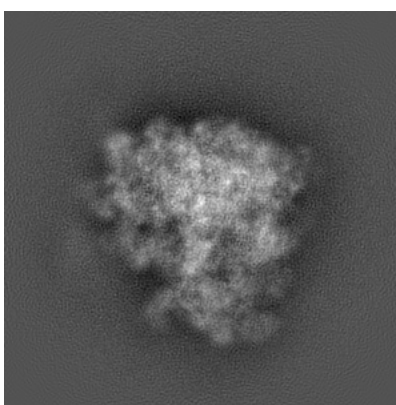
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

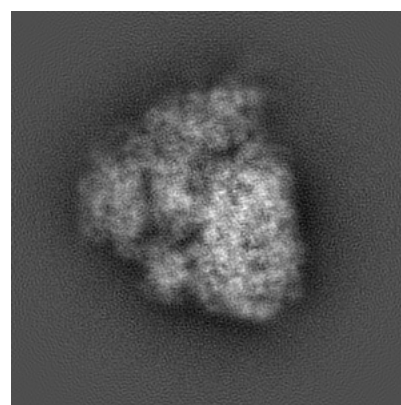
6.1.1 Primary map



X



Y

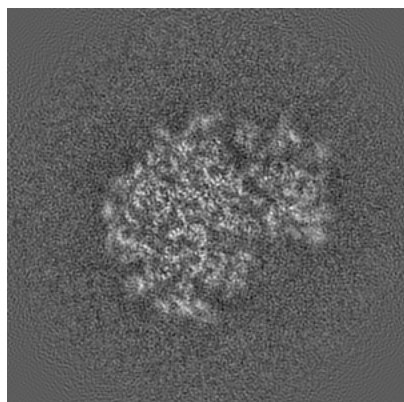


Z

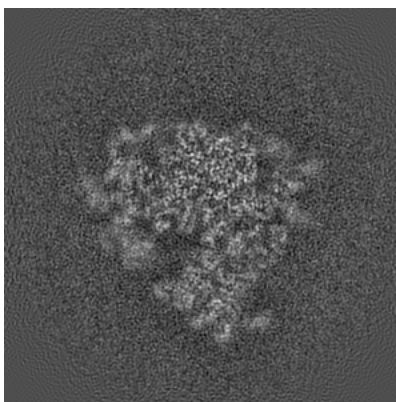
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

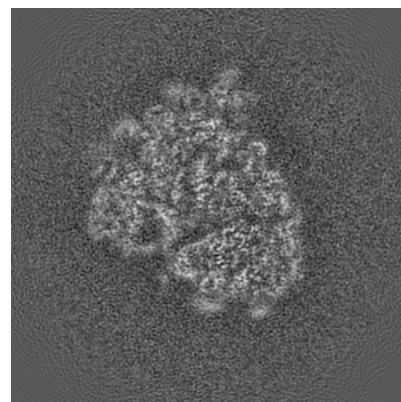
6.2.1 Primary map



X Index: 210



Y Index: 210

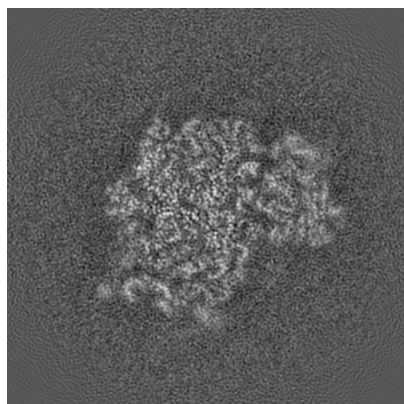


Z Index: 210

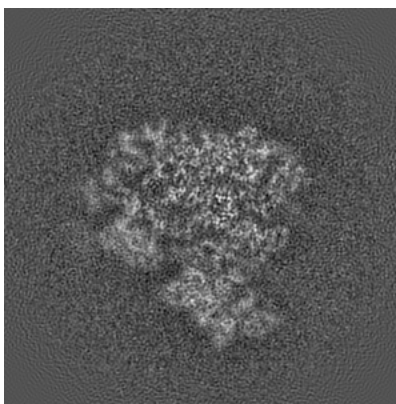
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

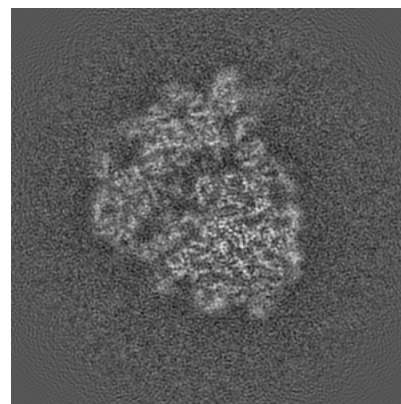
6.3.1 Primary map



X Index: 221



Y Index: 217

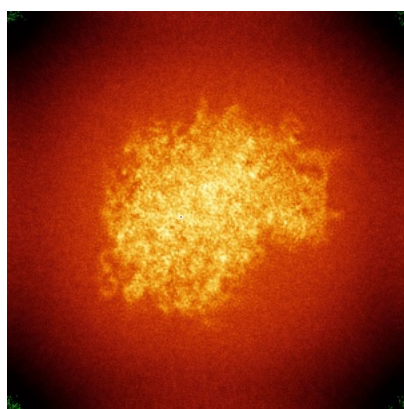


Z Index: 204

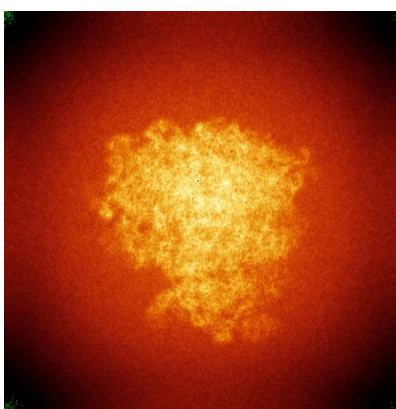
The images above show the largest variance slices of the map in three orthogonal directions.

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

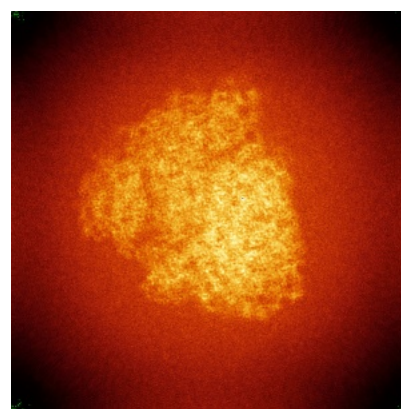
6.4.1 Primary map



X



Y

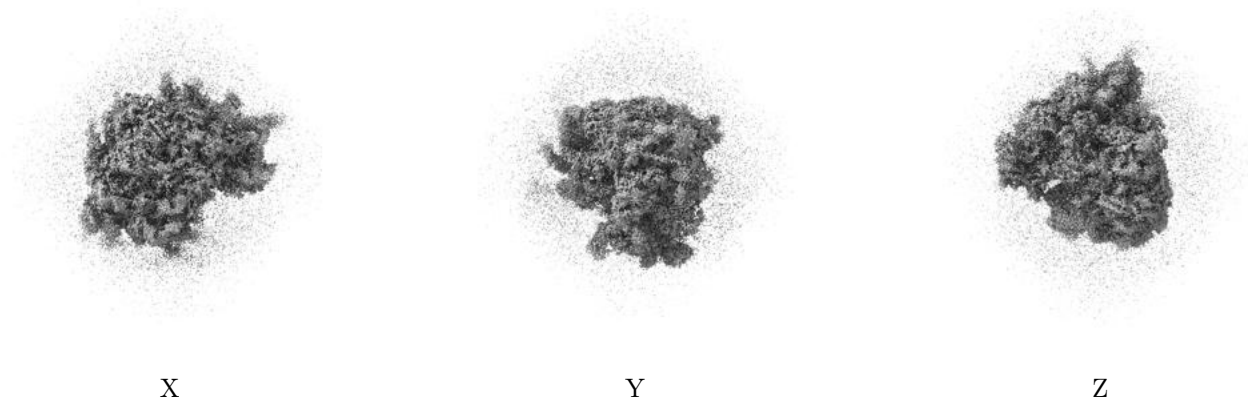


Z

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

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.0008. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

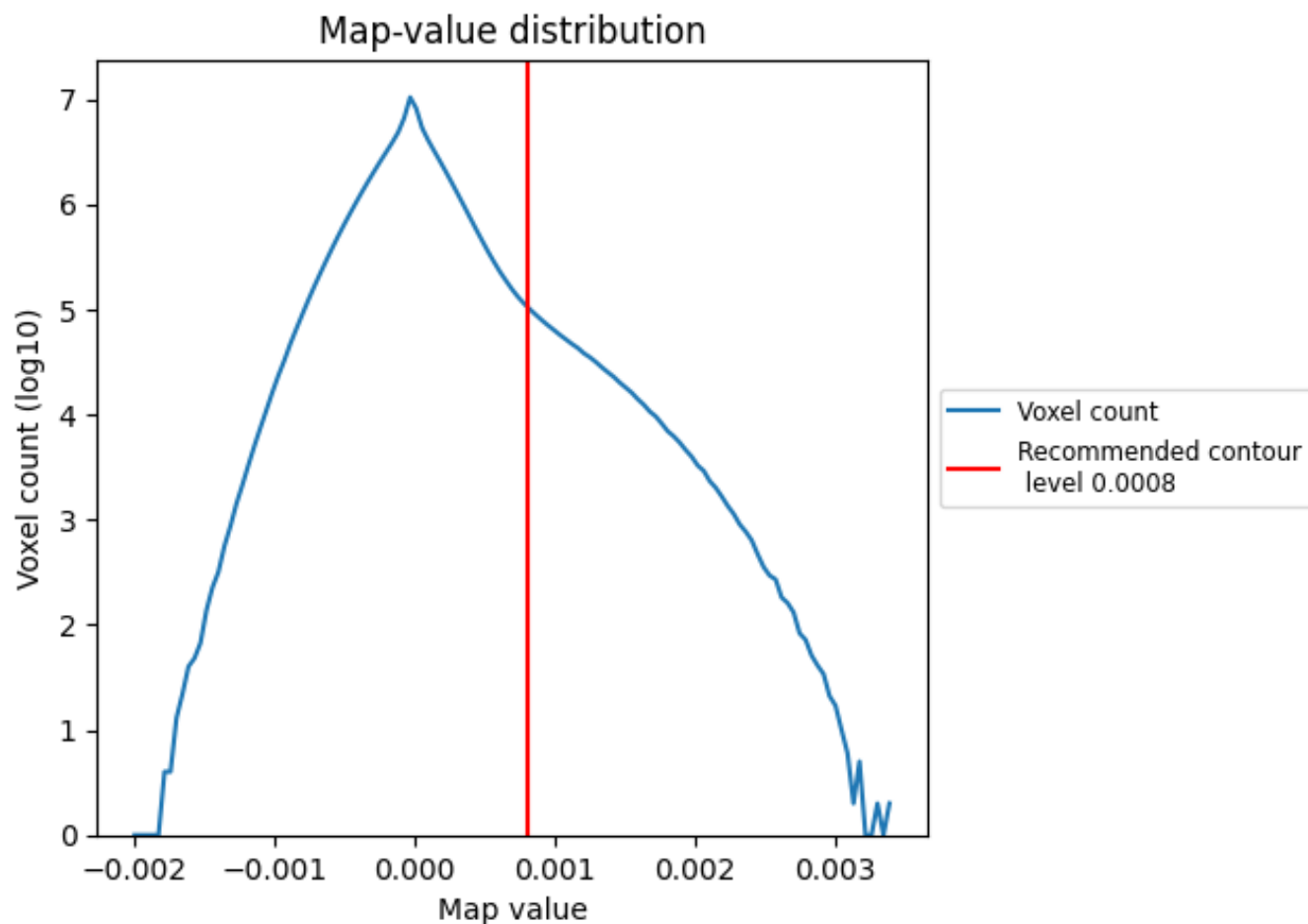
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

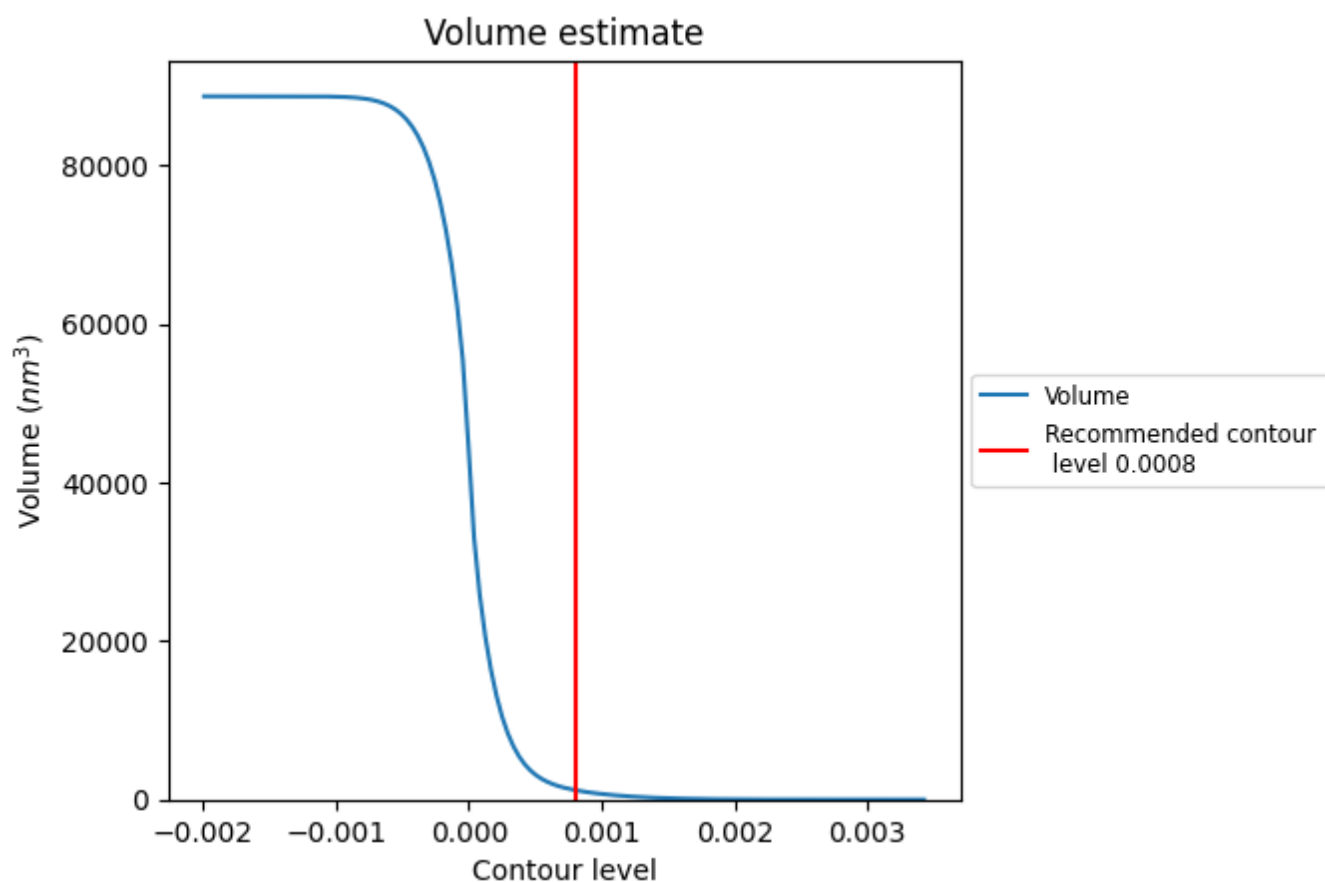
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

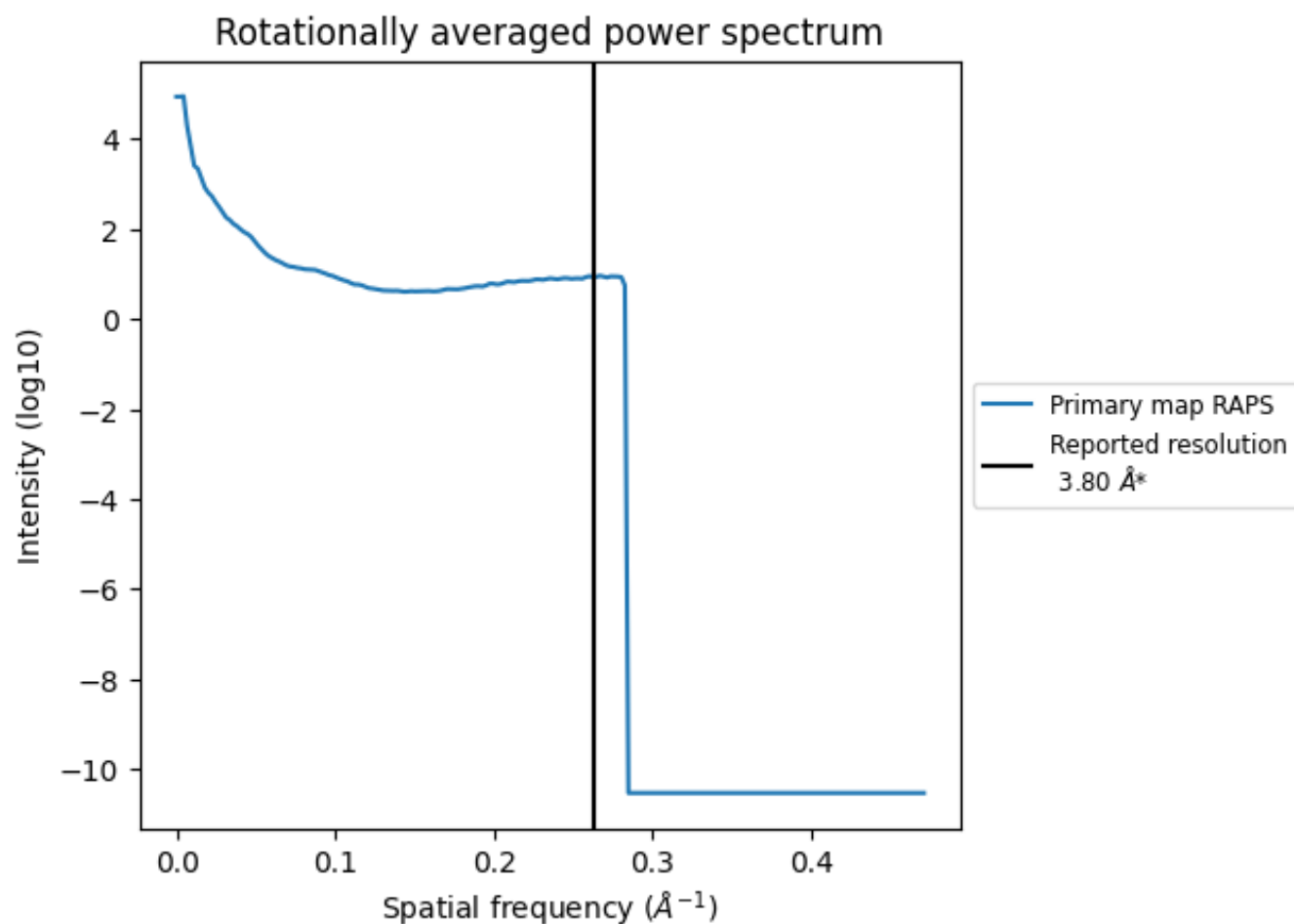
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1191 nm³; this corresponds to an approximate mass of 1076 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 0.263 $\text{\AA}^{-1}$

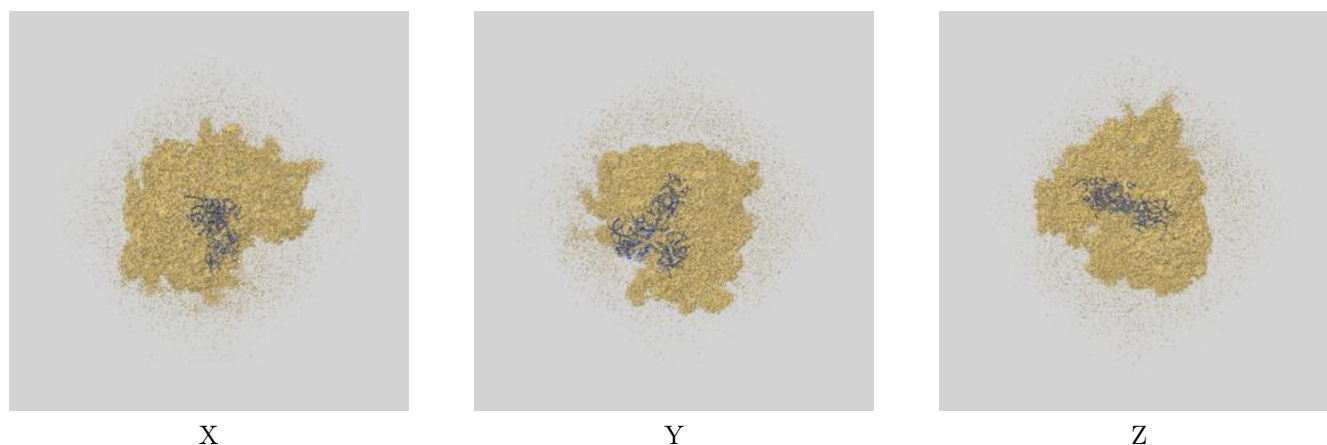
8 Fourier-Shell correlation ⓘ

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-3099 and PDB model 5A8L. Per-residue inclusion information can be found in [section 3](#) on [page 6](#).

9.1 Map-model overlay [i](#)



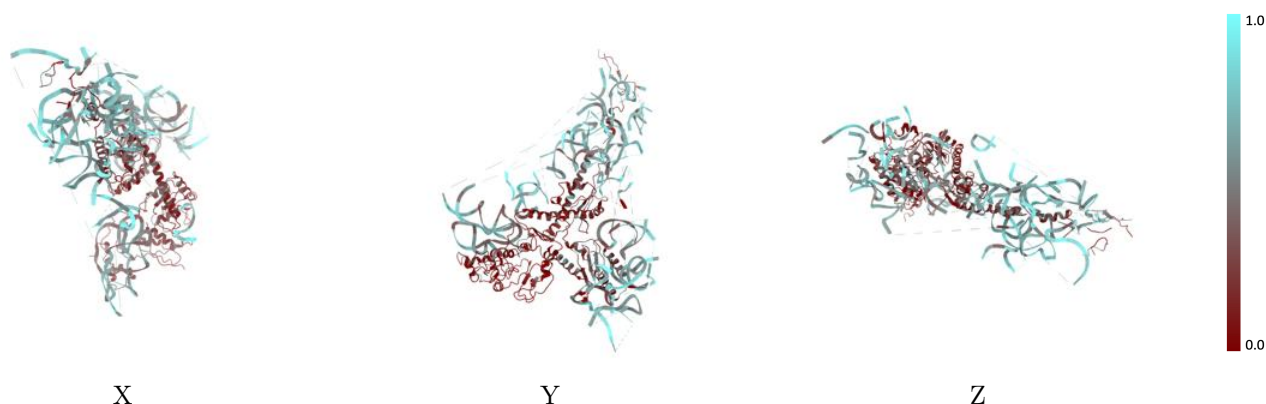
The images above show the 3D surface view of the map at the recommended contour level 0.0008 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



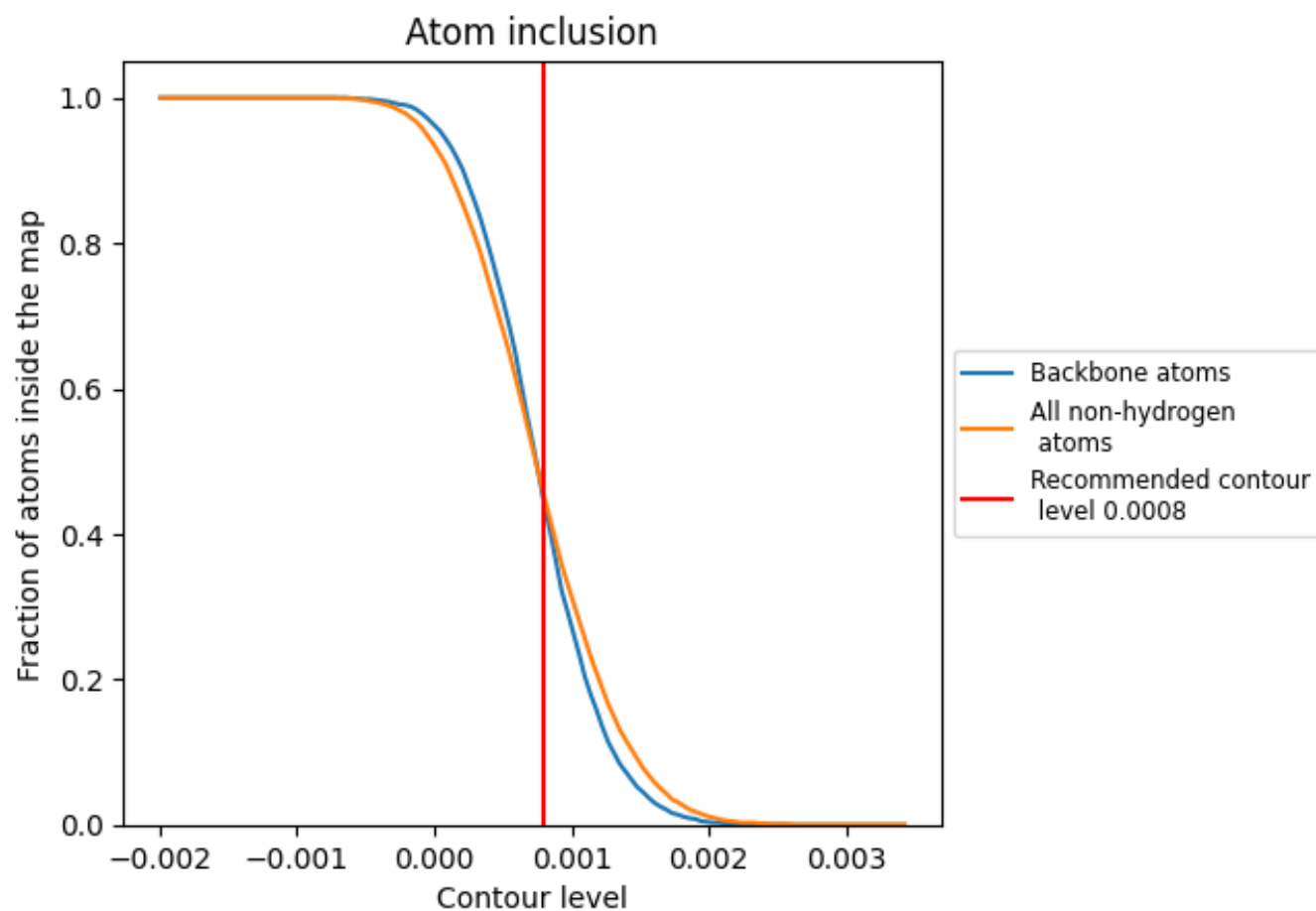
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.0008).

9.4 Atom inclusion [i](#)



At the recommended contour level, 44% of all backbone atoms, 45% 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 (0.0008) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.4520	0.2280
A	0.6180	0.2860
B	0.6340	0.2920
D	0.4430	0.2660
G	0.0910	0.0210
H	0.4150	0.2690
P	0.6100	0.2940
Q	0.2090	0.1560
R	0.5160	0.2830
Z	0.2430	0.3280

1.0

0.0

<0.0