



wwPDB EM Validation Summary Report ⓘ

Mar 26, 2026 – 04:30 PM UTC

PDB ID : 6N7X / pdb_00006n7x
EMDB ID : EMD-8622
Title : S. cerevisiae U1 snRNP
Authors : Li, X.; Liu, S.; Jiang, J.; Zhang, L.; Espinosa, S.; Hill, R.C.; Hansen, K.C.;
Zhou, Z.H.; Zhao, R.
Deposited on : 2018-11-28
Resolution : 3.60 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:

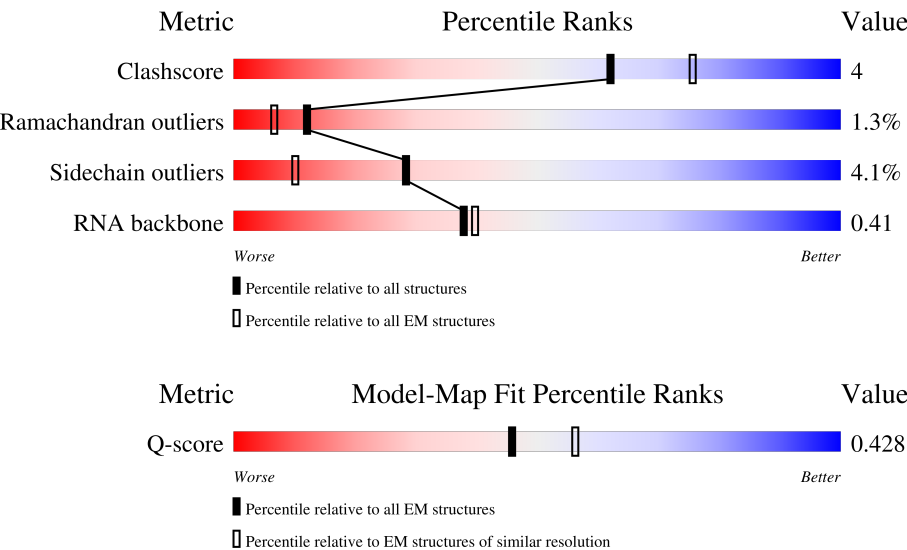
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 3.60 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	11569 (3.20 - 4.20)

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	300	
2	B	231	
3	C	298	

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Mol	Chain	Length	Quality of chain
4	D	544	
5	E	629	
6	F	523	
7	G	492	
8	H	620	
9	K	196	
10	L	146	
11	M	110	
12	N	101	
13	O	94	
14	P	86	
15	Q	77	
16	R	568	

2 Entry composition

There are 16 unique types of molecules in this entry. The entry contains 27183 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 U1 small nuclear ribonucleoprotein 70 kDa homolog.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	186	Total	C	N	O	S	0	0
			1219	752	232	233	2		

- Molecule 2 is a protein called U1 small nuclear ribonucleoprotein C.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	176	Total	C	N	O	S	0	0
			1412	882	271	255	4		

- Molecule 3 is a protein called U1 small nuclear ribonucleoprotein A.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	131	Total	C	N	O	S	0	0
			1041	663	190	185	3		

- Molecule 4 is a protein called U1 small nuclear ribonucleoprotein component PRP42.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	539	Total	C	N	O	S	0	0
			4447	2918	702	813	14		

- Molecule 5 is a protein called Pre-mRNA-processing factor 39.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	583	Total	C	N	O	S	0	0
			3908	2477	680	744	7		

- Molecule 6 is a protein called Protein NAM8.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	187	Total	C	N	O	S	0	0
			1414	889	244	271	10		

- Molecule 7 is a protein called 56 kDa U1 small nuclear ribonucleoprotein component.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	G	229	Total	C	N	O	S	0	0
			1841	1202	300	328	11		

- Molecule 8 is a protein called U1 small nuclear ribonucleoprotein component SNU71.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	H	52	Total	C	N	O	S	0	0
			420	266	74	79	1		

- Molecule 9 is a protein called Small nuclear ribonucleoprotein-associated protein B.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	K	124	Total	C	N	O	S	0	0
			1009	637	191	178	3		

- Molecule 10 is a protein called Small nuclear ribonucleoprotein Sm D1.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	L	119	Total	C	N	O	S	0	0
			917	575	163	176	3		

- Molecule 11 is a protein called Small nuclear ribonucleoprotein Sm D2.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	M	96	Total	C	N	O	S	0	0
			739	476	138	121	4		

- Molecule 12 is a protein called Small nuclear ribonucleoprotein Sm D3.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	N	92	Total	C	N	O	S	0	0
			692	441	123	127	1		

- Molecule 13 is a protein called Small nuclear ribonucleoprotein E.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	O	74	Total	C	N	O	S	0	0
			526	346	87	90	3		

- Molecule 14 is a protein called Small nuclear ribonucleoprotein F.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	P	68	Total	C	N	O	S	0	0
			518	337	96	84	1		

- Molecule 15 is a protein called Small nuclear ribonucleoprotein G.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	Q	75	Total	C	N	O	S	0	0
			543	346	94	102	1		

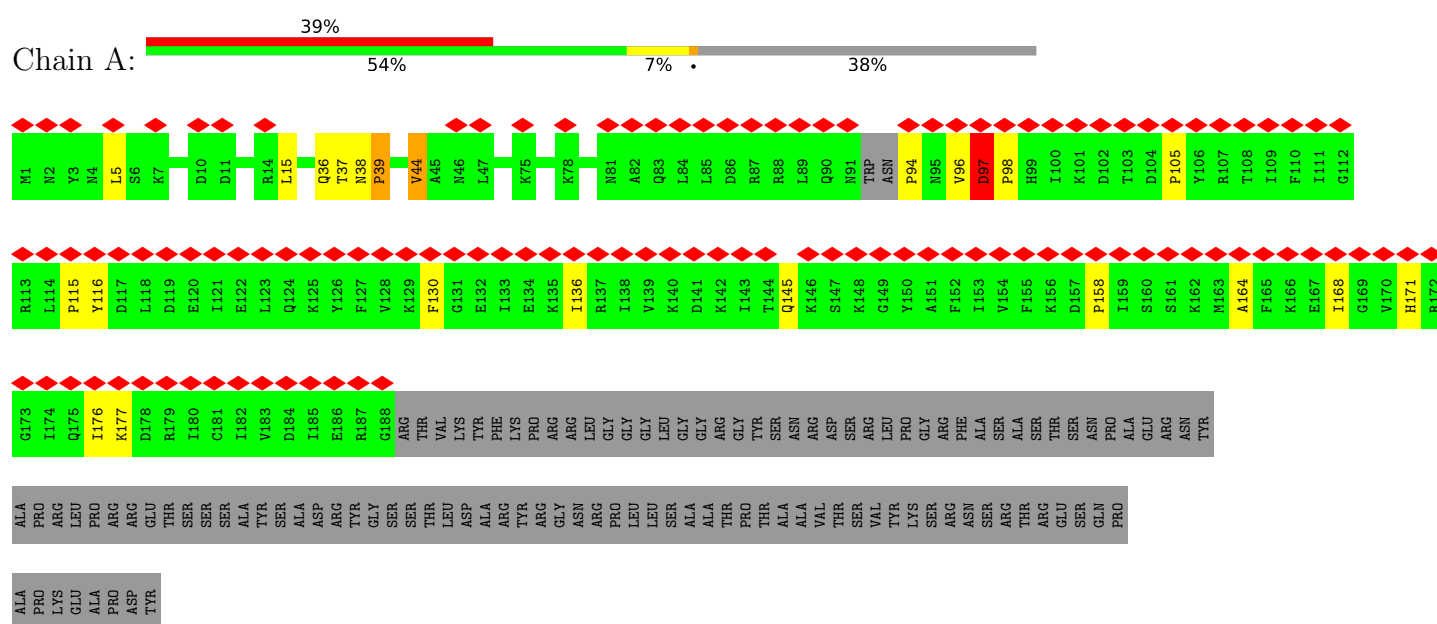
- Molecule 16 is a RNA chain called U1 snRNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	R	308	Total	C	N	O	P	0	0
			6537	2923	1123	2183	308		

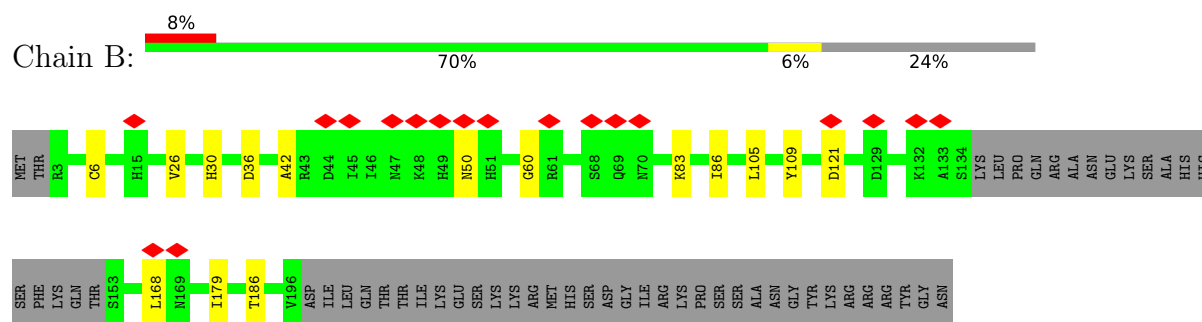
3 Residue-property plots [i](#)

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.

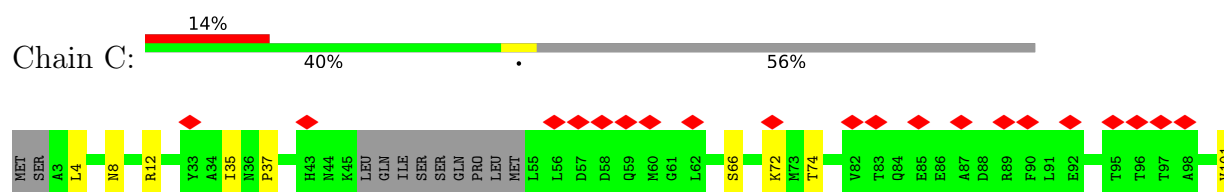
- Molecule 1: U1 small nuclear ribonucleoprotein 70 kDa homolog

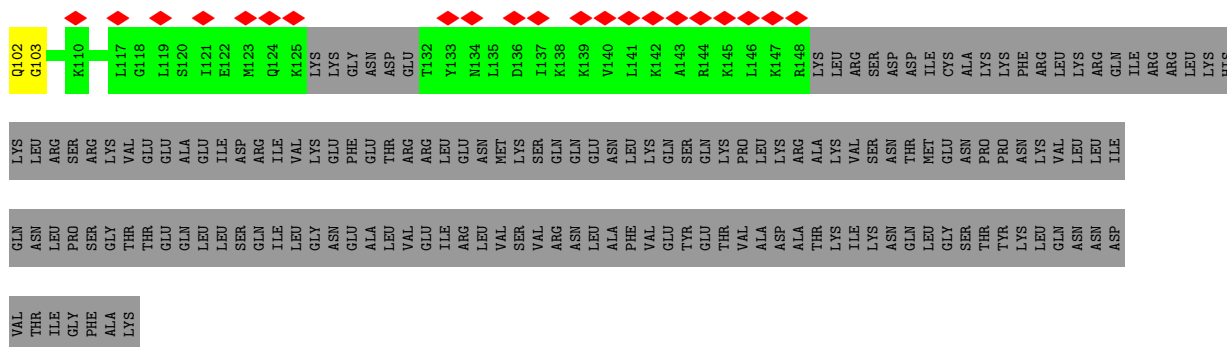


- Molecule 2: U1 small nuclear ribonucleoprotein C



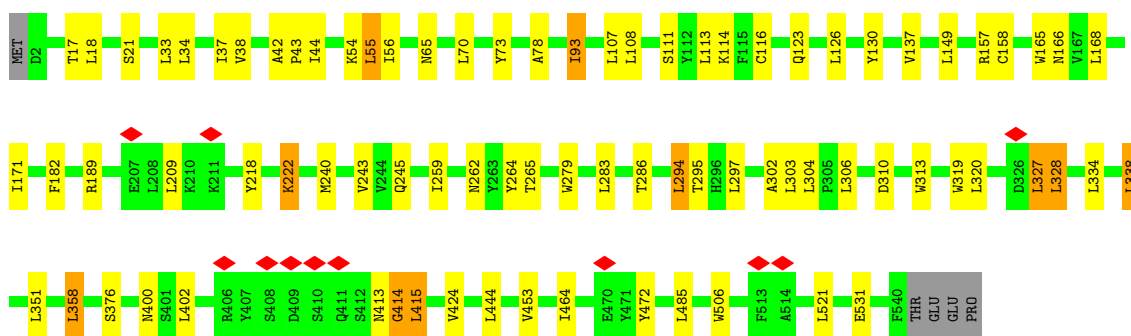
- Molecule 3: U1 small nuclear ribonucleoprotein A





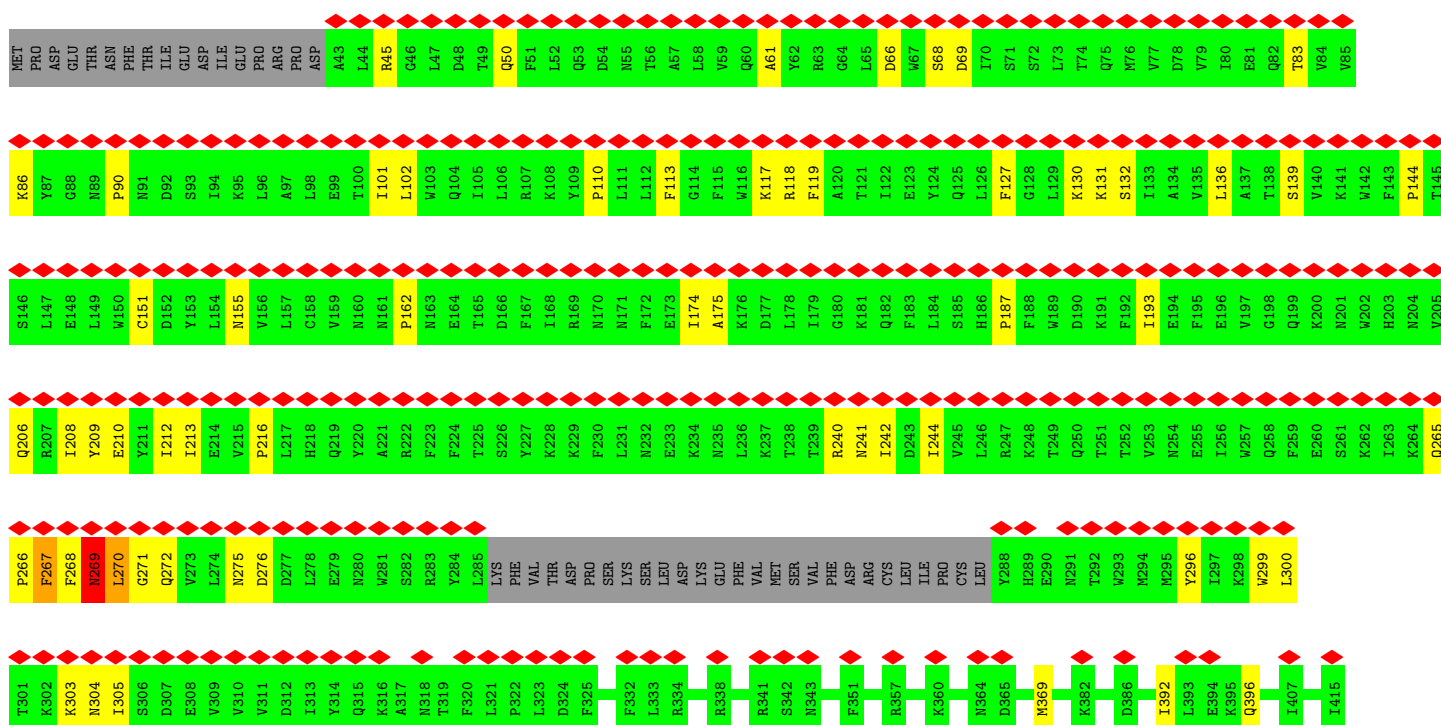
• Molecule 4: U1 small nuclear ribonucleoprotein component PRP42

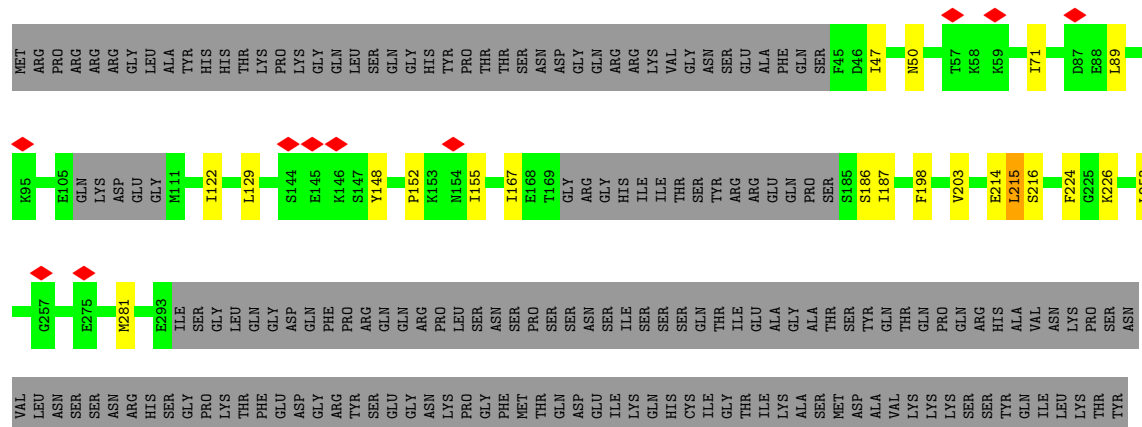
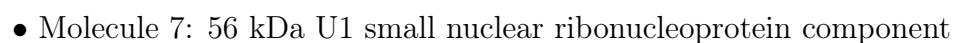
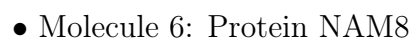
Chain D: 84% 13% ..

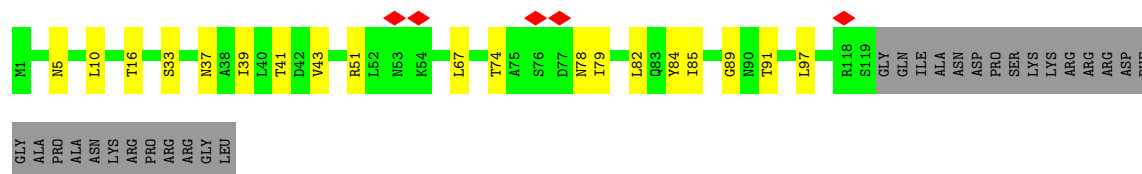


• Molecule 5: Pre-mRNA-processing factor 39

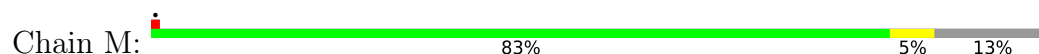
Chain E: 51% 80% 12% 7%



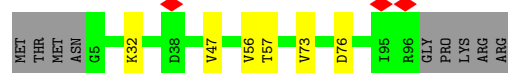
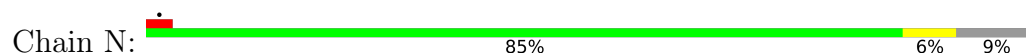




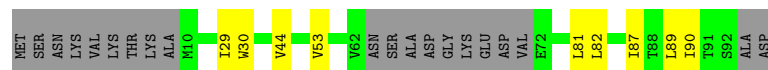
- Molecule 11: Small nuclear ribonucleoprotein Sm D2



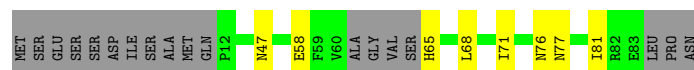
- Molecule 12: Small nuclear ribonucleoprotein Sm D3



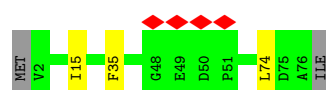
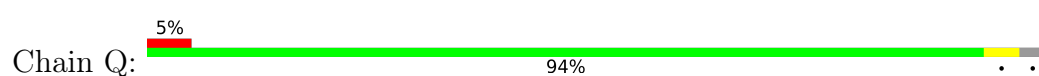
- Molecule 13: Small nuclear ribonucleoprotein E



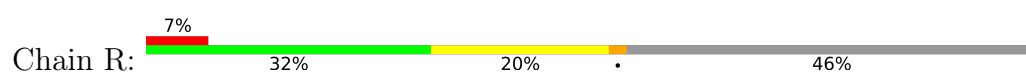
- Molecule 14: Small nuclear ribonucleoprotein F



- Molecule 15: Small nuclear ribonucleoprotein G



- Molecule 16: U1 snRNA





4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	352900	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 ($e^-/\text{\AA}^2$)	51.9	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	0.343	Depositor
Minimum map value	-0.213	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	0.007	Depositor
Recommended contour level	0.06	Depositor
Map size ($\text{\AA}$)	435.2, 435.2, 435.2	wwPDB
Map dimensions	320, 320, 320	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.36, 1.36, 1.36	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	0.66	1/1237 (0.1%)	1.13	9/1691 (0.5%)
2	B	0.58	0/1441	0.78	0/1941
3	C	0.59	0/1055	0.75	0/1417
4	D	0.58	0/4561	0.80	1/6186 (0.0%)
5	E	0.69	0/3971	1.10	29/5462 (0.5%)
6	F	0.61	0/1441	0.84	1/1957 (0.1%)
7	G	0.58	0/1882	0.75	2/2537 (0.1%)
8	H	0.61	0/428	0.69	0/575
9	K	0.65	0/1016	0.87	0/1355
10	L	0.64	0/926	0.89	1/1257 (0.1%)
11	M	0.53	0/751	0.76	0/1014
12	N	0.65	0/704	0.82	0/957
13	O	0.69	0/535	0.84	0/730
14	P	0.63	0/529	0.82	0/715
15	Q	0.62	0/548	0.88	0/745
16	R	0.41	0/7295	0.78	12/11341 (0.1%)
All	All	0.57	1/28320 (0.0%)	0.86	55/39880 (0.1%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	5
5	E	0	11
All	All	0	16

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	97	ASP	CA-C	5.45	1.59	1.52

The worst 5 of 55 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	115	PRO	N-CA-CB	8.64	110.85	103.25
16	R	66	U	C4'-C3'-O3'	8.28	121.83	109.40
16	R	258	U	C2'-C3'-O3'	7.46	120.68	109.50
16	R	64	A	C4'-C3'-O3'	7.27	120.31	109.40
1	A	97	ASP	CA-C-N	7.16	128.78	119.84

There are no chirality outliers.

5 of 16 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	136	ILE	Peptide
1	A	145	GLN	Peptide
1	A	176	ILE	Peptide
1	A	96	VAL	Peptide
1	A	97	ASP	Peptide

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1219	0	943	5	0
2	B	1412	0	1394	10	0
3	C	1041	0	1087	3	0
4	D	4447	0	4337	54	0
5	E	3908	0	3102	37	0
6	F	1414	0	1363	12	0
7	G	1841	0	1813	9	0
8	H	420	0	421	32	0
9	K	1009	0	1100	15	0
10	L	917	0	961	8	0
11	M	739	0	756	3	0
12	N	692	0	710	1	0
13	O	526	0	515	6	0
14	P	518	0	502	5	0
15	Q	543	0	538	1	0
16	R	6537	0	3294	60	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	27183	0	22836	217	0

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 4.

The worst 5 of 217 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)
8:H:45:HIS:CD2	8:H:46:ILE:HG13	1.50	1.43
8:H:45:HIS:NE2	8:H:46:ILE:HG13	1.57	1.19
4:D:376:SER:HB2	8:H:50:ARG:NH1	1.62	1.15
4:D:54:LYS:CE	8:H:1:MET:HE2	1.79	1.12
4:D:54:LYS:HE3	8:H:1:MET:CE	1.82	1.09

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	A	182/300 (61%)	155 (85%)	21 (12%)	6 (3%)	3	24
2	B	172/231 (74%)	153 (89%)	18 (10%)	1 (1%)	21	54
3	C	125/298 (42%)	112 (90%)	9 (7%)	4 (3%)	3	24
4	D	537/544 (99%)	494 (92%)	39 (7%)	4 (1%)	18	51
5	E	579/629 (92%)	517 (89%)	46 (8%)	16 (3%)	4	26
6	F	181/523 (35%)	168 (93%)	11 (6%)	2 (1%)	11	43
7	G	223/492 (45%)	208 (93%)	15 (7%)	0	100	100
8	H	50/620 (8%)	48 (96%)	2 (4%)	0	100	100
9	K	120/196 (61%)	107 (89%)	13 (11%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
10	L	117/146 (80%)	103 (88%)	12 (10%)	2 (2%)	7	35
11	M	94/110 (86%)	90 (96%)	4 (4%)	0	100	100
12	N	90/101 (89%)	87 (97%)	3 (3%)	0	100	100
13	O	70/94 (74%)	69 (99%)	1 (1%)	0	100	100
14	P	64/86 (74%)	59 (92%)	5 (8%)	0	100	100
15	Q	73/77 (95%)	64 (88%)	9 (12%)	0	100	100
All	All	2677/4447 (60%)	2434 (91%)	208 (8%)	35 (1%)	12	39

5 of 35 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	97	ASP
1	A	98	PRO
3	C	102	GLN
4	D	413	ASN
5	E	86	LYS

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	A	84/265 (32%)	82 (98%)	2 (2%)	43	63
2	B	154/214 (72%)	149 (97%)	5 (3%)	34	58
3	C	113/273 (41%)	110 (97%)	3 (3%)	39	61
4	D	483/519 (93%)	456 (94%)	27 (6%)	19	47
5	E	291/603 (48%)	284 (98%)	7 (2%)	43	63
6	F	154/451 (34%)	150 (97%)	4 (3%)	40	62
7	G	199/448 (44%)	195 (98%)	4 (2%)	48	66
8	H	48/568 (8%)	42 (88%)	6 (12%)	4	22
9	K	113/176 (64%)	106 (94%)	7 (6%)	16	44
10	L	106/129 (82%)	97 (92%)	9 (8%)	10	35

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
11	M	77/103 (75%)	75 (97%)	2 (3%)	40	62
12	N	76/89 (85%)	71 (93%)	5 (7%)	15	43
13	O	52/83 (63%)	52 (100%)	0	100	100
14	P	51/77 (66%)	50 (98%)	1 (2%)	48	66
15	Q	54/66 (82%)	52 (96%)	2 (4%)	30	56
All	All	2055/4064 (51%)	1971 (96%)	84 (4%)	28	53

5 of 84 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
8	H	50	ARG
10	L	82	LEU
9	K	31	ILE
10	L	5	ASN
11	M	105	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 48 such sidechains are listed below:

Mol	Chain	Res	Type
6	F	497	GLN
10	L	21	ASN
7	G	207	HIS
8	H	51	ASN
10	L	56	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
16	R	301/568 (52%)	69 (22%)	17 (5%)

5 of 69 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
16	R	17	A
16	R	51	A
16	R	54	C
16	R	55	G
16	R	56	C

5 of 17 RNA pucker outliers are listed below:

Mol	Chain	Res	Type
16	R	279	U
16	R	564	A
16	R	113	G
16	R	152	G
16	R	186	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.

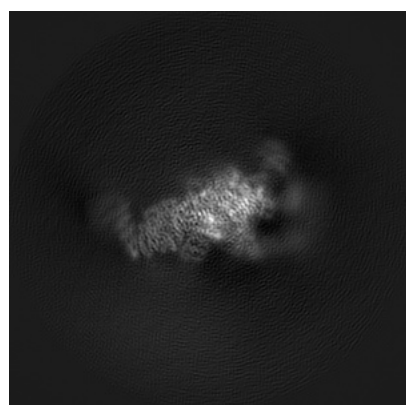
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-8622. These allow visual inspection of the internal detail of the map and identification of artifacts.

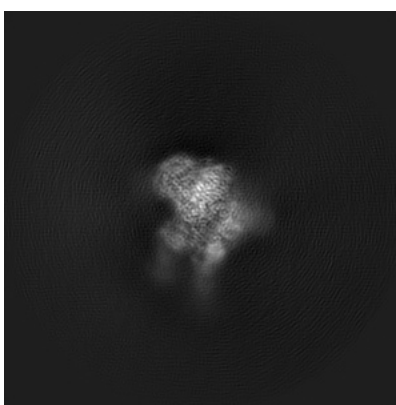
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

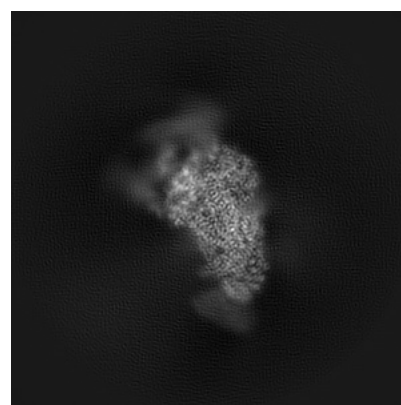
6.1.1 Primary map



X



Y

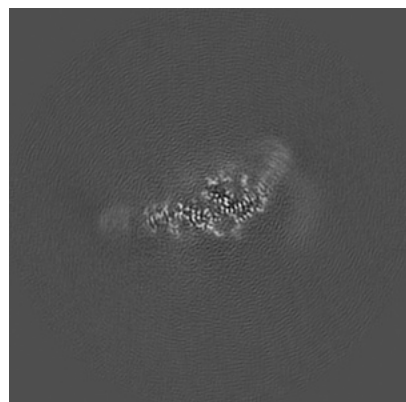


Z

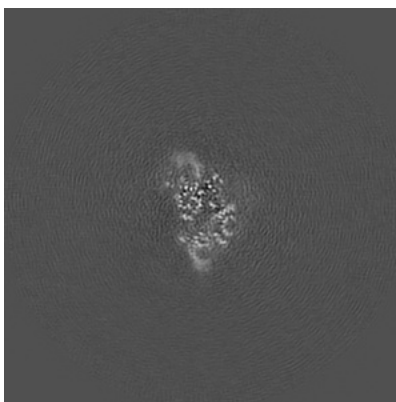
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

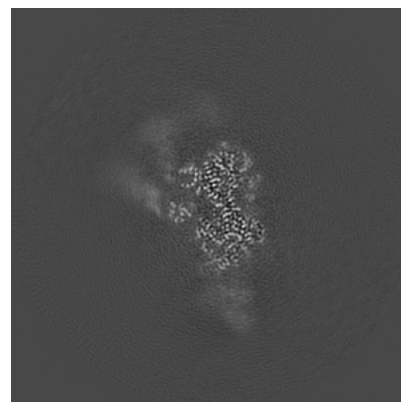
6.2.1 Primary map



X Index: 160



Y Index: 160

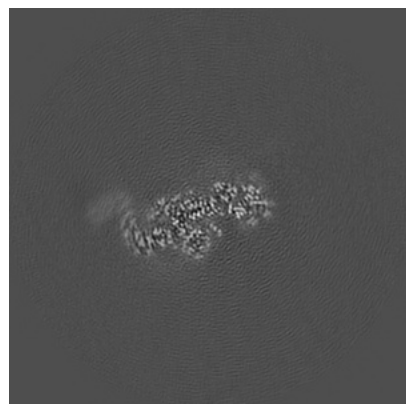


Z Index: 160

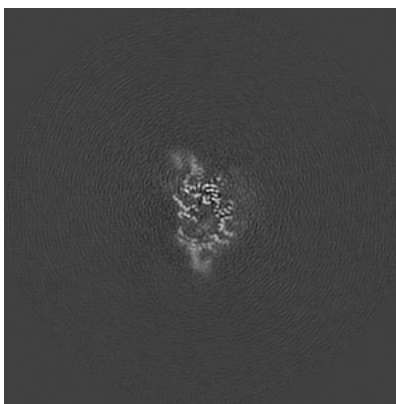
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

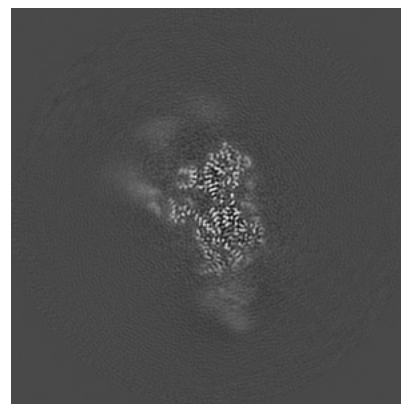
6.3.1 Primary map



X Index: 178



Y Index: 162

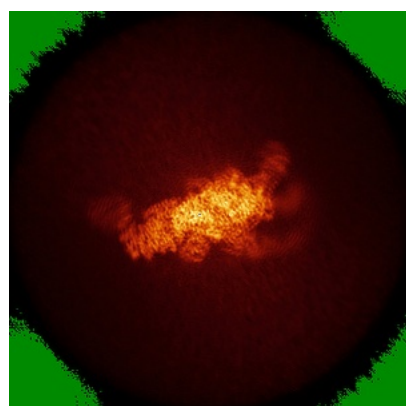


Z Index: 158

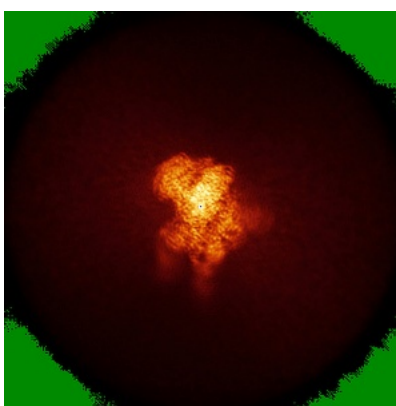
The images above show the largest variance slices of the map in three orthogonal directions.

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

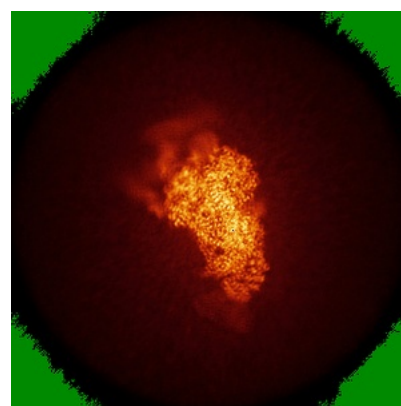
6.4.1 Primary map



X



Y

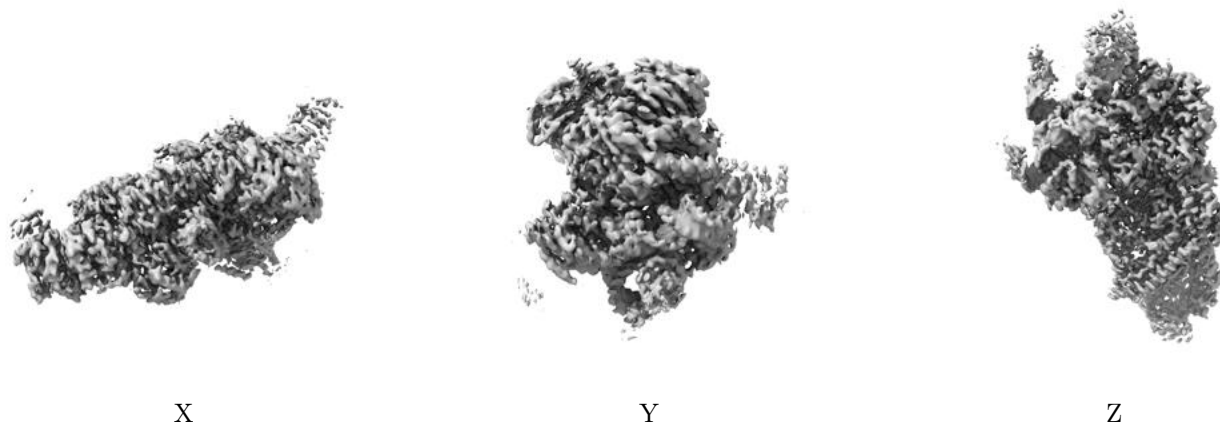


Z

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 0.06. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

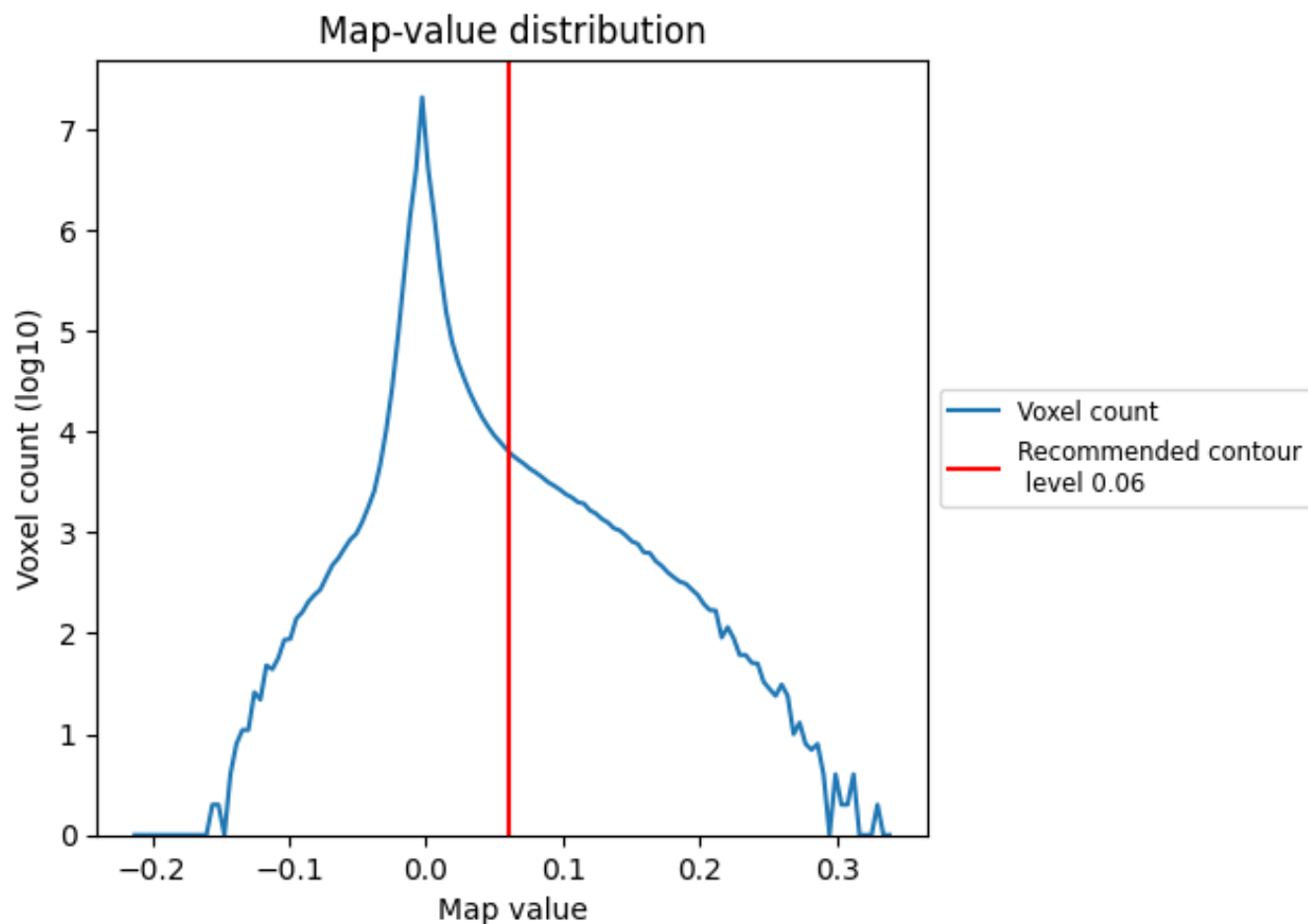
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

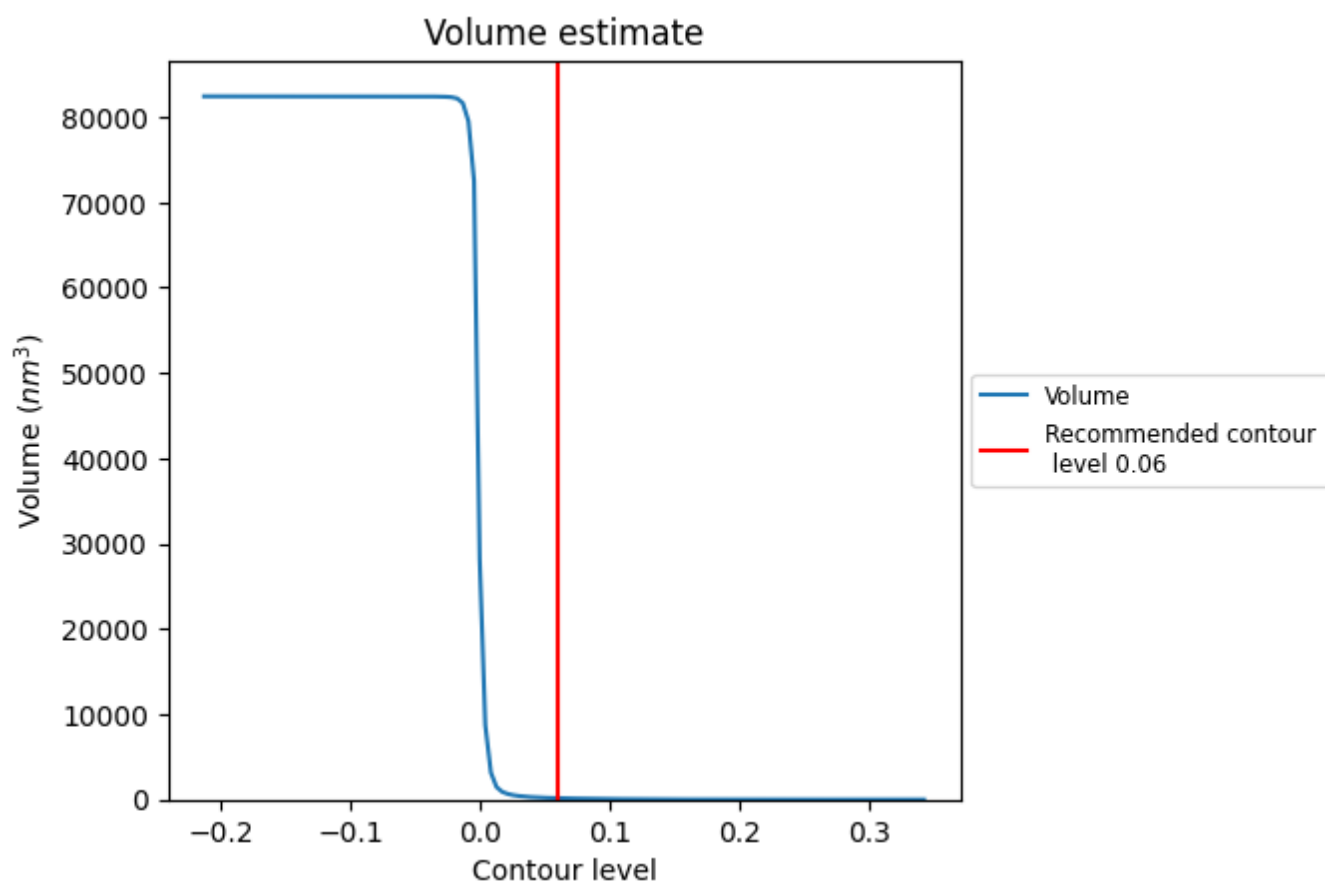
This section contains the results of statistical analysis of the map.

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.

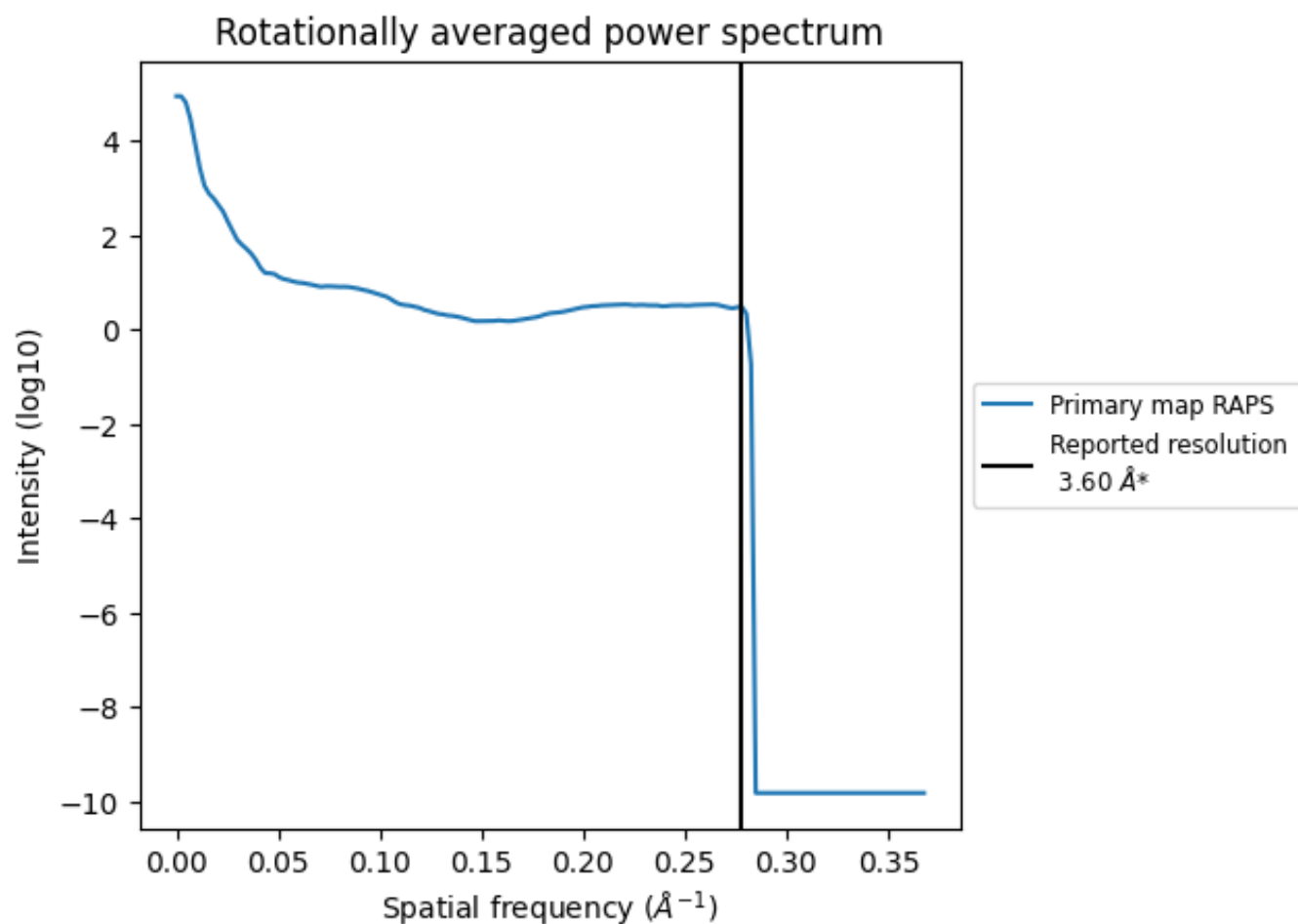
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 164 nm³; this corresponds to an approximate mass of 148 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 0.278 $\text{\AA}^{-1}$

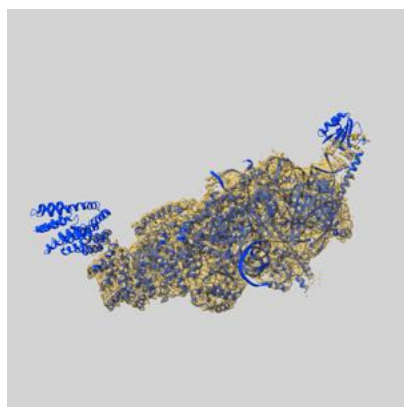
8 Fourier-Shell correlation

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

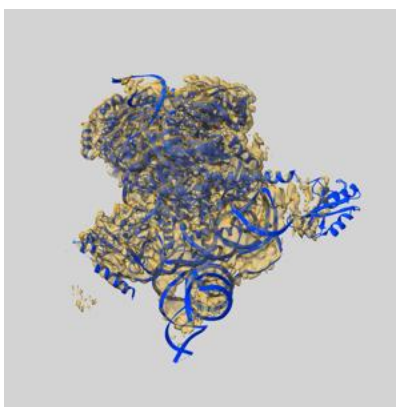
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-8622 and PDB model 6N7X. Per-residue inclusion information can be found in section [3](#) on page [7](#).

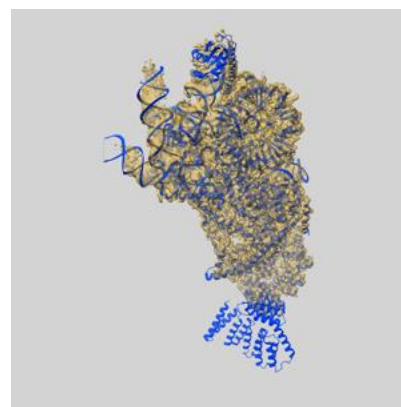
9.1 Map-model overlay [i](#)



X



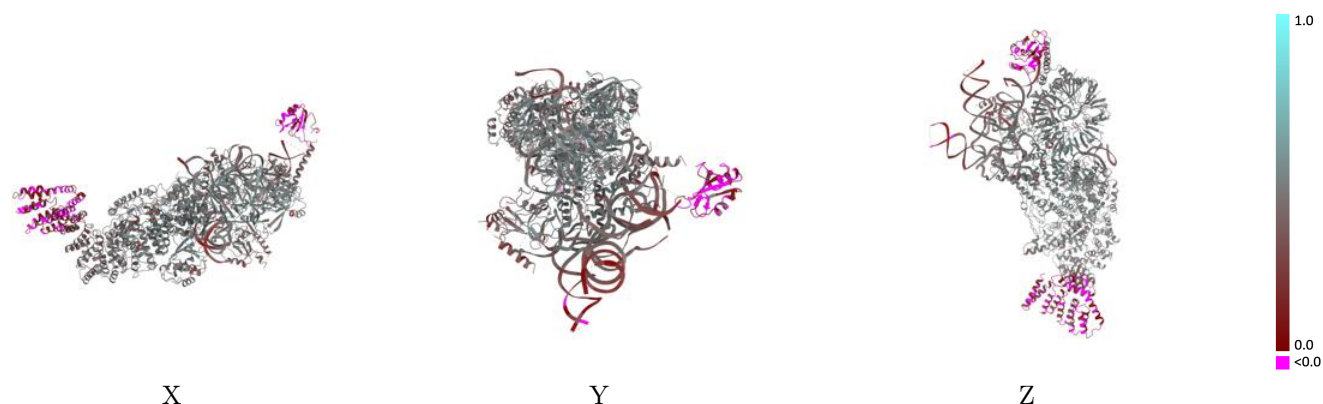
Y



Z

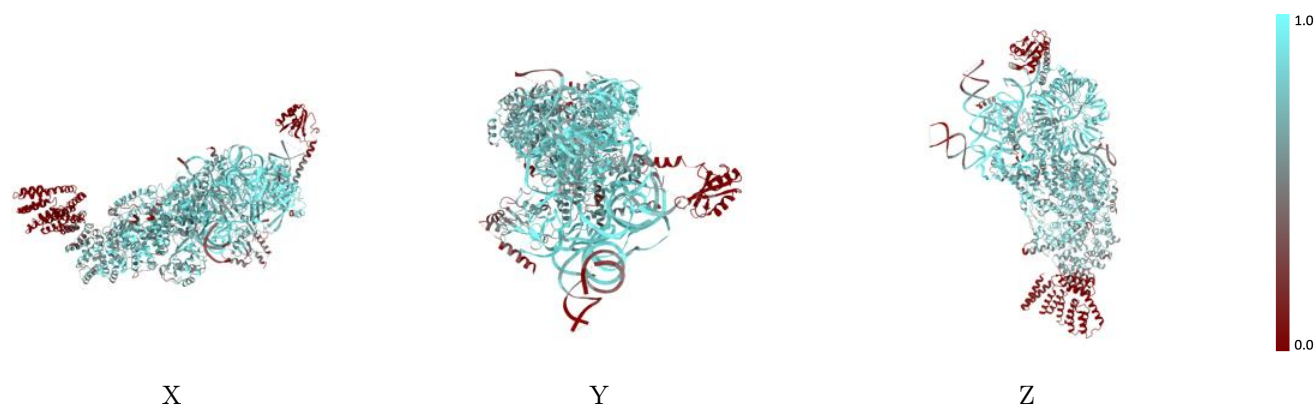
The images above show the 3D surface view of the map at the recommended contour level 0.06 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



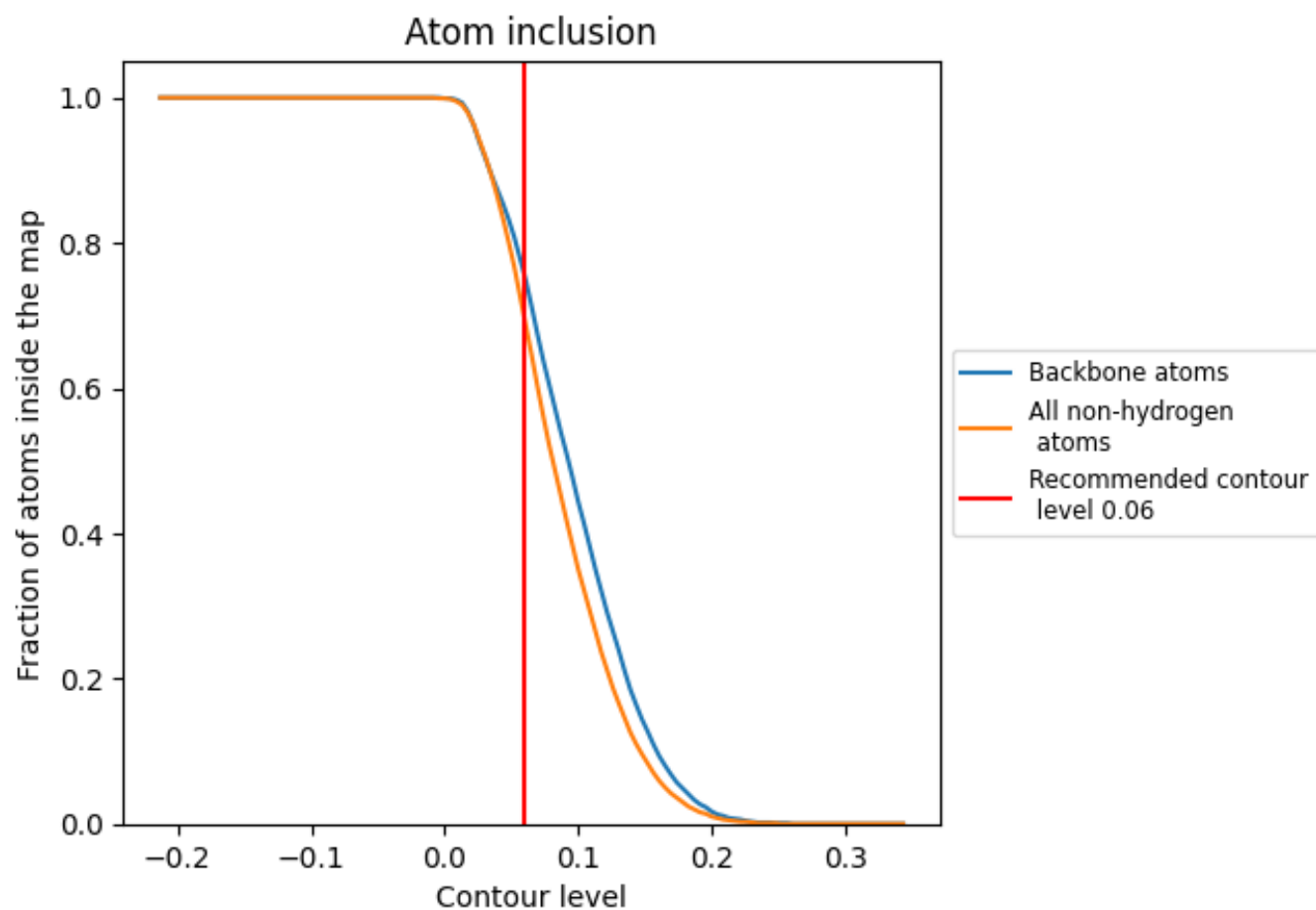
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.06).

9.4 Atom inclusion [i](#)



At the recommended contour level, 75% of all backbone atoms, 69% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.6930	 0.4280
A	 0.4160	 0.2810
B	 0.7180	 0.4670
C	 0.5510	 0.3960
D	 0.7960	 0.4950
E	 0.4310	 0.3250
F	 0.6700	 0.4750
G	 0.7140	 0.4750
H	 0.5700	 0.4320
K	 0.7560	 0.5070
L	 0.7460	 0.4920
M	 0.8120	 0.4890
N	 0.8270	 0.5200
O	 0.8570	 0.4920
P	 0.8820	 0.5010
Q	 0.7780	 0.5020
R	 0.7750	 0.3910

