



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 04:38 AM UTC

PDB ID : 8YHD / pdb_00008yhd
EMDB ID : EMD-39287
Title : Cryo-EM structure of CTR-bound type VII CRISPR-Cas complex at post-state I
Authors : Zhang, H.; Deng, Z.; Li, X.
Deposited on : 2024-02-28
Resolution : 2.93 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:

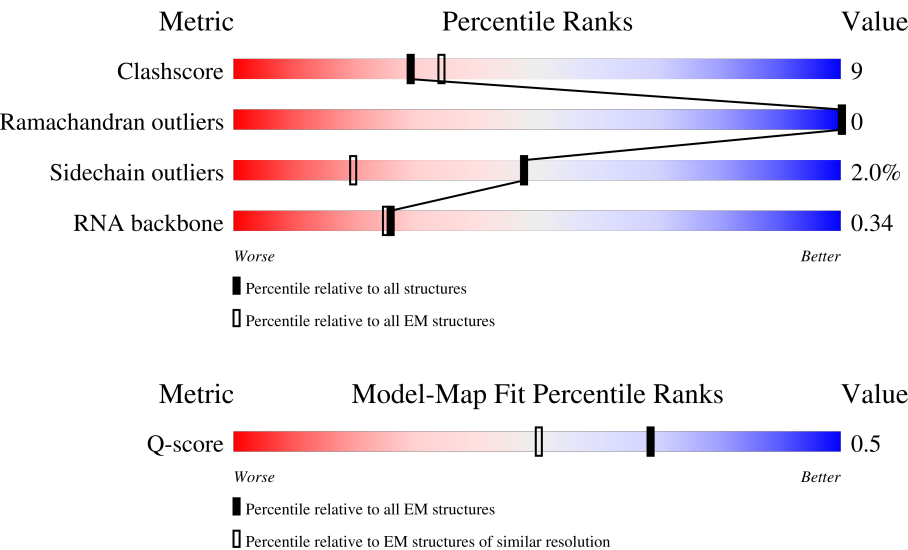
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & 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 2.93 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.



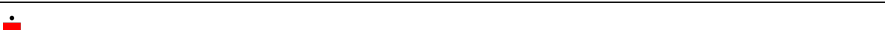
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	13037 (2.43 - 3.43)

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	200	26% 66% 15% 18%
1	B	200	6% 78% 20%
1	C	200	77% 21%

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Mol	Chain	Length	Quality of chain
1	D	200	 77% 22%
1	E	200	 80% 20%
1	F	200	 78% 20% .
1	J	200	 80% 18% .
1	O	200	 70% 24% 5%
2	G	240	 12% 67% 28% . .
3	H	609	 11% 61% 22% . 15%
3	I	609	 21% 5% 74%
3	K	609	 5% 29% 7% 64%
3	L	609	 21% . 74%
4	M	66	 27% 44% 9% 20%
5	N	52	 42% 23% . 33%

2 Entry composition [i](#)

There are 6 unique types of molecules in this entry. The entry contains 23859 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 a protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	163	Total	C	N	O	S	0	0
			1157	724	196	229	8		
1	B	199	Total	C	N	O	S	0	0
			1531	954	274	295	8		
1	C	199	Total	C	N	O	S	0	0
			1547	965	278	296	8		
1	D	199	Total	C	N	O	S	0	0
			1537	960	274	295	8		
1	E	199	Total	C	N	O	S	0	0
			1541	962	275	296	8		
1	F	199	Total	C	N	O	S	0	0
			1523	950	269	296	8		
1	J	199	Total	C	N	O	S	0	0
			1543	962	277	296	8		
1	O	190	Total	C	N	O	S	0	0
			1425	891	246	280	8		

- Molecule 2 is a protein called a protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	G	234	Total	C	N	O	S	0	0
			1703	1086	280	332	5		

- Molecule 3 is a protein called a protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	H	518	Total	C	N	O	S	0	0
			4083	2618	678	773	14		
3	I	160	Total	C	N	O	S	0	0
			1318	846	221	245	6		
3	K	219	Total	C	N	O	S	0	0
			1781	1147	297	334	3		
3	L	160	Total	C	N	O	S	0	0
			1301	835	221	240	5		

- Molecule 4 is a RNA chain called RNA (53-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
4	M	53	Total	C	N	O	P	0	0
			1110	496	184	377	53		

- Molecule 5 is a RNA chain called RNA (35-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
5	N	35	Total	C	N	O	P	0	0
			748	335	139	239	35		

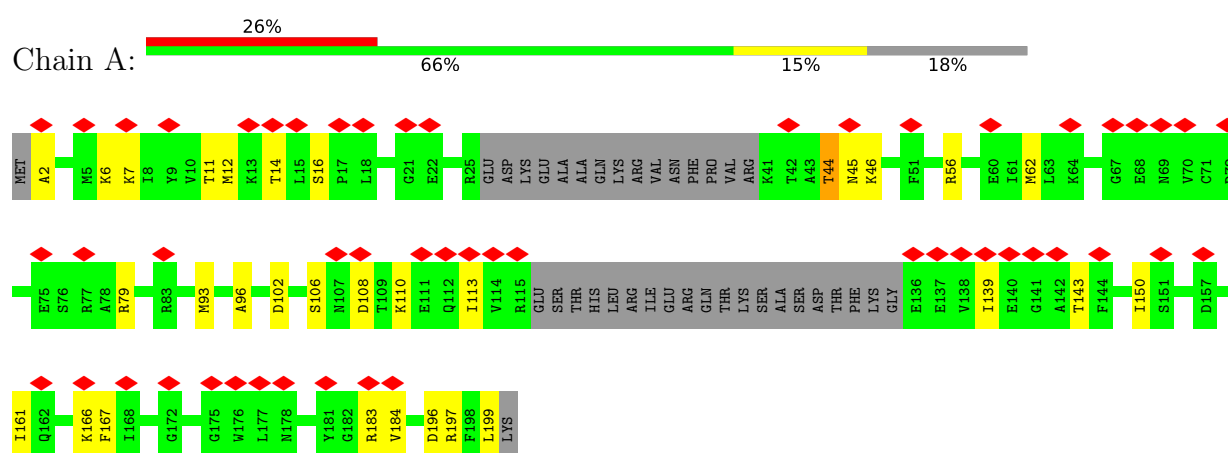
- Molecule 6 is ZINC ION (CCD ID: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		AltConf
6	A	1	Total	Zn	0
			1	1	
6	B	1	Total	Zn	0
			1	1	
6	C	1	Total	Zn	0
			1	1	
6	D	1	Total	Zn	0
			1	1	
6	E	1	Total	Zn	0
			1	1	
6	F	1	Total	Zn	0
			1	1	
6	G	1	Total	Zn	0
			1	1	
6	H	2	Total	Zn	0
			2	2	
6	J	1	Total	Zn	0
			1	1	
6	O	1	Total	Zn	0
			1	1	

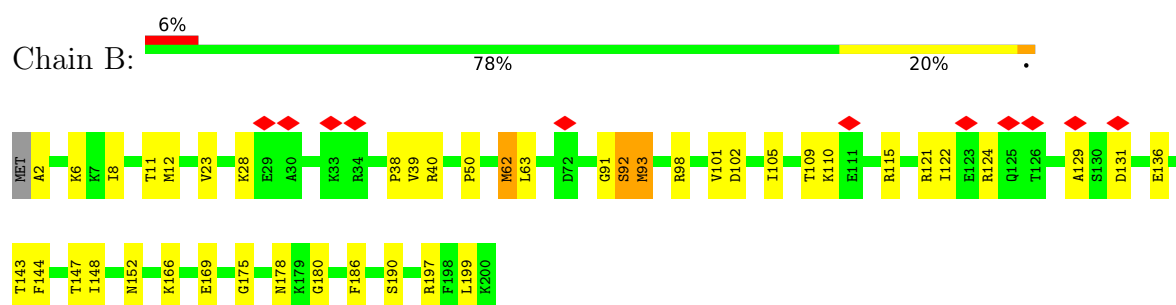
3 Residue-property plots

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

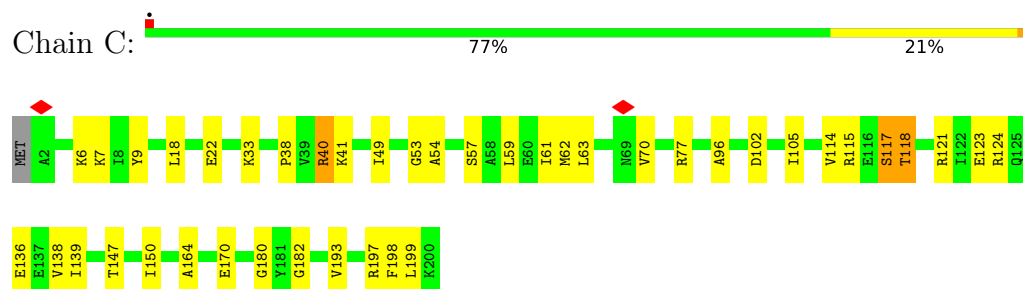
- Molecule 1: a protein



- Molecule 1: a protein

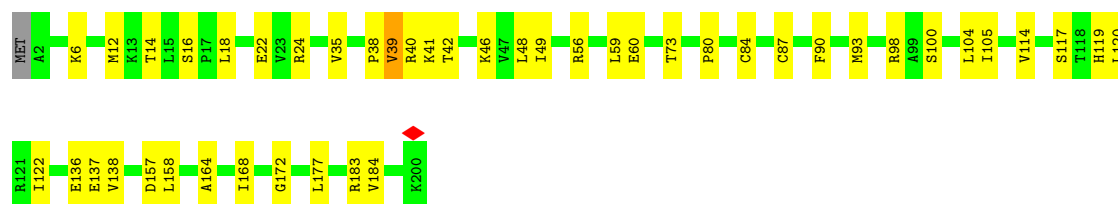


- Molecule 1: a protein



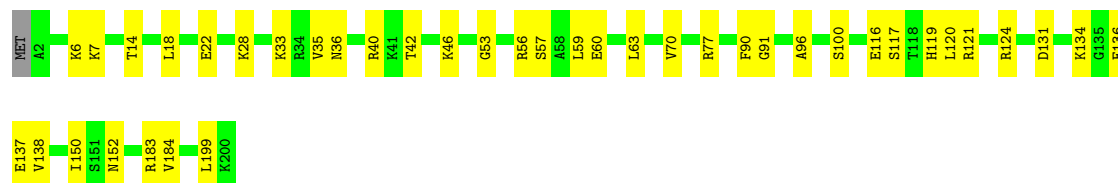
- Molecule 1: a protein

Chain D:  77% 22%



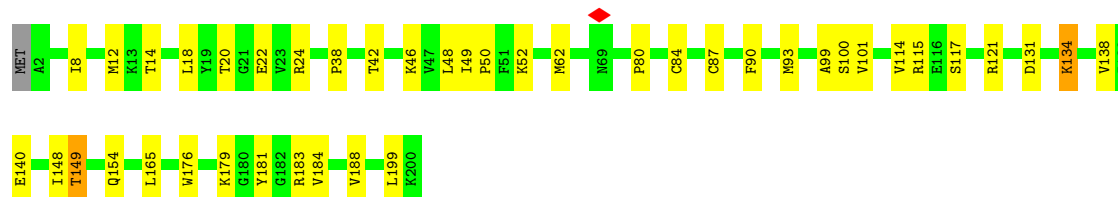
- Molecule 1: a protein

Chain E:  80% 20%



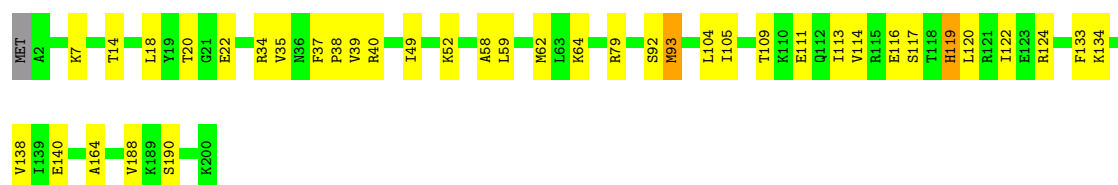
- Molecule 1: a protein

Chain F:  78% 20%



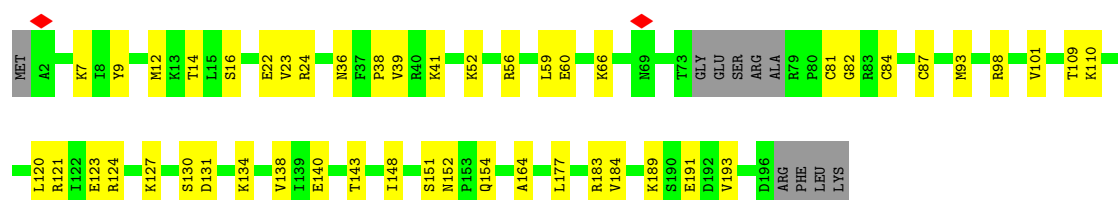
- Molecule 1: a protein

Chain J:  80% 18%

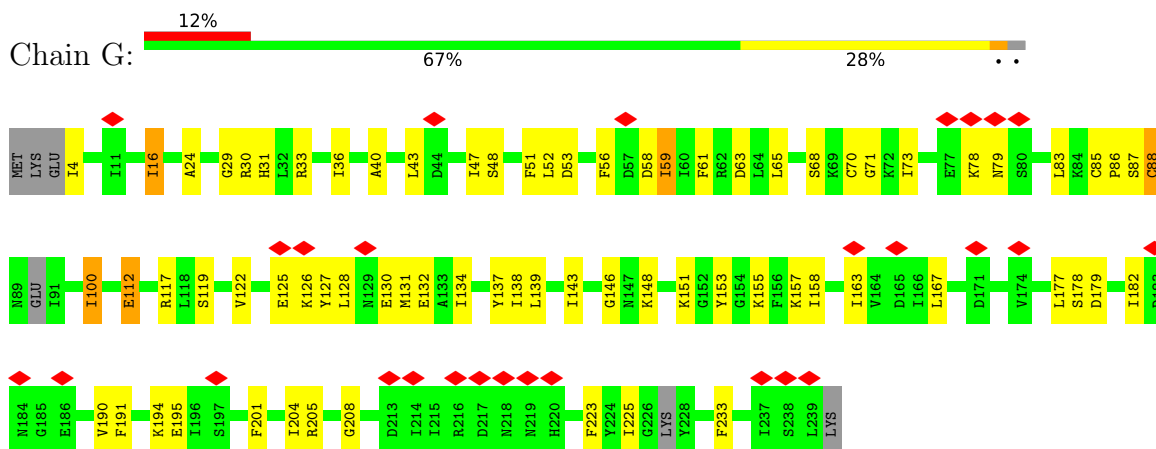


- Molecule 1: a protein

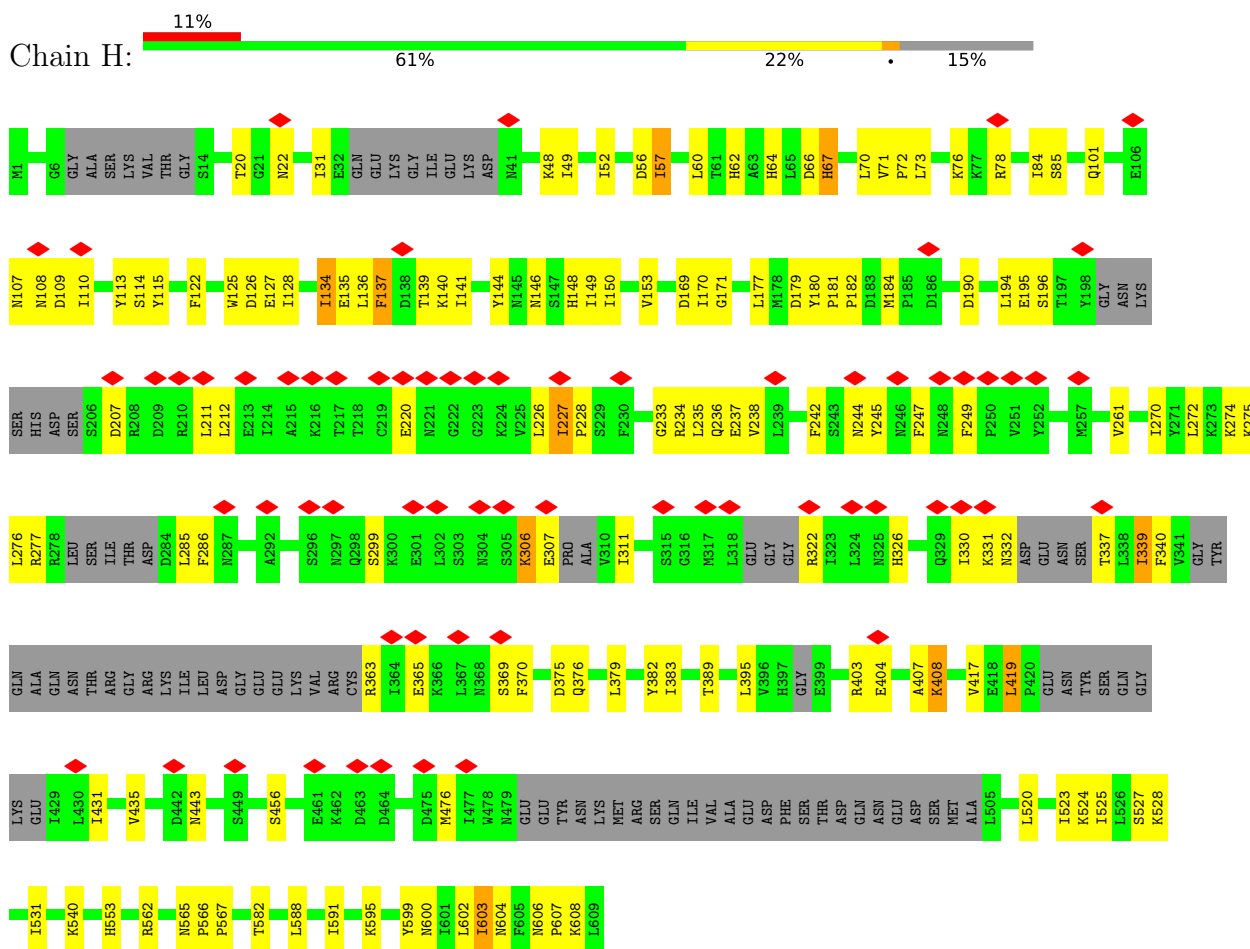
Chain O:  70% 24% 5%



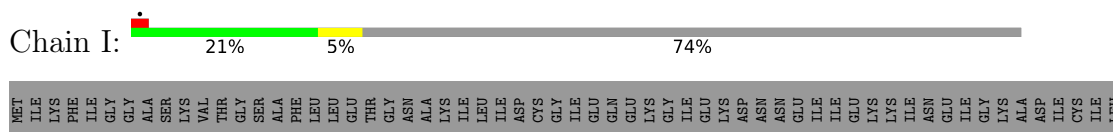
- Molecule 2: a protein

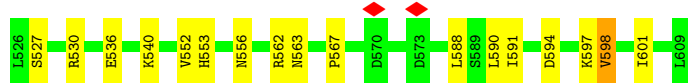


- Molecule 3: a protein

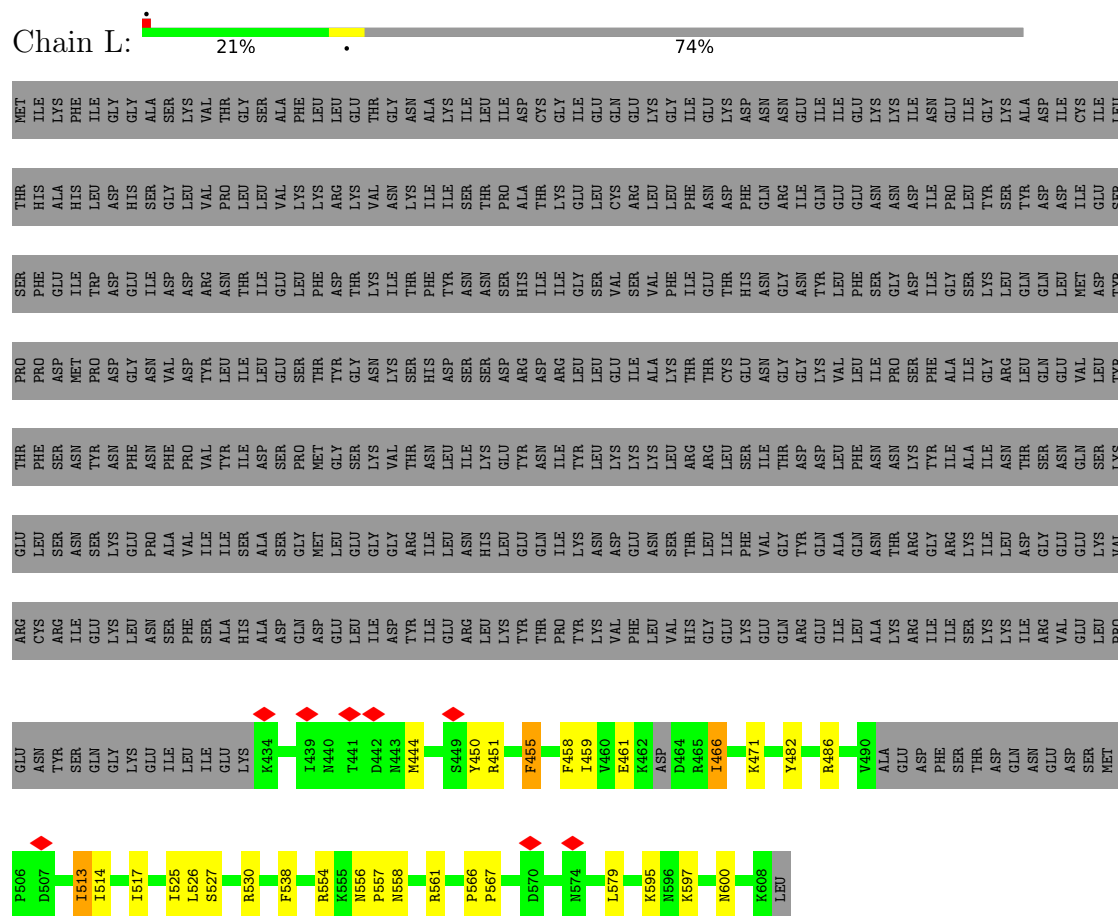


- Molecule 3: a protein

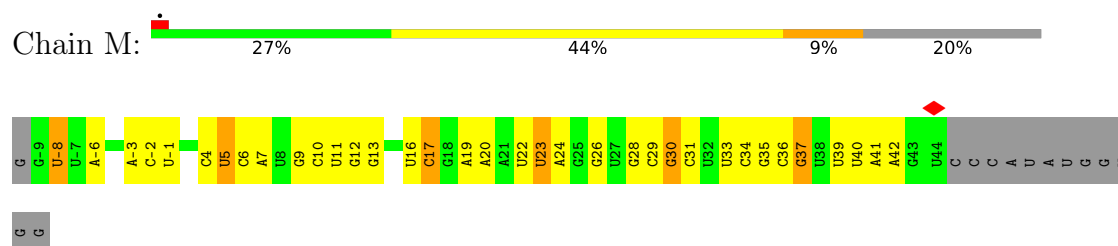




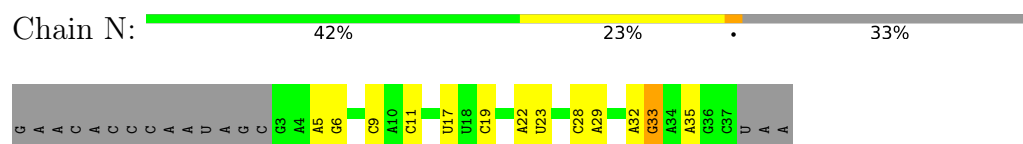
- Molecule 3: a protein



- Molecule 4: RNA (53-MER)



- Molecule 5: RNA (35-MER)



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	100000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING ONLY	Depositor
Microscope	JEOL CRYO ARM 300	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	40	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	1.985	Depositor
Minimum map value	-0.996	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.048	Depositor
Recommended contour level	0.18	Depositor
Map size ($\text{\AA}$)	285.0, 285.0, 285.0	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.95, 0.95, 0.95	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: ZN

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	0.40	1/1169 (0.1%)	0.77	2/1582 (0.1%)
1	B	0.50	4/1550 (0.3%)	0.73	0/2082
1	C	0.59	3/1566 (0.2%)	0.79	2/2099 (0.1%)
1	D	0.44	1/1556 (0.1%)	0.69	0/2087
1	E	0.40	0/1560	0.76	1/2092 (0.0%)
1	F	0.40	0/1542	0.67	1/2073 (0.0%)
1	J	0.40	0/1562	0.74	1/2095 (0.0%)
1	O	0.45	0/1442	0.76	0/1944
2	G	0.64	1/1730 (0.1%)	1.00	5/2337 (0.2%)
3	H	0.62	3/4153 (0.1%)	0.96	6/5616 (0.1%)
3	I	0.44	1/1348 (0.1%)	0.76	2/1824 (0.1%)
3	K	0.39	2/1817 (0.1%)	0.73	3/2459 (0.1%)
3	L	0.60	2/1331 (0.2%)	0.86	0/1805
4	M	0.32	0/1237	0.57	0/1923
5	N	0.25	0/837	0.41	0/1302
All	All	0.50	18/24400 (0.1%)	0.79	23/33320 (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	B	0	1
3	H	0	1
3	K	0	1
3	L	0	1
All	All	0	4

The worst 5 of 18 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	C	40	ARG	C-N	-7.94	1.22	1.33
2	G	40	ALA	C-N	7.54	1.44	1.33
3	H	407	ALA	C-N	-7.22	1.24	1.33
1	C	117	SER	C-N	7.13	1.43	1.33
1	B	40	ARG	C-N	-6.93	1.23	1.33

The worst 5 of 23 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	H	67	HIS	N-CA-C	-7.47	103.74	114.12
1	C	54	ALA	N-CA-C	-6.94	104.97	113.50
1	C	53	GLY	N-CA-C	-6.72	106.39	113.58
2	G	86	PRO	N-CA-C	-6.49	104.40	113.81
1	E	33	LYS	N-CA-C	-6.41	104.51	112.90

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	B	115	ARG	Sidechain
3	H	78	ARG	Sidechain
3	K	486	ARG	Sidechain
3	L	554	ARG	Sidechain

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	1157	0	1086	19	0
1	B	1531	0	1547	31	0
1	C	1547	0	1586	32	0
1	D	1537	0	1569	25	0
1	E	1541	0	1575	30	0
1	F	1523	0	1531	28	0
1	J	1543	0	1575	24	0
1	O	1425	0	1411	35	0
2	G	1703	0	1574	64	0
3	H	4083	0	3933	100	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	I	1318	0	1249	25	0
3	K	1781	0	1688	30	0
3	L	1301	0	1214	22	0
4	M	1110	0	561	25	0
5	N	748	0	380	12	0
6	A	1	0	0	0	0
6	B	1	0	0	0	0
6	C	1	0	0	0	0
6	D	1	0	0	0	0
6	E	1	0	0	0	0
6	F	1	0	0	0	0
6	G	1	0	0	1	0
6	H	2	0	0	0	0
6	J	1	0	0	0	0
6	O	1	0	0	0	0
All	All	23859	0	22479	431	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.

The worst 5 of 431 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:G:56:PHE:CZ	2:G:58:ASP:HB3	1.93	1.04
2:G:56:PHE:CE2	2:G:58:ASP:HB3	1.95	1.02
3:L:444:MET:HG2	3:L:451:ARG:HG2	1.52	0.90
3:K:192:LEU:HB3	3:K:393:VAL:HG12	1.54	0.89
2:G:56:PHE:CE2	2:G:59:ILE:HG12	2.08	0.87

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	157/200 (78%)	147 (94%)	10 (6%)	0	100	100
1	B	197/200 (98%)	193 (98%)	4 (2%)	0	100	100
1	C	197/200 (98%)	191 (97%)	6 (3%)	0	100	100
1	D	197/200 (98%)	191 (97%)	6 (3%)	0	100	100
1	E	197/200 (98%)	192 (98%)	5 (2%)	0	100	100
1	F	197/200 (98%)	190 (96%)	7 (4%)	0	100	100
1	J	197/200 (98%)	195 (99%)	2 (1%)	0	100	100
1	O	186/200 (93%)	181 (97%)	5 (3%)	0	100	100
2	G	228/240 (95%)	216 (95%)	12 (5%)	0	100	100
3	H	494/609 (81%)	459 (93%)	35 (7%)	0	100	100
3	I	154/609 (25%)	149 (97%)	5 (3%)	0	100	100
3	K	207/609 (34%)	196 (95%)	11 (5%)	0	100	100
3	L	154/609 (25%)	148 (96%)	6 (4%)	0	100	100
All	All	2762/4276 (65%)	2648 (96%)	114 (4%)	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	A	114/170 (67%)	113 (99%)	1 (1%)	70	82
1	B	165/170 (97%)	162 (98%)	3 (2%)	51	72
1	C	169/170 (99%)	167 (99%)	2 (1%)	63	78
1	D	167/170 (98%)	164 (98%)	3 (2%)	51	72
1	E	168/170 (99%)	167 (99%)	1 (1%)	78	89
1	F	164/170 (96%)	161 (98%)	3 (2%)	51	72
1	J	168/170 (99%)	164 (98%)	4 (2%)	43	67
1	O	152/170 (89%)	151 (99%)	1 (1%)	76	86

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	G	164/209 (78%)	158 (96%)	6 (4%)	30	55
3	H	437/558 (78%)	422 (97%)	15 (3%)	32	58
3	I	141/558 (25%)	138 (98%)	3 (2%)	47	70
3	K	187/558 (34%)	186 (100%)	1 (0%)	81	90
3	L	136/558 (24%)	132 (97%)	4 (3%)	37	62
All	All	2332/3801 (61%)	2285 (98%)	47 (2%)	48	70

5 of 47 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	H	244	ASN
3	I	606	ASN
3	H	339	ILE
3	H	476	MET
1	J	52	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 22 such sidechains are listed below:

Mol	Chain	Res	Type
3	K	585	ASN
3	L	556	ASN
3	L	472	ASN
3	L	600	ASN
3	H	161	ASN

5.3.3 RNA ⓘ

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	M	52/66 (78%)	16 (30%)	2 (3%)
5	N	34/52 (65%)	3 (8%)	0
All	All	86/118 (72%)	19 (22%)	2 (2%)

5 of 19 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	M	-8	U
4	M	-3	A
4	M	-2	C

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Mol	Chain	Res	Type
4	M	-1	U
4	M	5	U

All (2) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	M	23	U
4	M	30	G

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](#)

Of 11 ligands modelled in this entry, 11 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.

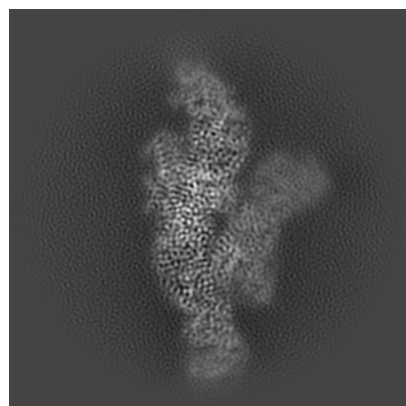
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-39287. These allow visual inspection of the internal detail of the map and identification of artifacts.

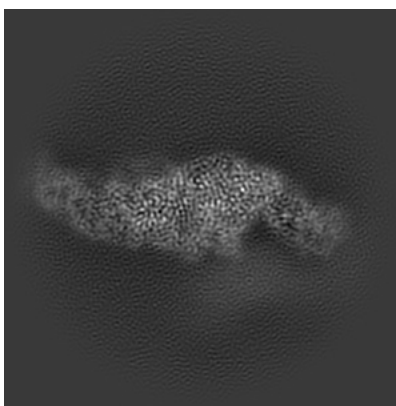
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

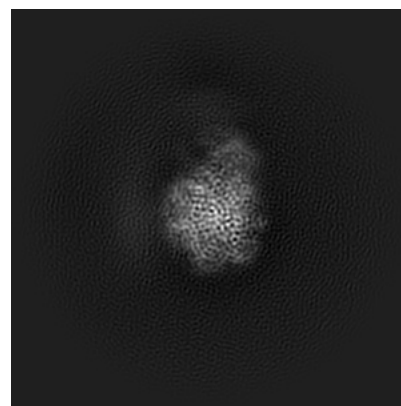
6.1.1 Primary map



X

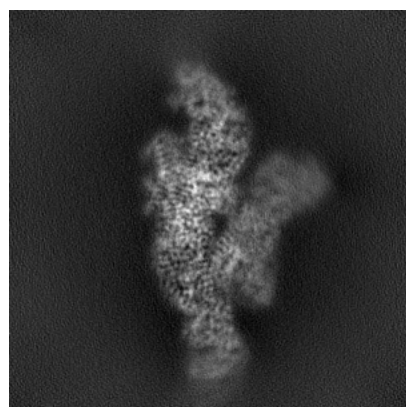


Y

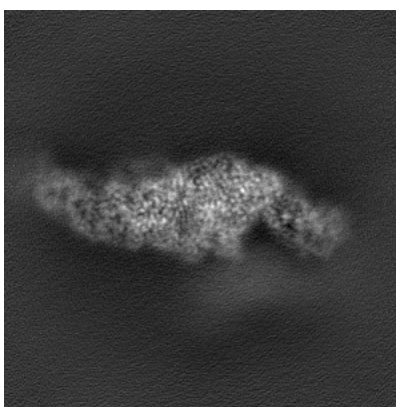


Z

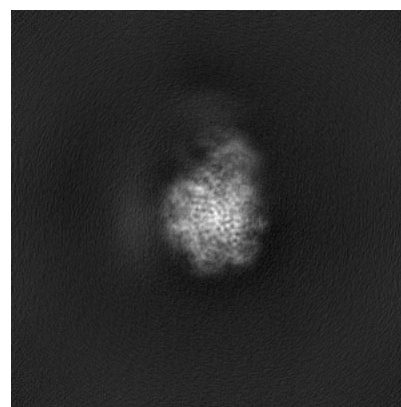
6.1.2 Raw map



X



Y

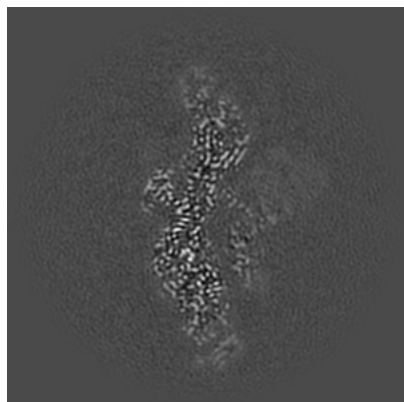


Z

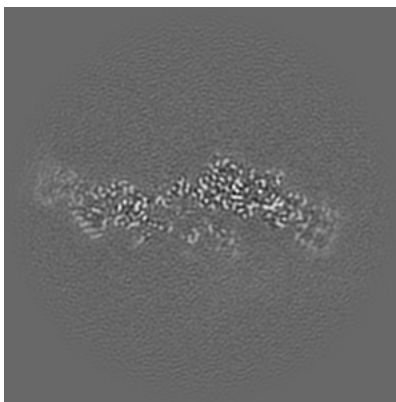
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

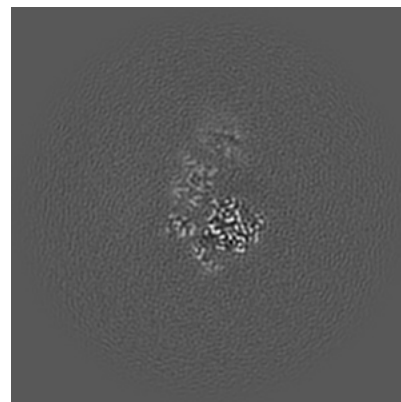
6.2.1 Primary map



X Index: 150

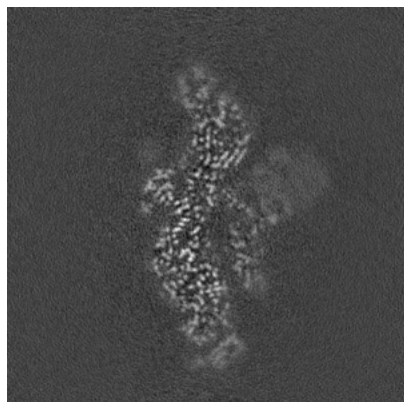


Y Index: 150

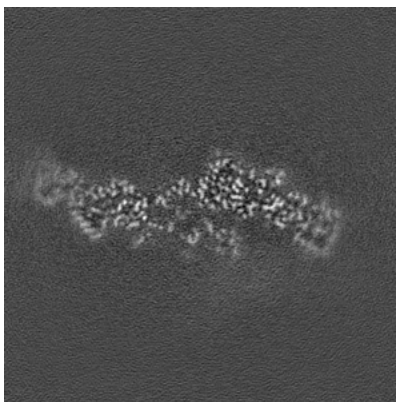


Z Index: 150

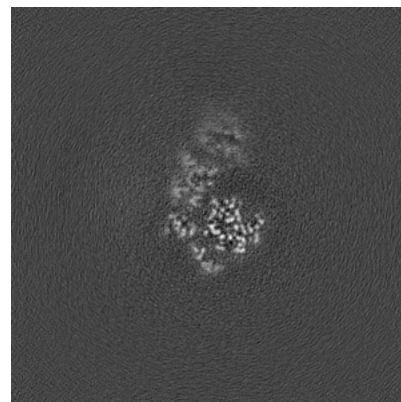
6.2.2 Raw map



X Index: 150



Y Index: 150

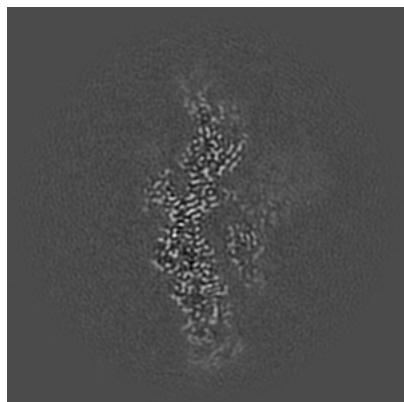


Z Index: 150

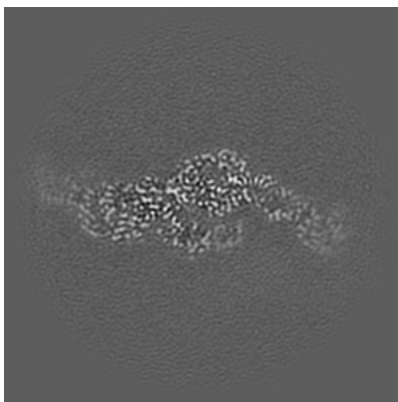
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

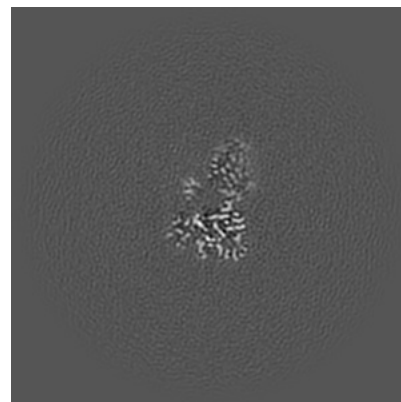
6.3.1 Primary map



X Index: 152

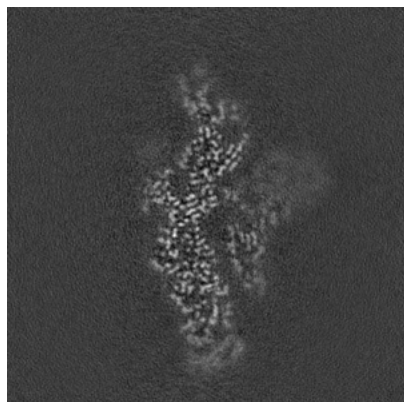


Y Index: 138

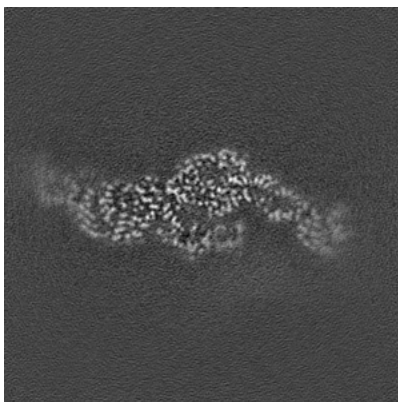


Z Index: 127

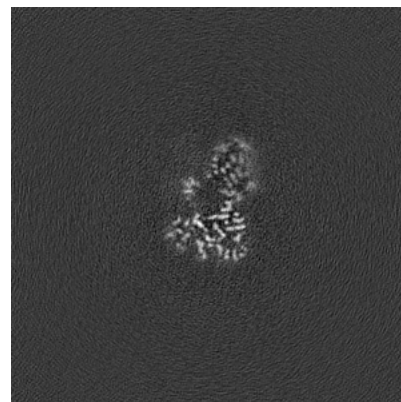
6.3.2 Raw map



X Index: 152



Y Index: 138

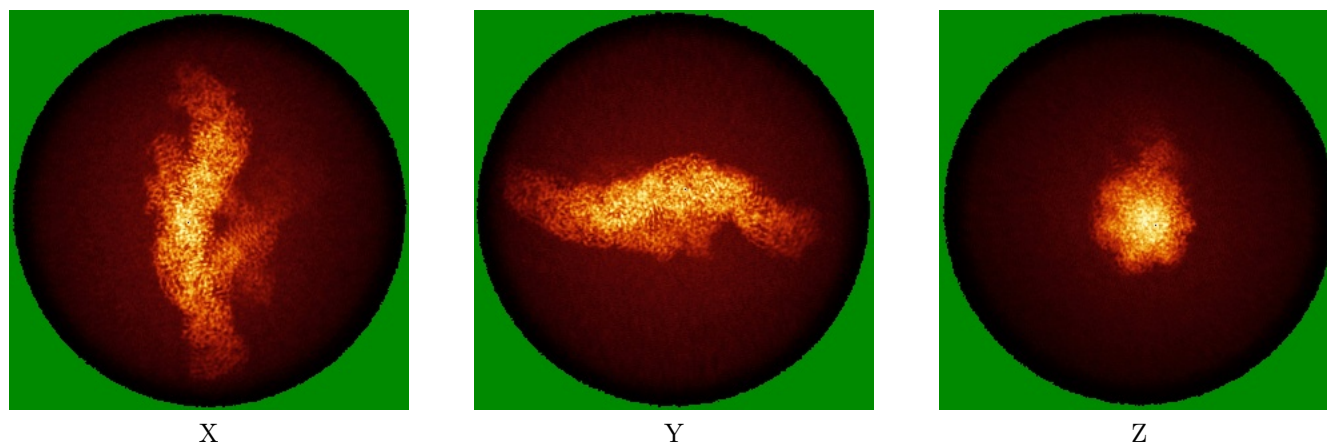


Z Index: 127

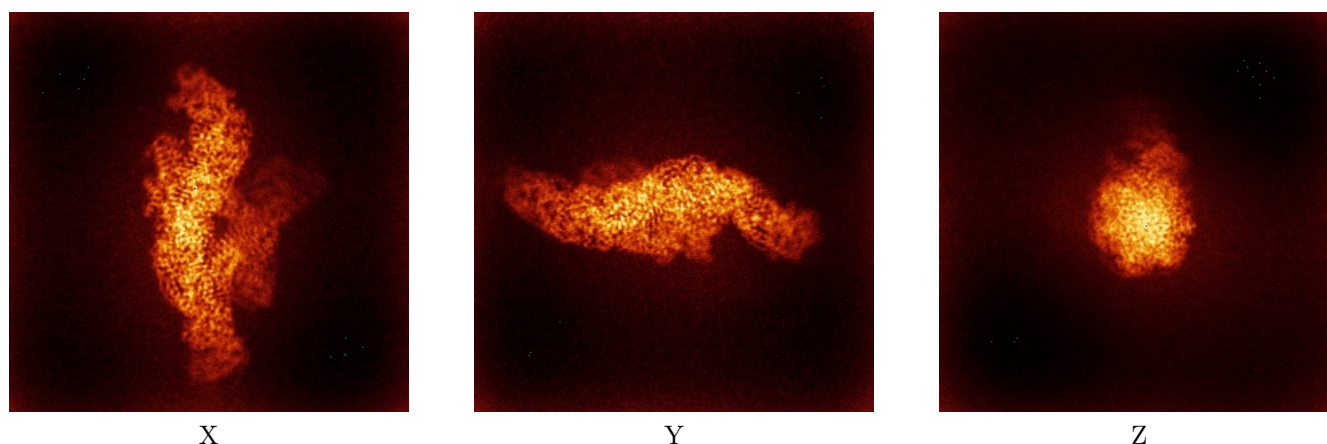
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map



6.4.2 Raw map



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](#)

This section was not generated.

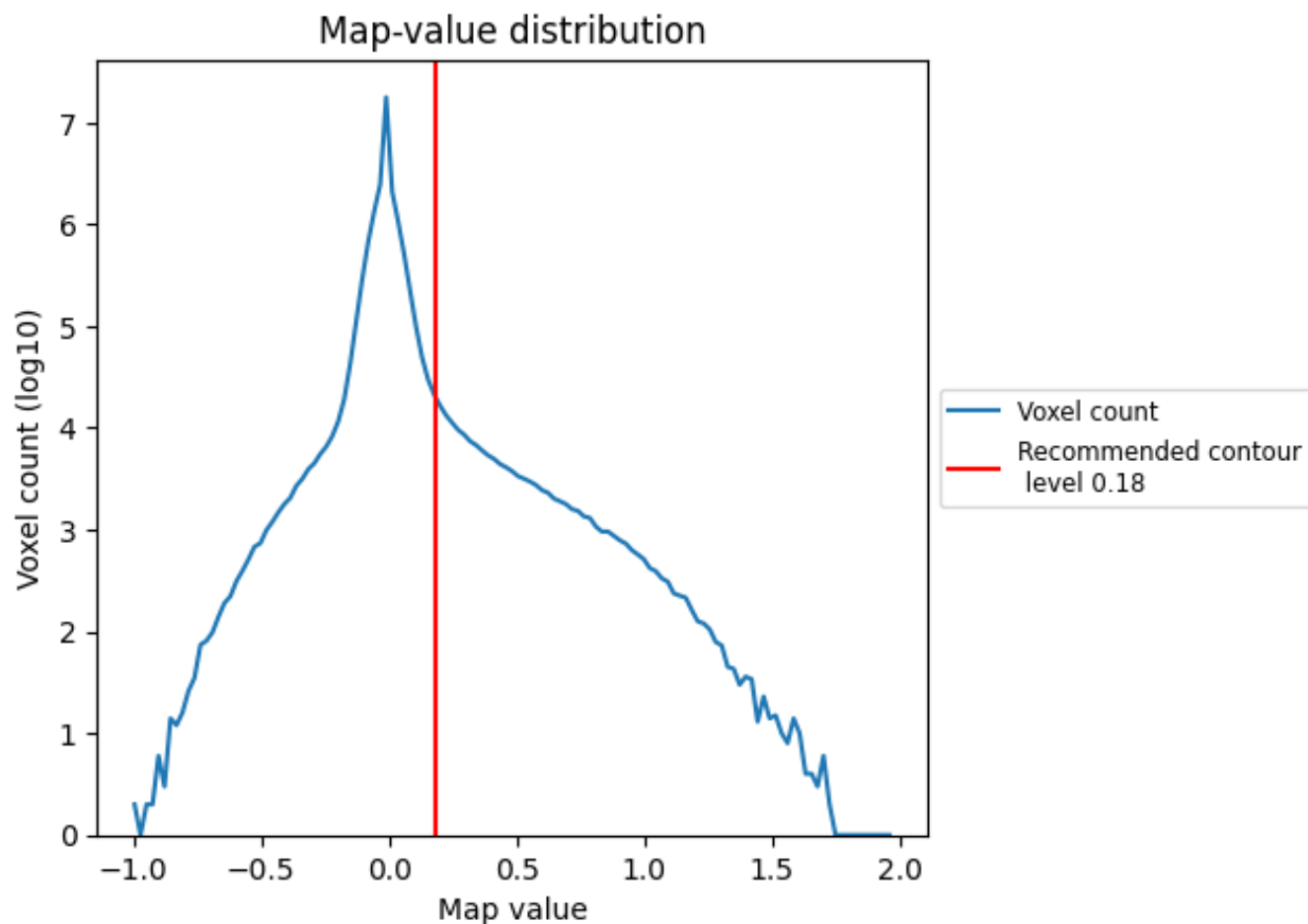
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

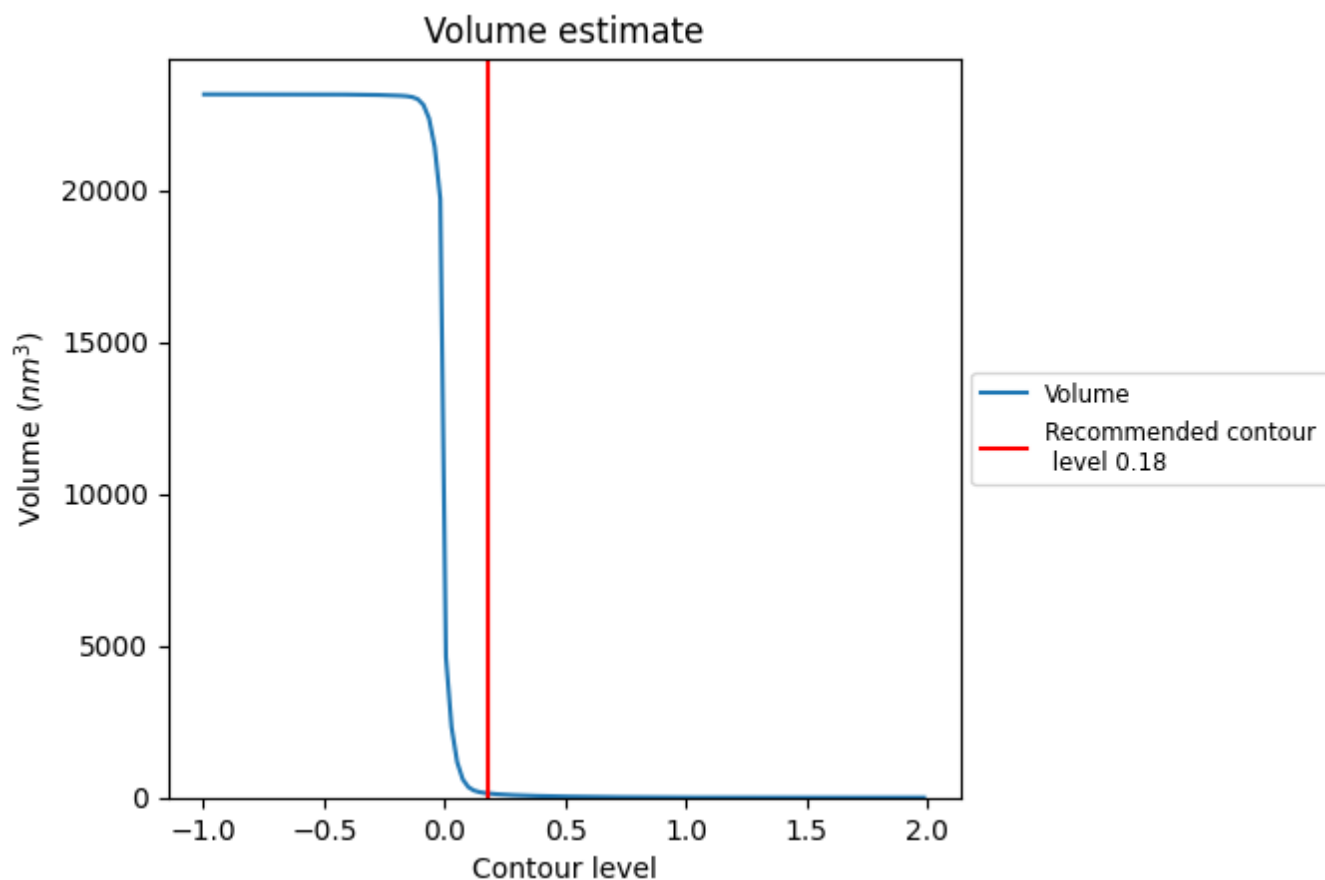
This section contains the results of statistical analysis of the map.

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.

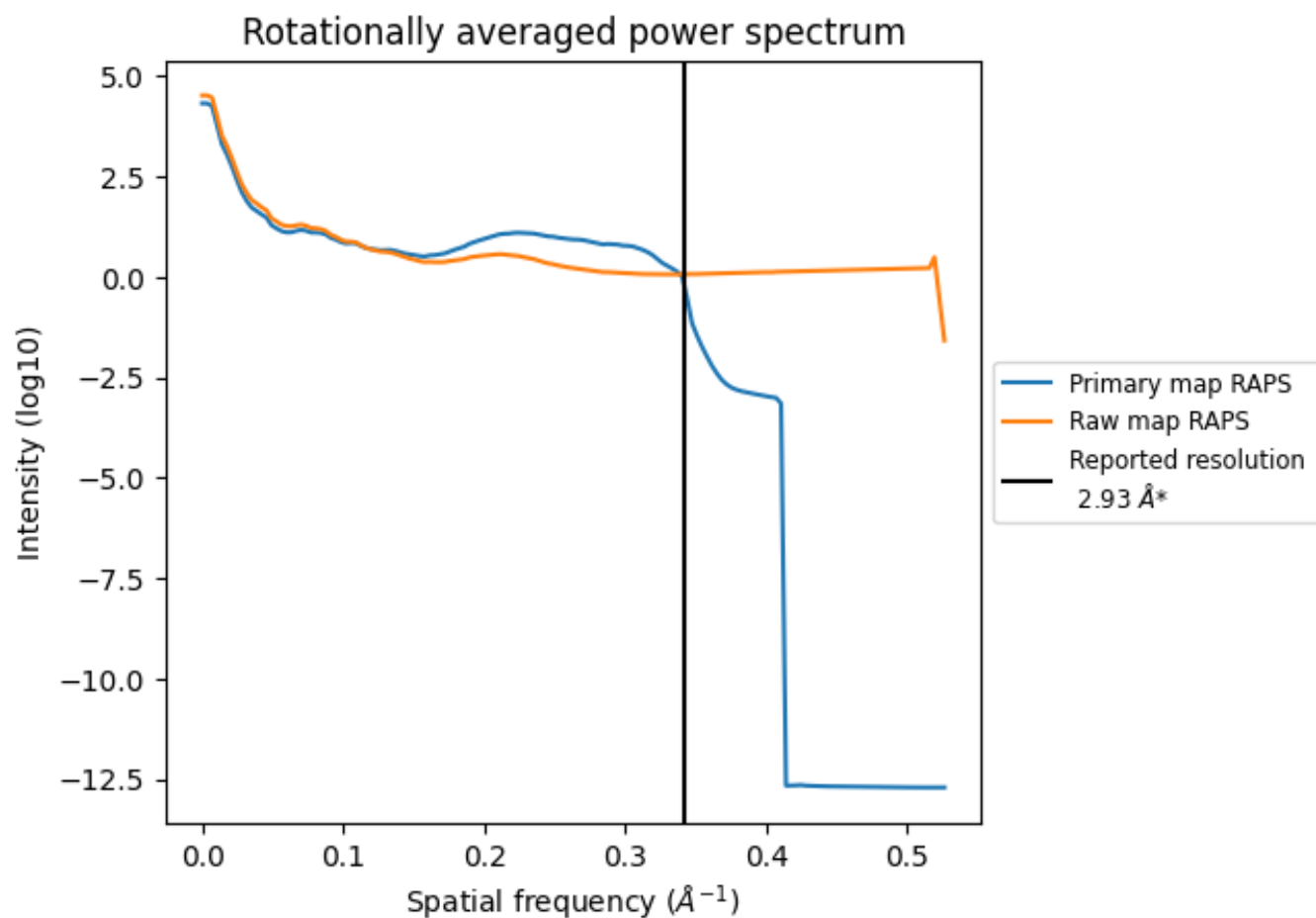
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 138 nm³; this corresponds to an approximate mass of 124 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 ⓘ

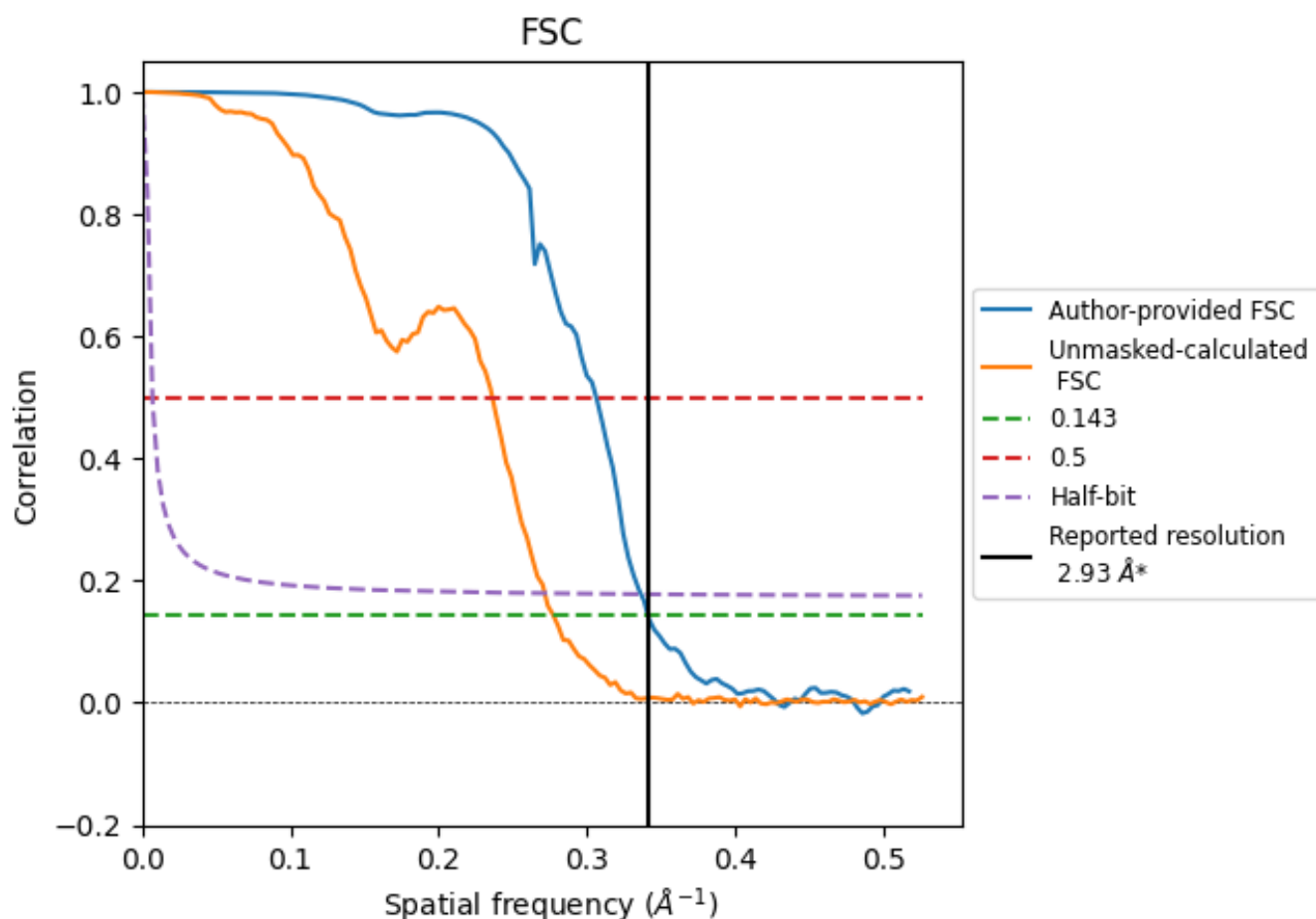


*Reported resolution corresponds to spatial frequency of 0.341 Å⁻¹

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 0.341 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.93	-	-
Author-provided FSC curve	2.93	3.26	2.97
Unmasked-calculated*	3.60	4.23	3.68

*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 3.60 differs from the reported value 2.93 by more than 10 %

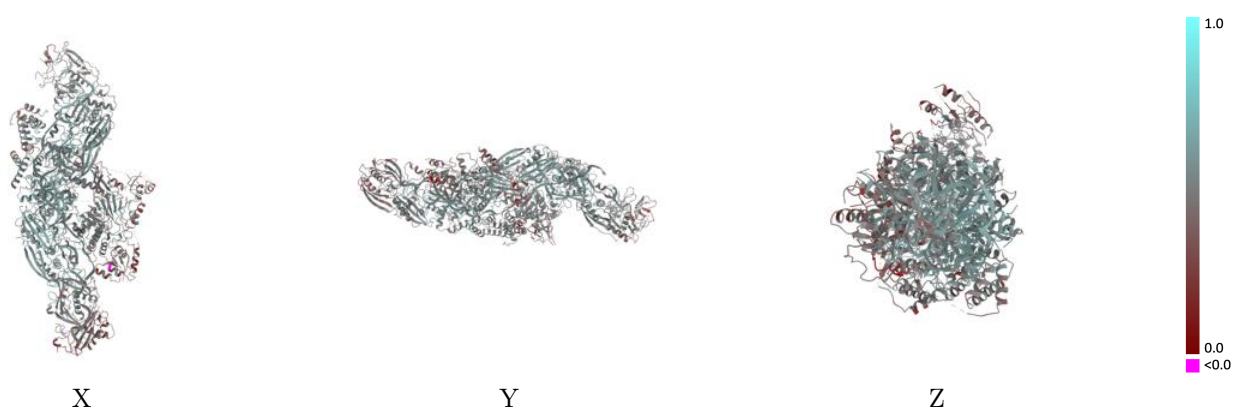
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-39287 and PDB model 8YHD. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)

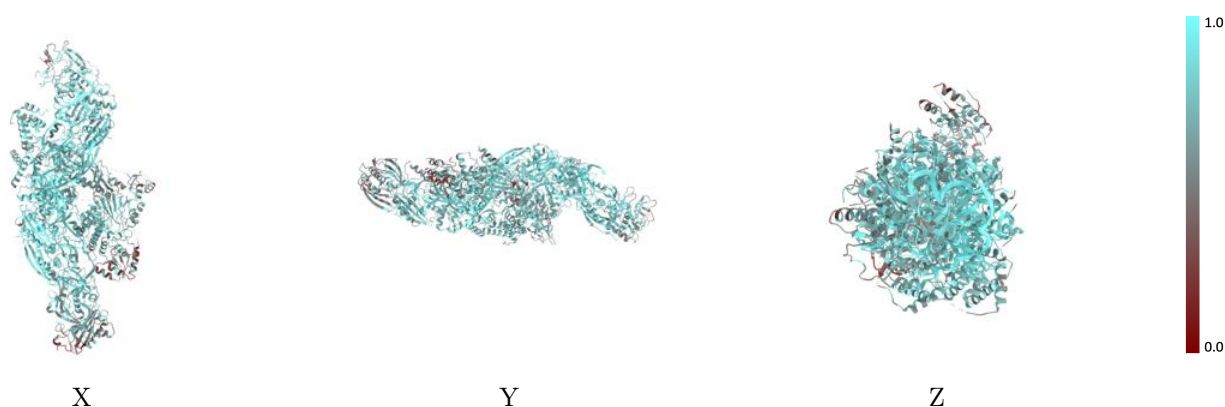
This section was not generated.

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



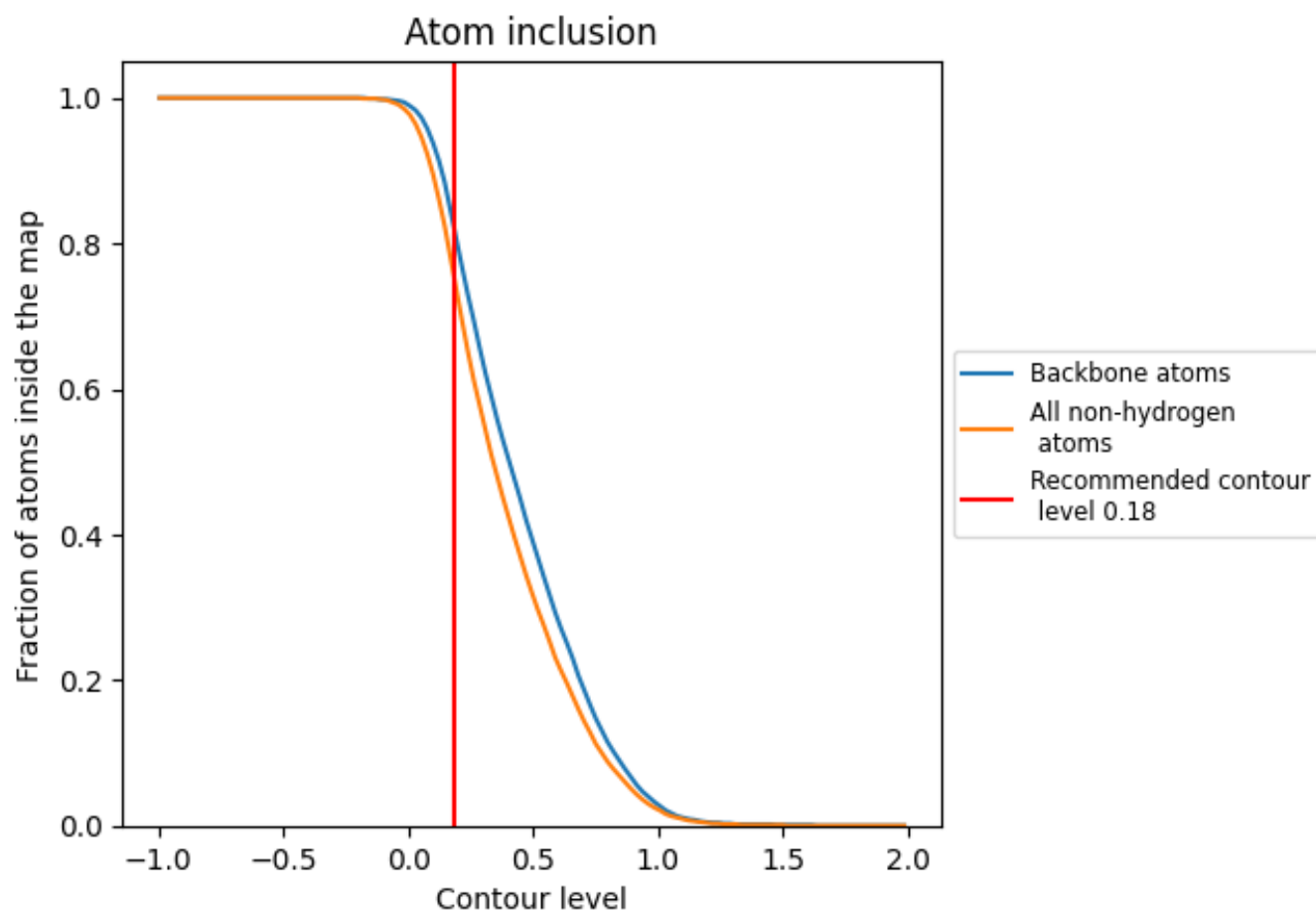
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.18).

9.4 Atom inclusion [i](#)



At the recommended contour level, 82% of all backbone atoms, 76% 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 (0.18) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7580	 0.5000
A	 0.5130	 0.3880
B	 0.7360	 0.4940
C	 0.8270	 0.5410
D	 0.8590	 0.5540
E	 0.8650	 0.5650
F	 0.8310	 0.5520
G	 0.6570	 0.4390
H	 0.6690	 0.4590
I	 0.7460	 0.4810
J	 0.8510	 0.5580
K	 0.6830	 0.4490
L	 0.7120	 0.4570
M	 0.9130	 0.5630
N	 0.9440	 0.5650
O	 0.8000	 0.5250

