



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 02:28 PM UTC

PDB ID : 3J8C / pdb_00003j8c
EMDB ID : EMD-5658
Title : Model of the human eIF3 PCI-MPN octamer docked into the 43S EM map
Authors : Erzberger, J.P.; Ban, N.
Deposited on : 2014-10-08
Resolution : 11.60 Å (reported)
Based on initial models : 3CHM, 3J47, 4O8X, 4LCT, 4B4T, 4U1C, 4U1D, 1RZ4

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:

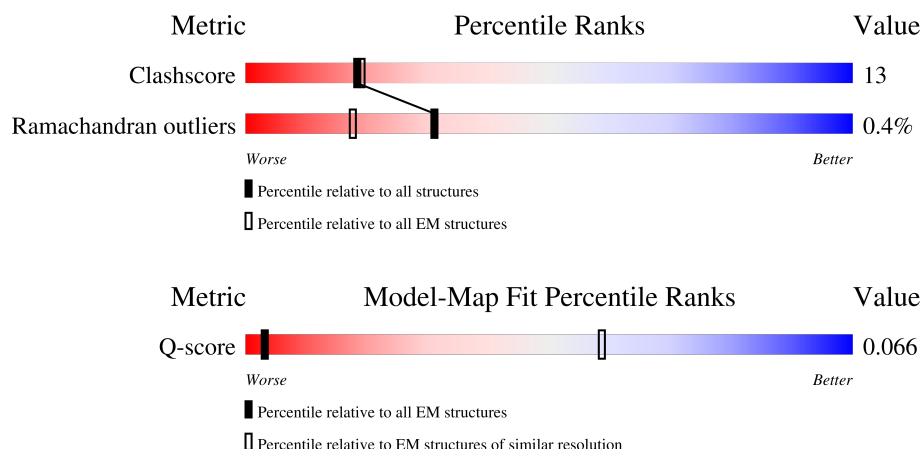
EMDB validation analysis : 0.0.1.dev132
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDb archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh &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

1 Overall quality at a glance

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

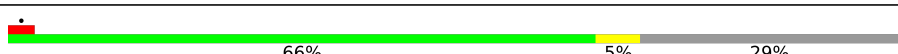
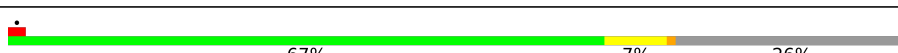
The reported resolution of this entry is 11.60 Å.

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	-
Q-score	-	25397	94 (11.10 - 12.10)

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	529	
2	C	547	
3	E	422	
4	F	327	
5	H	343	

Continued on next page...

Mol	Chain	Length	Quality of chain
6	K	208	
7	L	540	
8	M	364	

2 Entry composition [i](#)

There are 8 unique types of molecules in this entry. The entry contains 12633 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 Eukaryotic translation initiation factor 3 subunit A.

Mol	Chain	Residues	Atoms				AltConf	Trace
1	A	488	Total	C	N	O	0	0
			2432	1456	488	488		

- Molecule 2 is a protein called Eukaryotic translation initiation factor 3 subunit C.

Mol	Chain	Residues	Atoms				AltConf	Trace
2	C	546	Total	C	N	O	0	0
			2711	1619	546	546		

- Molecule 3 is a protein called Eukaryotic translation initiation factor 3 subunit E.

Mol	Chain	Residues	Atoms				AltConf	Trace
3	E	384	Total	C	N	O	0	0
			1911	1143	384	384		

- Molecule 4 is a protein called Eukaryotic translation initiation factor 3 subunit F.

Mol	Chain	Residues	Atoms				AltConf	Trace
4	F	232	Total	C	N	O	0	0
			1149	685	232	232		

- Molecule 5 is a protein called Eukaryotic translation initiation factor 3 subunit H.

Mol	Chain	Residues	Atoms				AltConf	Trace
5	H	254	Total	C	N	O	0	0
			1262	754	254	254		

- Molecule 6 is a protein called Eukaryotic translation initiation factor 3 subunit K.

Mol	Chain	Residues	Atoms				AltConf	Trace
6	K	204	Total	C	N	O	0	0
			1012	604	204	204		

- Molecule 7 is a protein called Eukaryotic translation initiation factor 3 subunit L.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	L	258	Total	C	N	O	0	0
			1281	765	258	258		

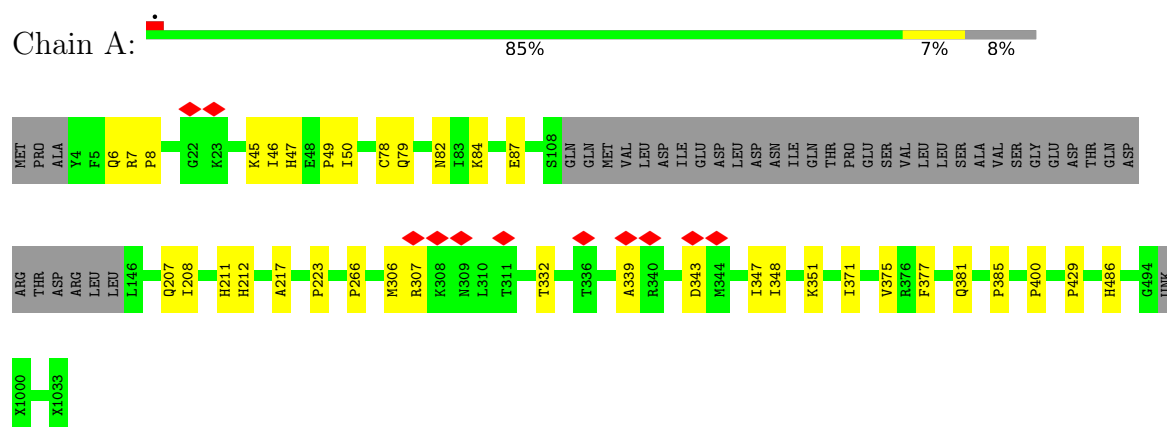
- Molecule 8 is a protein called Eukaryotic translation initiation factor 3 subunit M.

Mol	Chain	Residues	Atoms				AltConf	Trace
8	M	176	Total	C	N	O	0	0
			875	523	176	176		

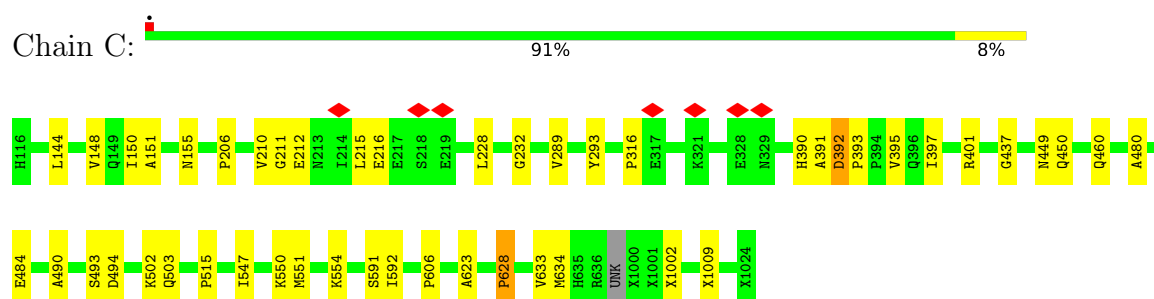
3 Residue-property plots [i](#)

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.

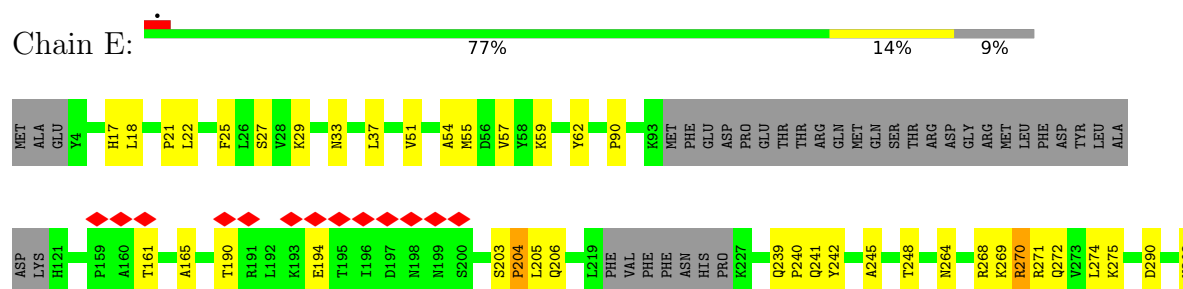
- Molecule 1: Eukaryotic translation initiation factor 3 subunit A

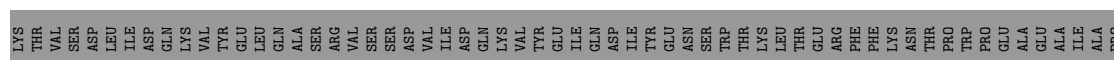


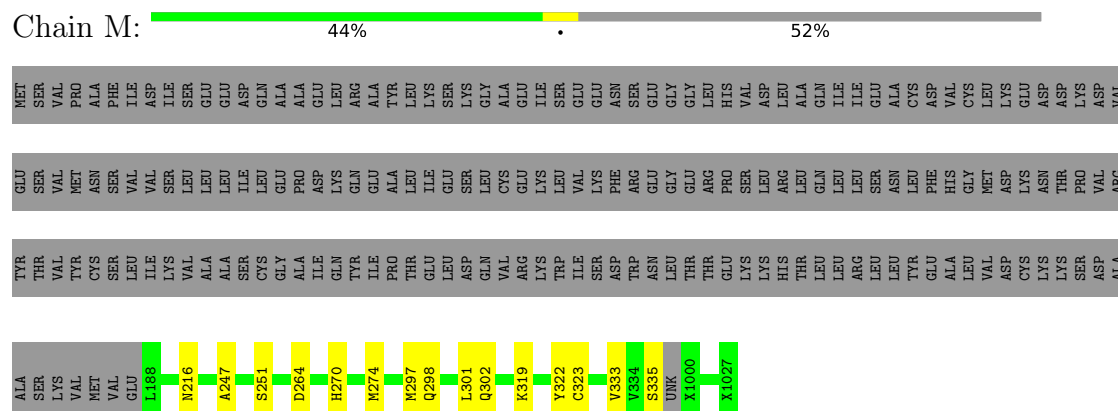
- Molecule 2: Eukaryotic translation initiation factor 3 subunit C



- Molecule 3: Eukaryotic translation initiation factor 3 subunit E







4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	29000	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	Not provided	
Microscope	FEI TECNAI 20	Depositor
Voltage (kV)	110	Depositor
Electron dose ($e^-/\text{\AA}^2$)	12	Depositor
Minimum defocus (nm)	-1000	Depositor
Maximum defocus (nm)	-4000	Depositor
Magnification	Not provided	
Image detector	GATAN ULTRASCAN 4000 (4k x 4k)	Depositor
Maximum map value	0.116	Depositor
Minimum map value	-0.027	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.009	Depositor
Recommended contour level	0.019	Depositor
Map size ($\text{\AA}$)	525.32996, 525.32996, 525.32996	wwPDB
Map dimensions	234, 234, 234	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	2.245, 2.245, 2.245	Depositor

5 Model quality

5.1 Standard geometry

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	0.48	0/2260	1.07	5/3156 (0.2%)
2	C	0.44	0/2585	1.05	5/3606 (0.1%)
3	E	0.59	1/1778 (0.1%)	1.12	10/2479 (0.4%)
4	F	0.77	0/833	0.99	3/1158 (0.3%)
5	H	0.77	0/805	1.07	5/1119 (0.4%)
6	K	0.40	0/925	1.04	4/1287 (0.3%)
7	L	0.85	6/1159 (0.5%)	1.25	7/1614 (0.4%)
8	M	0.64	0/734	1.02	0/1023
All	All	0.59	7/11079 (0.1%)	1.08	39/15442 (0.3%)

The worst 5 of 7 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
7	L	490	LEU	C-N	7.59	1.43	1.33
7	L	408	VAL	N-CA	-6.14	1.39	1.46
7	L	332	ALA	CA-C	-6.05	1.45	1.52
7	L	333	ILE	N-CA	-6.04	1.39	1.46
7	L	492	VAL	CA-C	5.88	1.60	1.52

The worst 5 of 39 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	266	PRO	N-CA-CB	13.19	110.08	103.22
1	A	429	PRO	N-CA-CB	7.37	110.99	103.25
6	K	151	GLN	N-CA-C	-7.33	102.93	112.68
4	F	234	PRO	N-CA-CB	7.26	108.10	103.23
5	H	111	ASN	N-CA-C	6.97	118.88	111.28

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	2432	0	1010	28	0
2	C	2711	0	1169	42	0
3	E	1911	0	807	67	0
4	F	1149	0	435	17	0
5	H	1262	0	439	27	0
6	K	1012	0	439	17	0
7	L	1281	0	526	59	0
8	M	875	0	364	17	0
All	All	12633	0	5189	233	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 13.

The worst 5 of 233 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:C:484:GLU:CB	2:C:502:LYS:CB	1.75	1.62
1:A:45:LYS:CB	1:A:82:ASN:CB	1.88	1.50
2:C:628:PRO:CB	3:E:352:SER:CB	1.93	1.42
7:L:467:TYR:O	7:L:515:SER:CB	1.65	1.42
3:E:57:VAL:O	3:E:62:TYR:CB	1.67	1.42

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	
1	A	450/529 (85%)	440 (98%)	8 (2%)	2 (0%)	30	67
2	C	519/547 (95%)	492 (95%)	24 (5%)	3 (1%)	21	59
3	E	352/422 (83%)	332 (94%)	19 (5%)	1 (0%)	36	72
4	F	167/327 (51%)	162 (97%)	4 (2%)	1 (1%)	21	59
5	H	159/343 (46%)	153 (96%)	6 (4%)	0	100	100
6	K	183/208 (88%)	178 (97%)	5 (3%)	0	100	100
7	L	230/540 (43%)	218 (95%)	11 (5%)	1 (0%)	30	67
8	M	146/364 (40%)	143 (98%)	3 (2%)	0	100	100
All	All	2206/3280 (67%)	2118 (96%)	80 (4%)	8 (0%)	31	67

5 of 8 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	206	PRO
2	C	392	ASP
2	C	393	PRO
4	F	195	PRO
7	L	379	PRO

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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.

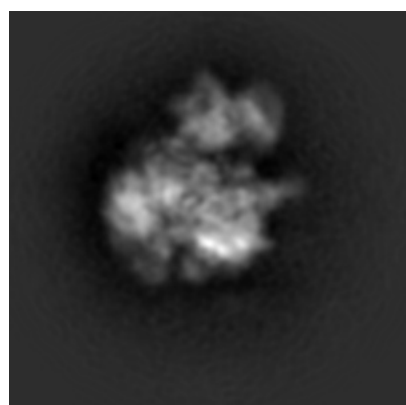
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-5658. These allow visual inspection of the internal detail of the map and identification of artifacts.

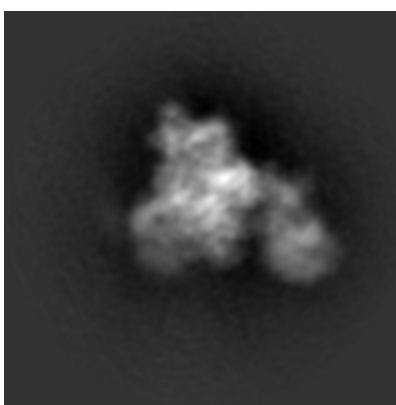
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

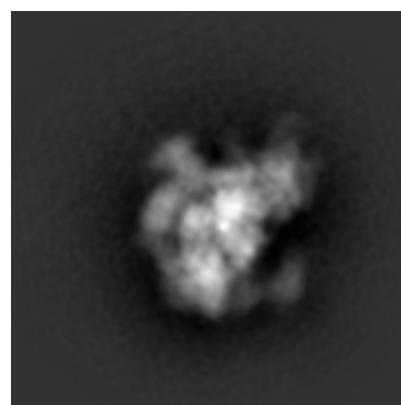
6.1.1 Primary map



X



Y

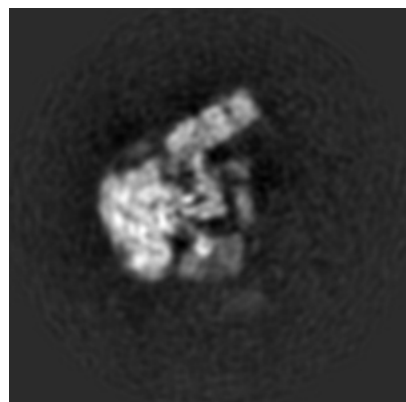


Z

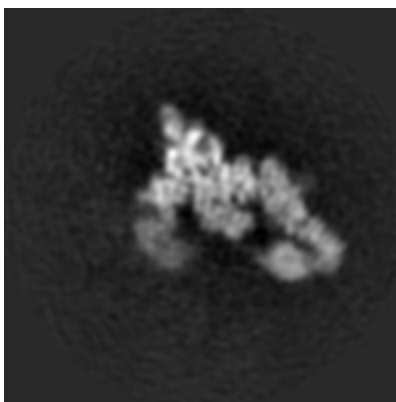
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

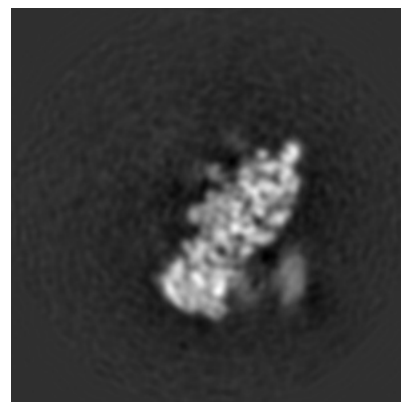
6.2.1 Primary map



X Index: 117



Y Index: 117

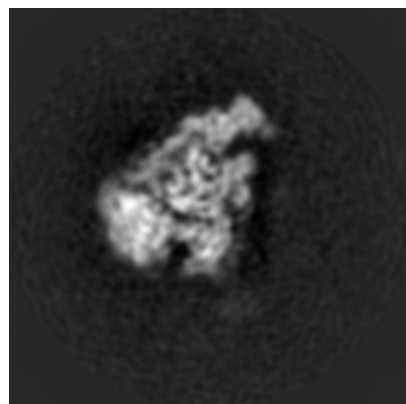


Z Index: 117

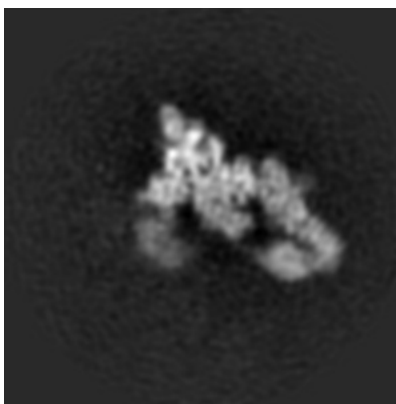
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

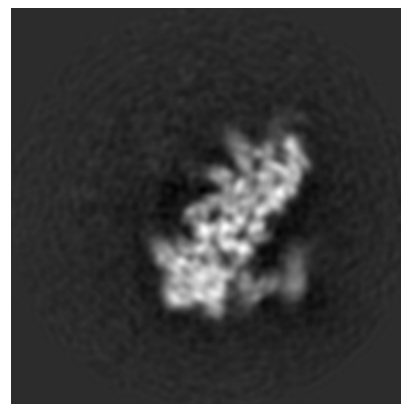
6.3.1 Primary map



X Index: 122



Y Index: 118

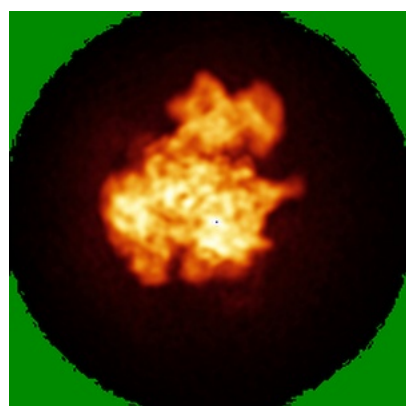


Z Index: 122

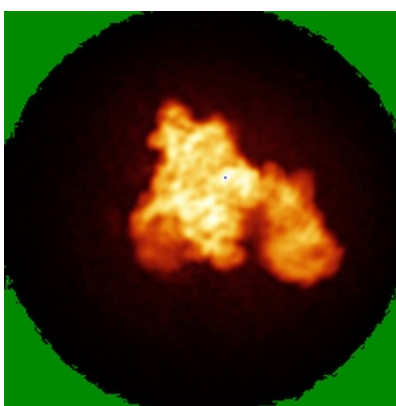
The images above show the largest variance slices of the map in three orthogonal directions.

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

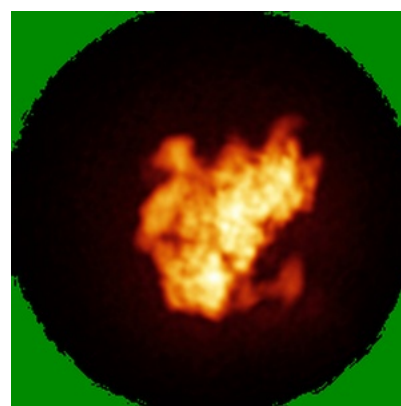
6.4.1 Primary map



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

This section was not generated.

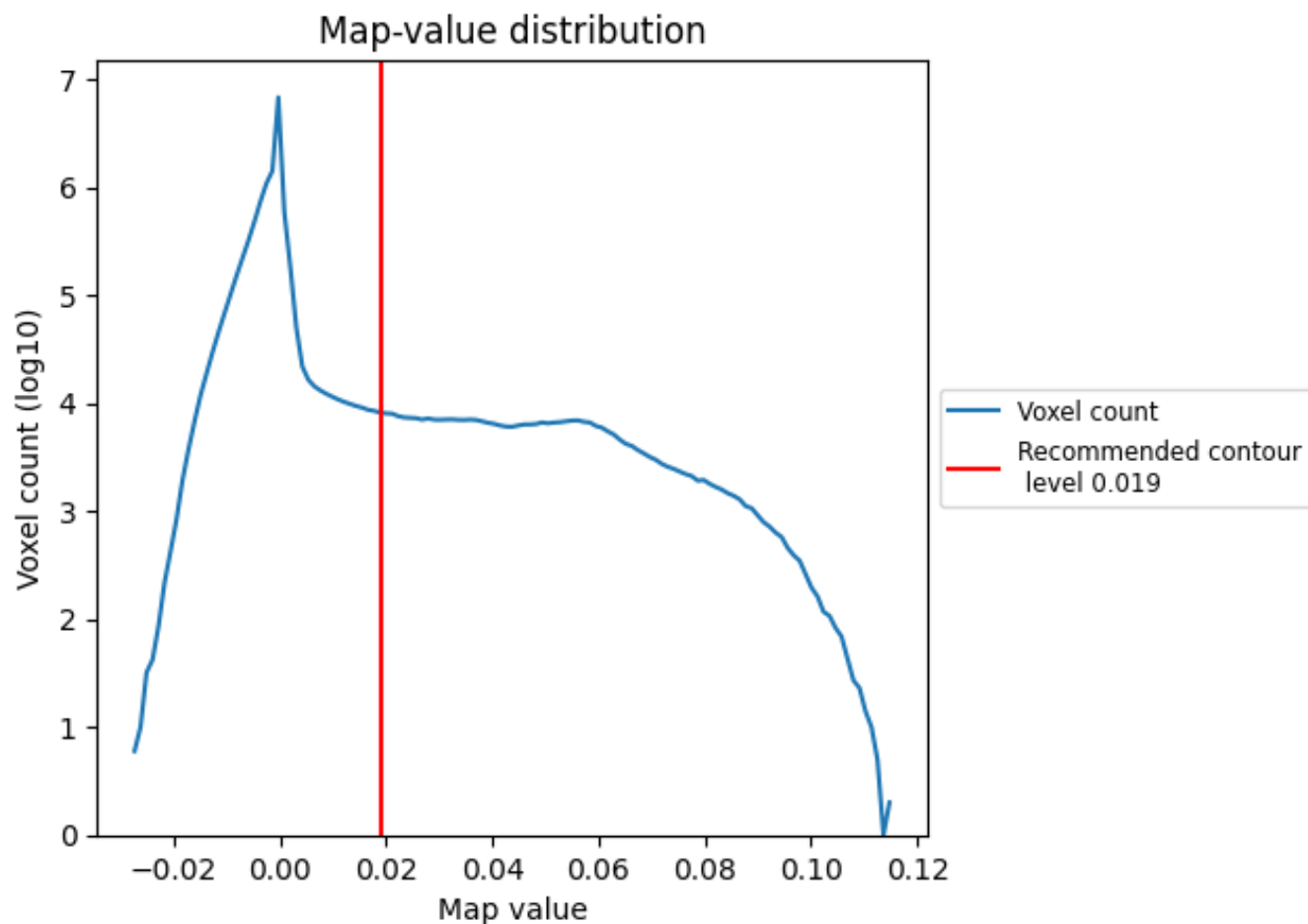
6.6 Mask visualisation

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

7 Map analysis [i](#)

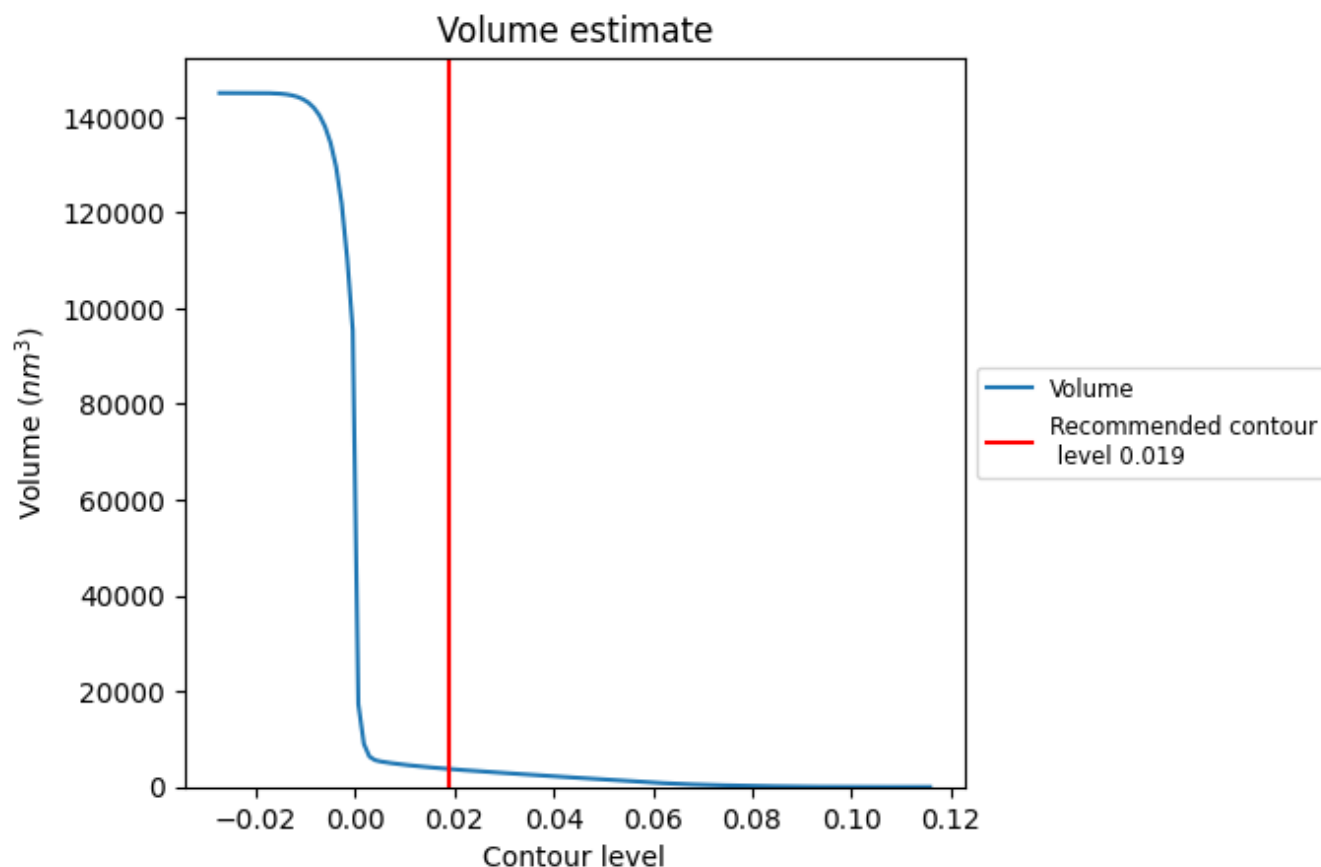
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

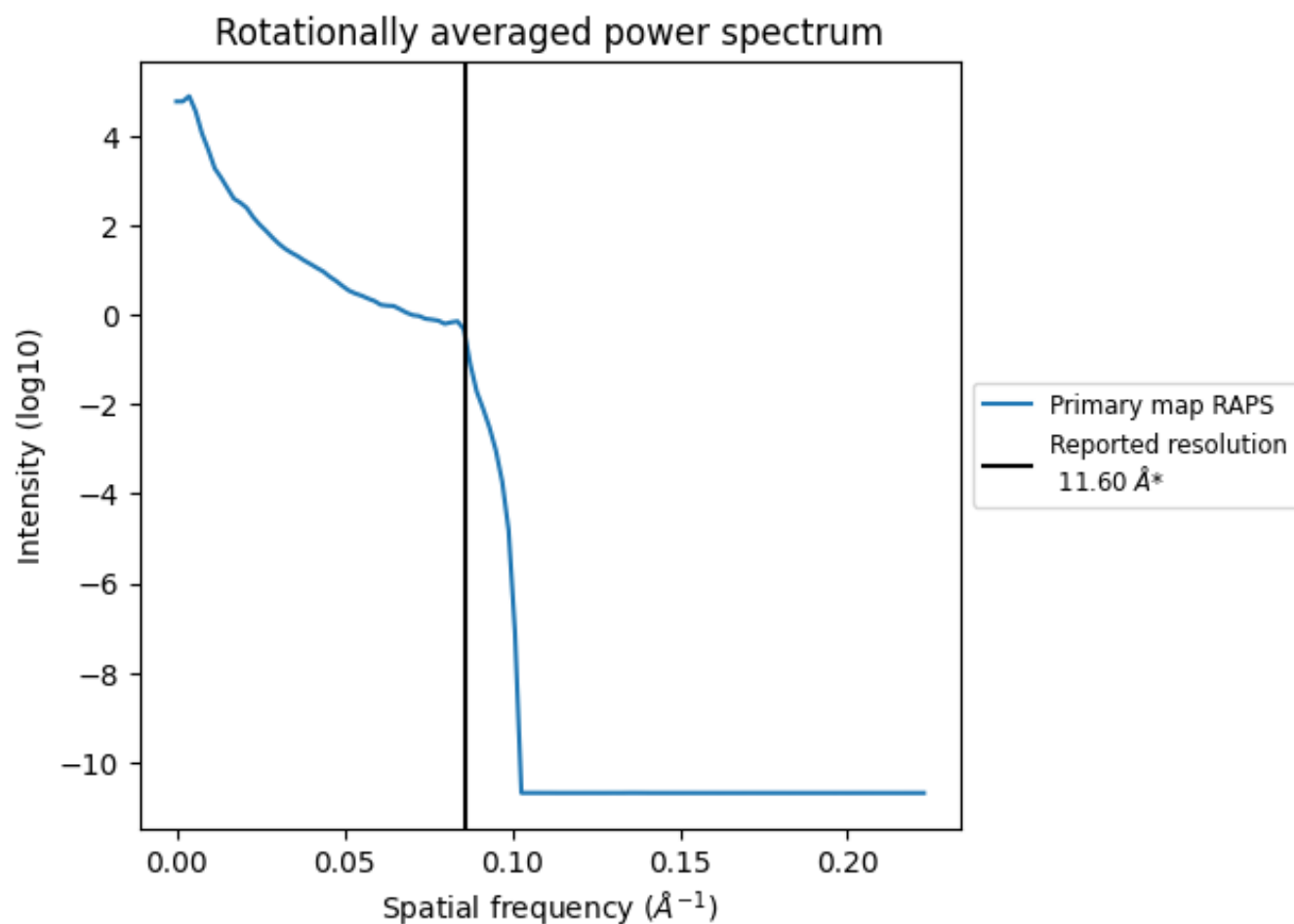
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 3751 nm^3 ; this corresponds to an approximate mass of 3388 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 $0.086 \text{ \AA}^{-1}$

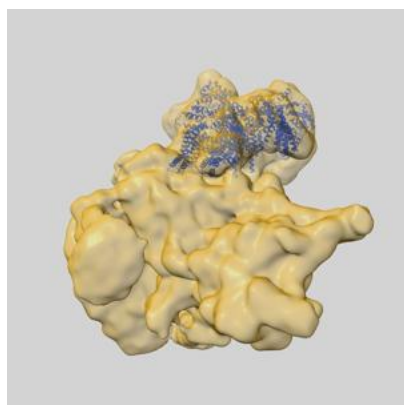
8 Fourier-Shell correlation

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

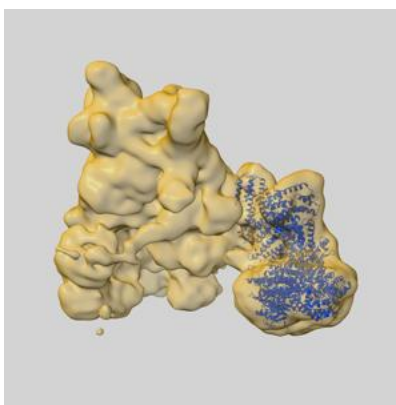
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-5658 and PDB model 3J8C. Per-residue inclusion information can be found in section [3](#) on page [6](#).

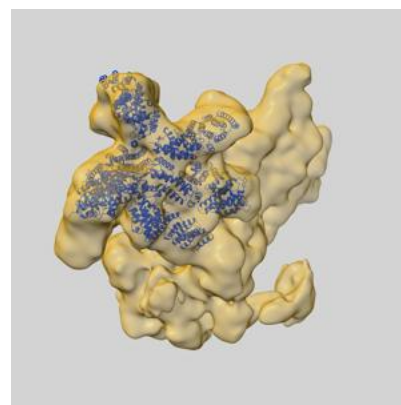
9.1 Map-model overlay [i](#)



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.019 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](#)

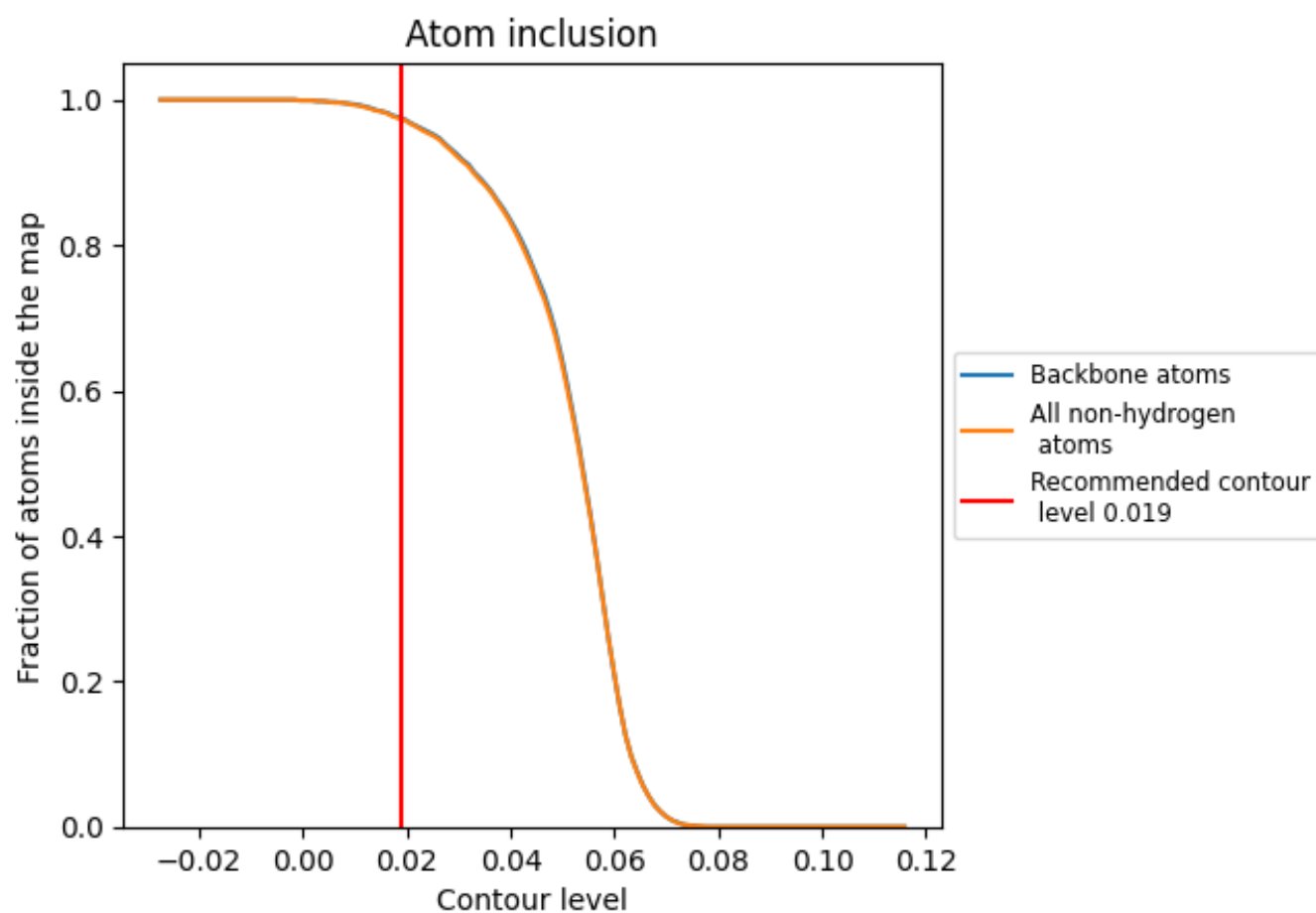


The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

This section was not generated.

9.4 Atom inclusion [i](#)



At the recommended contour level, 98% of all backbone atoms, 97% 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 (0.019) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9730	0.0660
A	0.9760	0.0700
C	0.9900	0.0870
E	0.9600	0.0720
F	0.9540	0.0270
H	0.9720	0.0500
K	0.9430	0.0650
L	0.9790	0.0460
M	0.9990	0.0850

1.0

0.0

<0.0