



## Full wwPDB EM Validation Report ⓘ

Nov 13, 2022 – 04:11 PM EST

PDB ID : 7JSI  
EMDB ID : EMD-22455  
Title : Adeno-Associated Virus 2 Rep68 HD Hexamer-ssDNA with ATPgS  
Authors : Escalante, C.R.  
Deposited on : 2020-08-14  
Resolution : 5.01 Å (reported)  
Based on initial model : 1S9H

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

We welcome your comments at [validation@mail.wwpdb.org](mailto: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>.

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

EMDB validation analysis : 0.0.1.dev43  
MolProbity : 4.02b-467  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.2

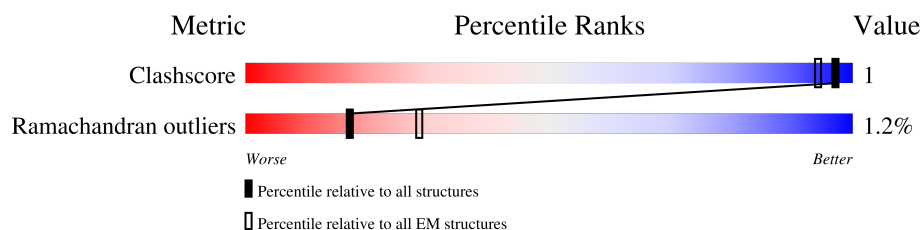
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 5.01 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023

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  $\geq 3$ , 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	536	<div> <div>11%</div> <div>51%</div> <div>49%</div> </div>
1	B	536	<div> <div>9%</div> <div>51%</div> <div>49%</div> </div>
1	C	536	<div> <div>7%</div> <div>51%</div> <div>49%</div> </div>
1	D	536	<div> <div>9%</div> <div>50%</div> <div>49%</div> </div>
1	E	536	<div> <div>12%</div> <div>51%</div> <div>49%</div> </div>
1	F	536	<div> <div>10%</div> <div>50%</div> <div>49%</div> </div>
2	H	8	<div> <div>75%</div> <div>100%</div> </div>

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 6784 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 Protein Rep68.

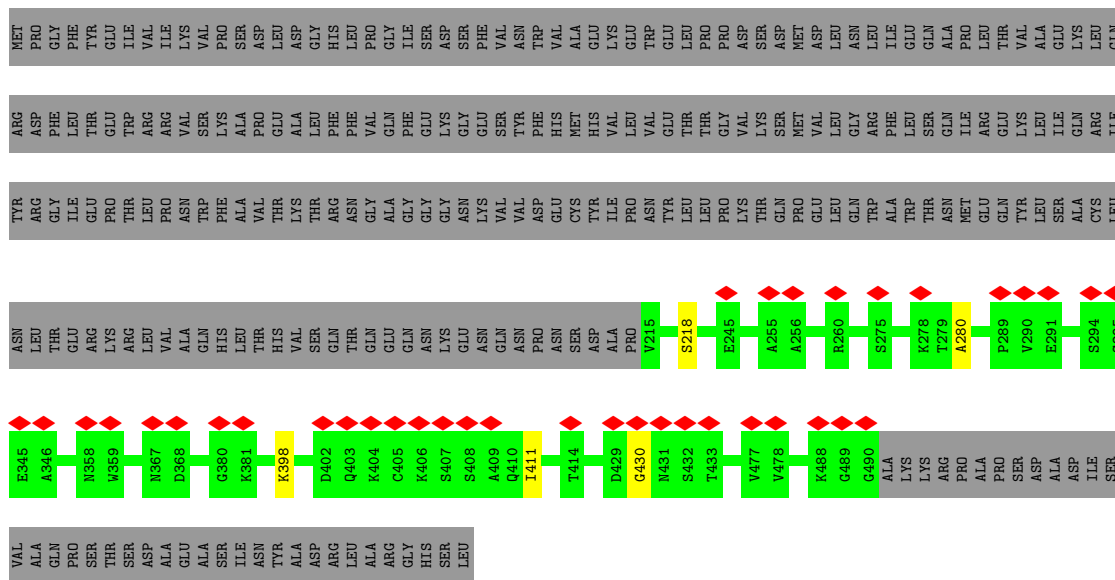
Mol	Chain	Residues	Atoms				AltConf	Trace
1	A	276	Total	C	N	O	0	0
			1104	552	276	276		
1	B	276	Total	C	N	O	0	0
			1104	552	276	276		
1	C	276	Total	C	N	O	0	0
			1104	552	276	276		
1	D	276	Total	C	N	O	0	0
			1104	552	276	276		
1	E	276	Total	C	N	O	0	0
			1104	552	276	276		
1	F	276	Total	C	N	O	0	0
			1104	552	276	276		

- Molecule 2 is a DNA chain called DNA (5'-D(P\*TP\*TP\*TP\*TP\*TP\*TP\*TP\*T)-3').

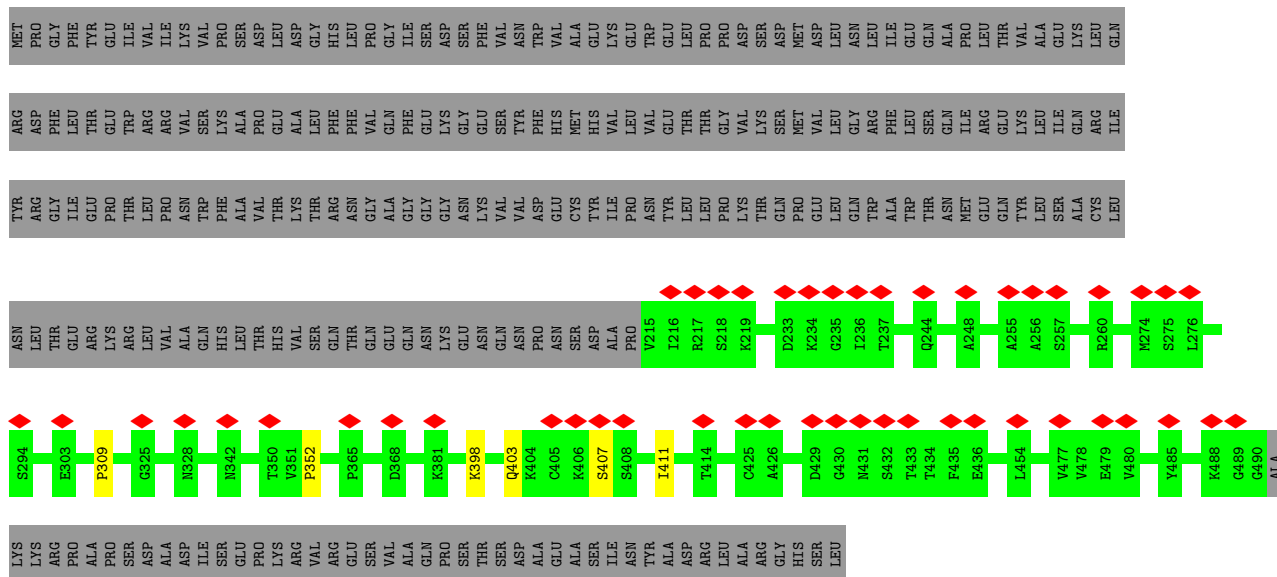
Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	8	Total	C	N	O	P	0	0
			160	80	16	56	8		



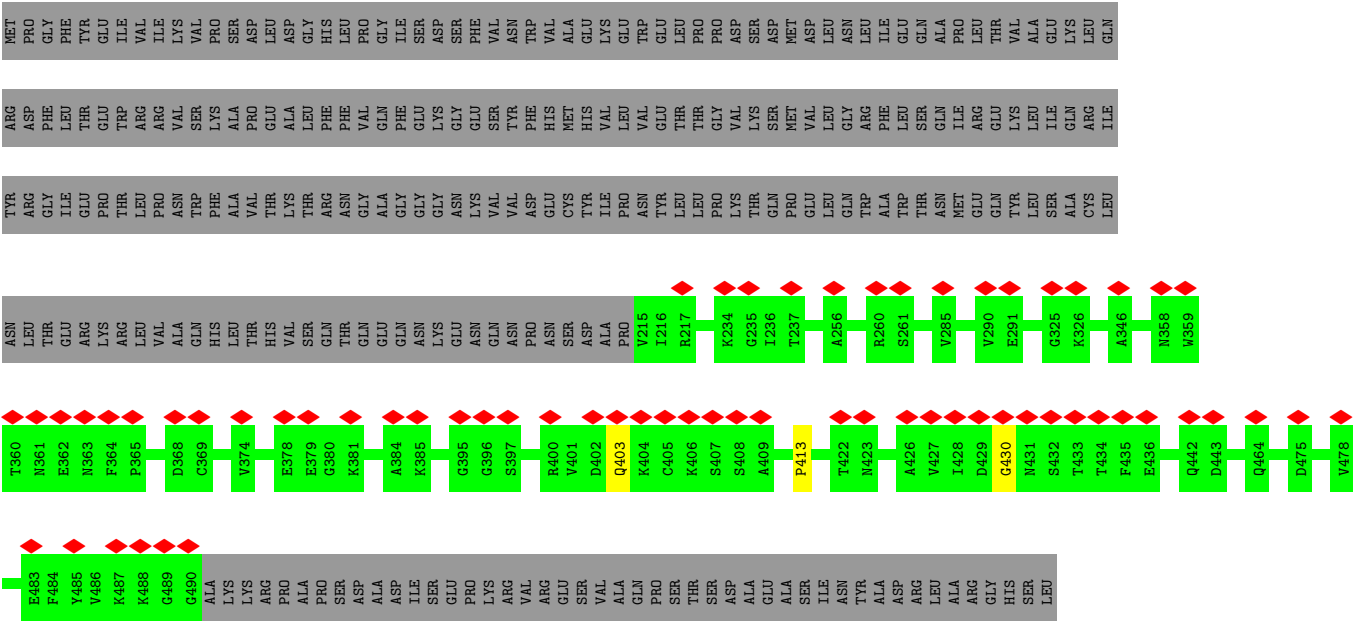
- Molecule 1: Protein Rep68



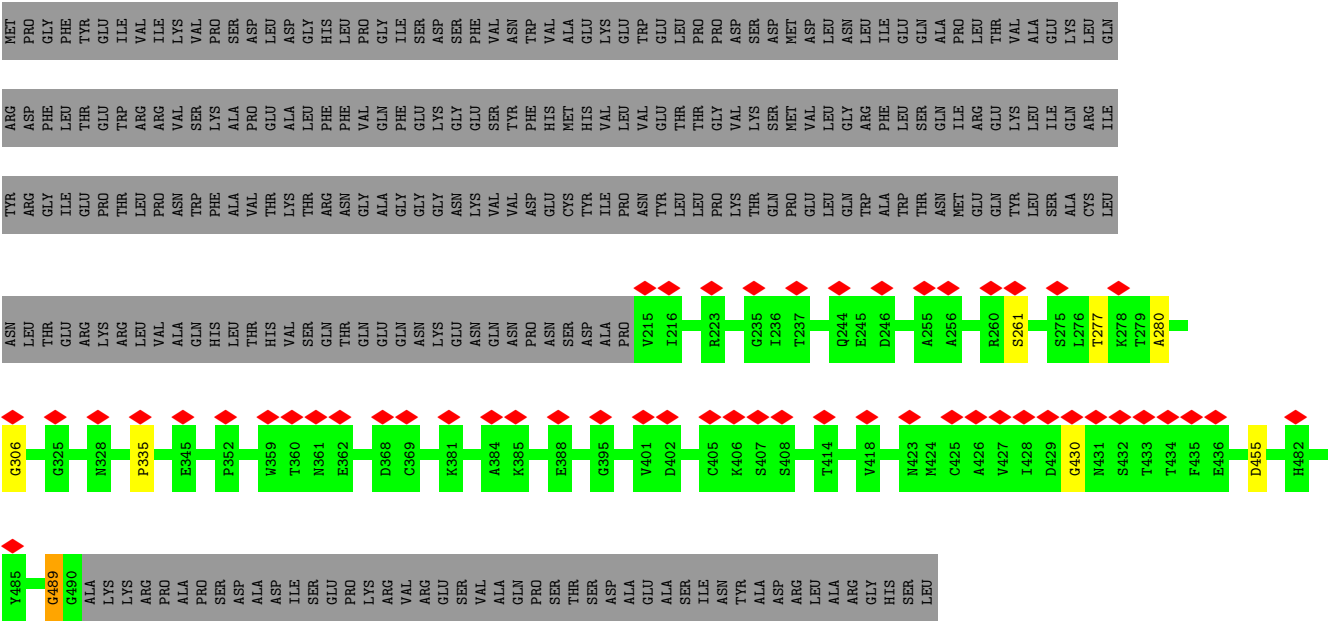
- Molecule 1: Protein Rep68



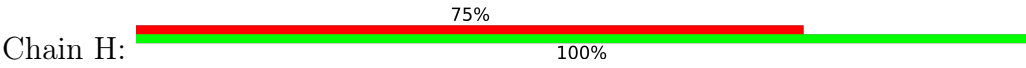
- Molecule 1: Protein Rep68



• Molecule 1: Protein Rep68



• Molecule 2: DNA (5'-D(P\*TP\*TP\*TP\*TP\*TP\*TP\*T)-3')



## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	61195	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 ( $e^-/\text{\AA}^2$ )	50	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	1.156	Depositor
Minimum map value	-0.403	Depositor
Average map value	0.007	Depositor
Map value standard deviation	0.064	Depositor
Recommended contour level	0.34	Depositor
Map size (Å)	214.59999, 214.59999, 214.59999	wwPDB
Map dimensions	200, 200, 200	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.073, 1.073, 1.073	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	0.25	0/1103	0.45	0/1377
1	B	0.26	0/1103	0.45	0/1377
1	C	0.23	0/1103	0.44	0/1377
1	D	0.30	0/1103	0.45	0/1377
1	E	0.26	0/1103	0.45	0/1377
1	F	0.23	0/1103	0.44	0/1377
2	H	0.43	0/175	1.24	0/268
All	All	0.26	0/6793	0.49	0/8530

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 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	1104	0	300	0	0
1	B	1104	0	300	1	0
1	C	1104	0	300	1	0
1	D	1104	0	300	2	0
1	E	1104	0	300	0	0
1	F	1104	0	300	1	0
2	H	160	0	97	0	0
All	All	6784	0	1897	5	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 1.

All (5) 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)
1:C:398:LYS:CA	1:C:411:ILE:O	2.53	0.57
1:B:398:LYS:CA	1:B:411:ILE:O	2.63	0.47
1:D:398:LYS:CA	1:D:411:ILE:O	2.65	0.45
1:D:403:GLN:N	1:D:407:SER:O	2.48	0.42
1:F:306:GLY:O	1:F:489:GLY:N	2.54	0.41

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	274/536 (51%)	243 (89%)	29 (11%)	2 (1%)	22	62
1	B	274/536 (51%)	235 (86%)	36 (13%)	3 (1%)	14	52
1	C	274/536 (51%)	240 (88%)	31 (11%)	3 (1%)	14	52
1	D	274/536 (51%)	239 (87%)	33 (12%)	2 (1%)	22	62
1	E	274/536 (51%)	244 (89%)	27 (10%)	3 (1%)	14	52
1	F	274/536 (51%)	243 (89%)	24 (9%)	7 (3%)	5	34
All	All	1644/3216 (51%)	1444 (88%)	180 (11%)	20 (1%)	17	50

All (20) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	457	ASP
1	B	456	HIS
1	C	218	SER

*Continued on next page...*

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Mol	Chain	Res	Type
1	E	403	GLN
1	F	277	THR
1	A	413	PRO
1	A	481	GLU
1	C	430	GLY
1	E	430	GLY
1	F	455	ASP
1	B	430	GLY
1	C	280	ALA
1	F	335	PRO
1	F	430	GLY
1	D	309	PRO
1	F	261	SER
1	F	280	ALA
1	F	489	GLY
1	D	352	PRO
1	E	413	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 monosaccharides 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.

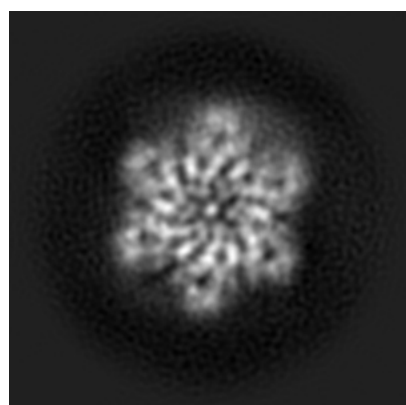
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-22455. These allow visual inspection of the internal detail of the map and identification of artifacts.

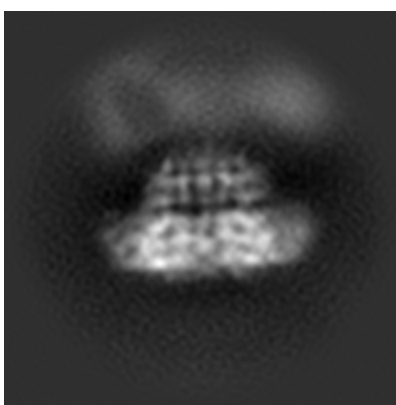
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

### 6.1 Orthogonal projections [i](#)

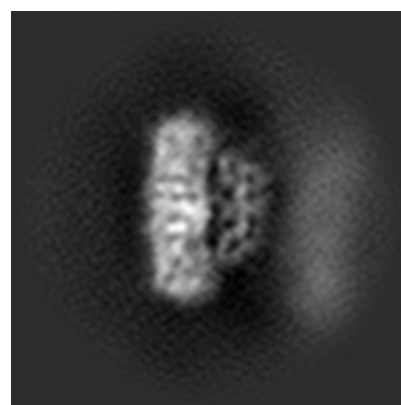
#### 6.1.1 Primary map



X



Y

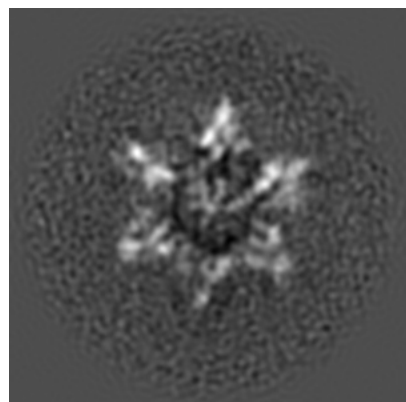


Z

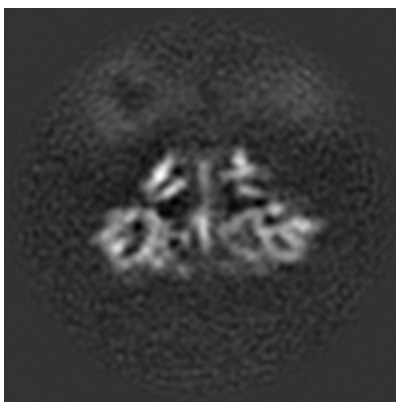
The images above show the map projected in three orthogonal directions.

### 6.2 Central slices [i](#)

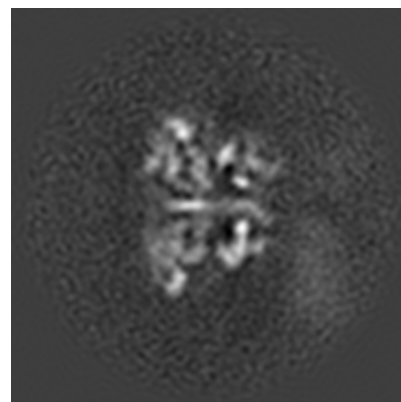
#### 6.2.1 Primary map



X Index: 100



Y Index: 100

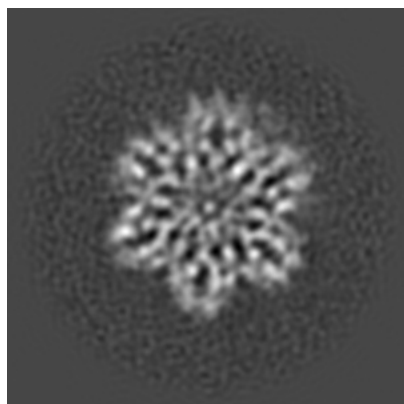


Z Index: 100

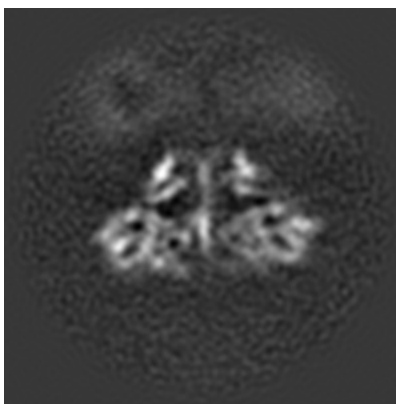
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

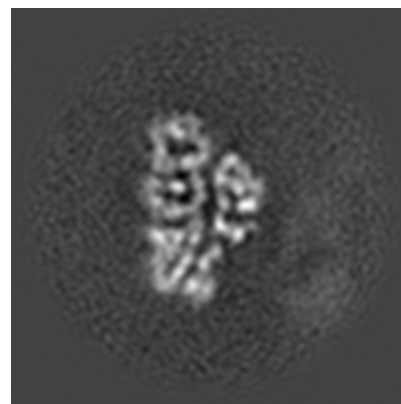
### 6.3.1 Primary map



X Index: 85



Y Index: 101

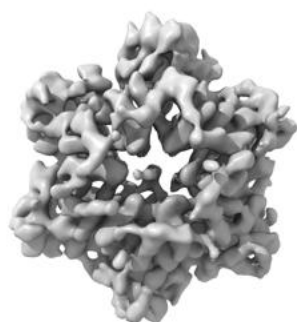


Z Index: 79

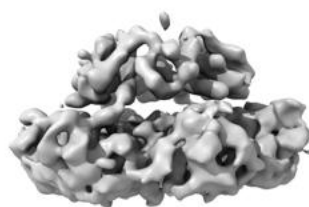
The images above show the largest variance slices of the map in three orthogonal directions.

## 6.4 Orthogonal surface views [i](#)

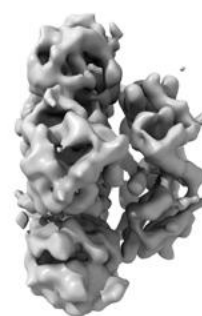
### 6.4.1 Primary map



X



Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.34. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

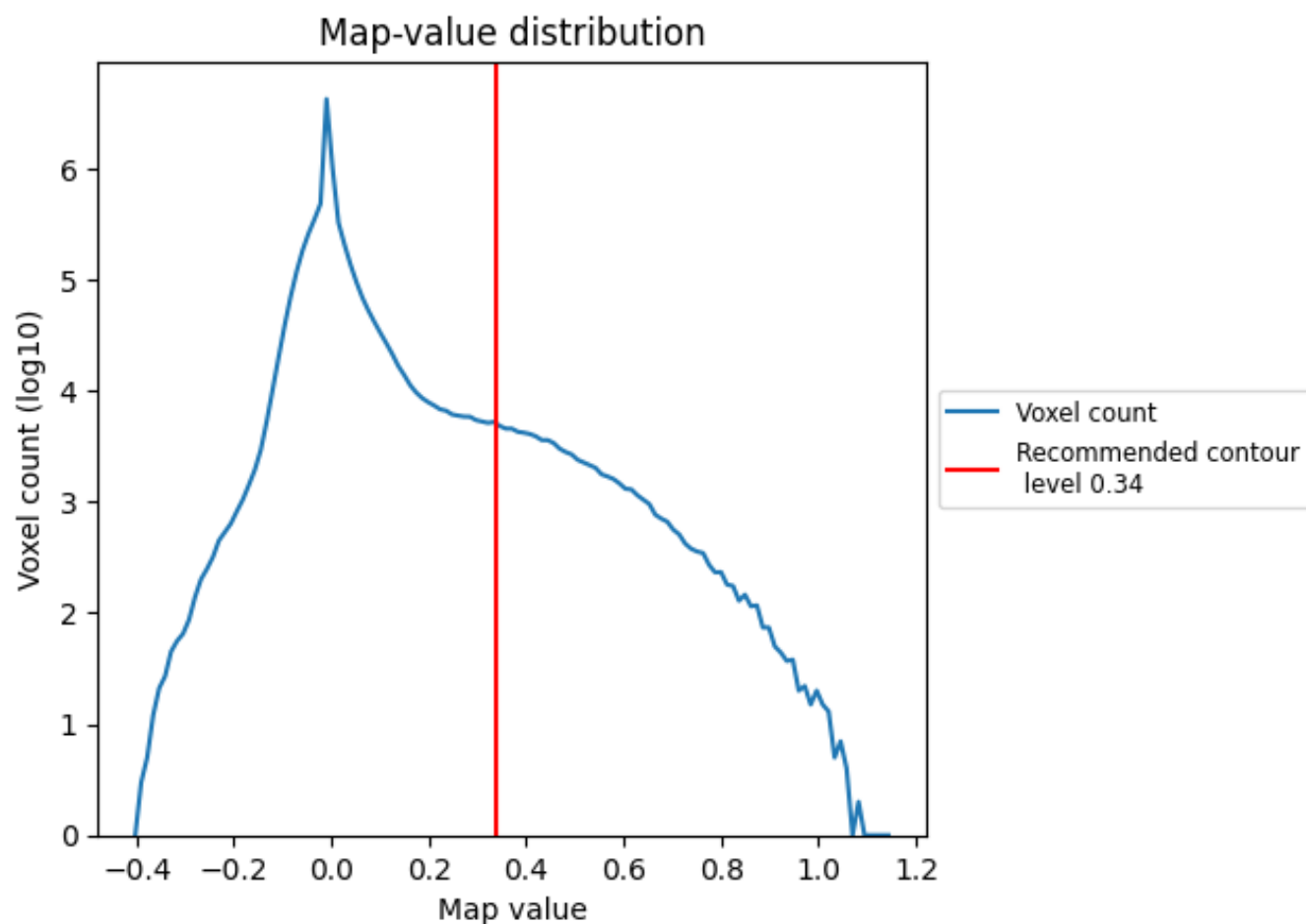
## 6.5 Mask visualisation

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

## 7 Map analysis [i](#)

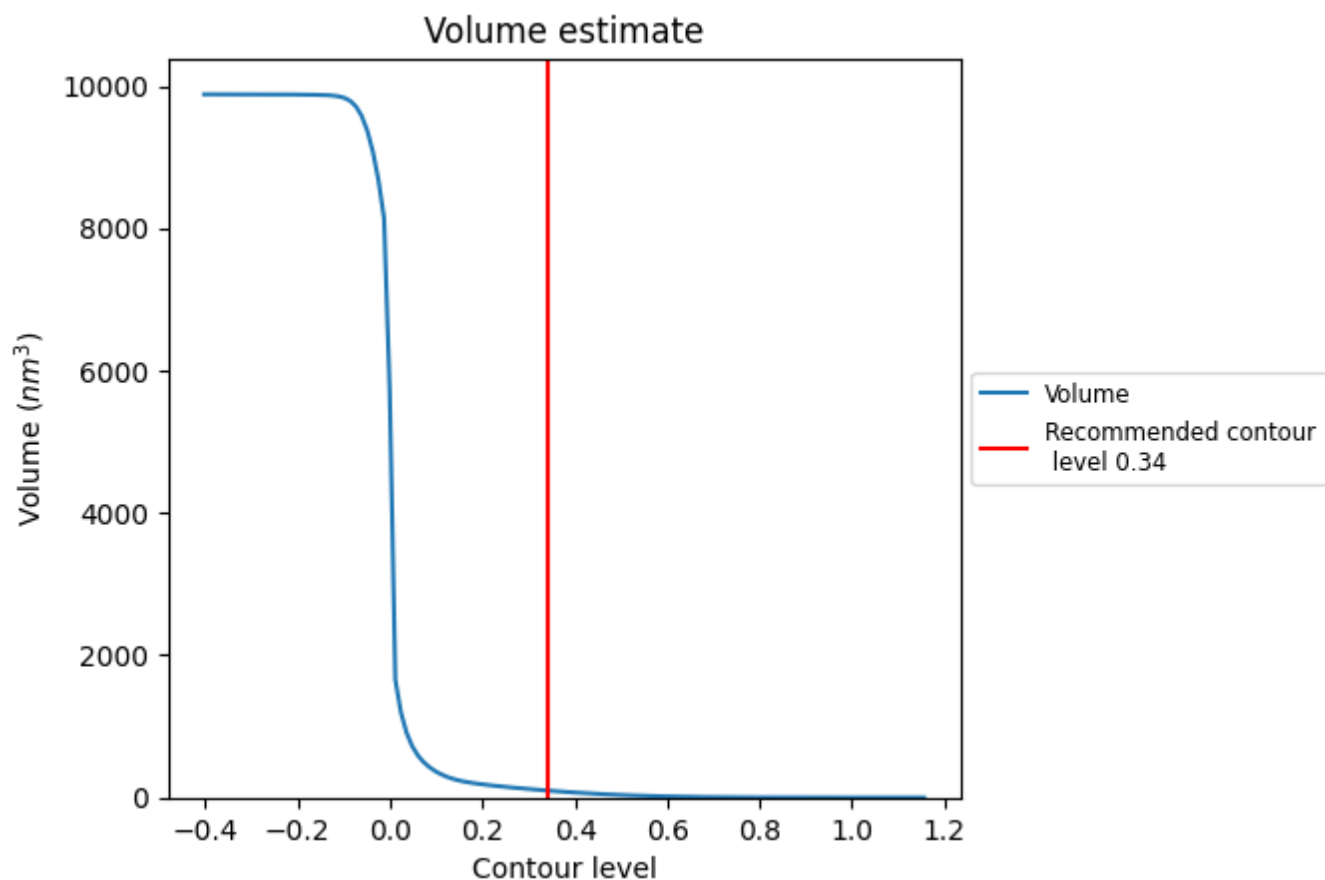
This section contains the results of statistical analysis of the map.

### 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](#)

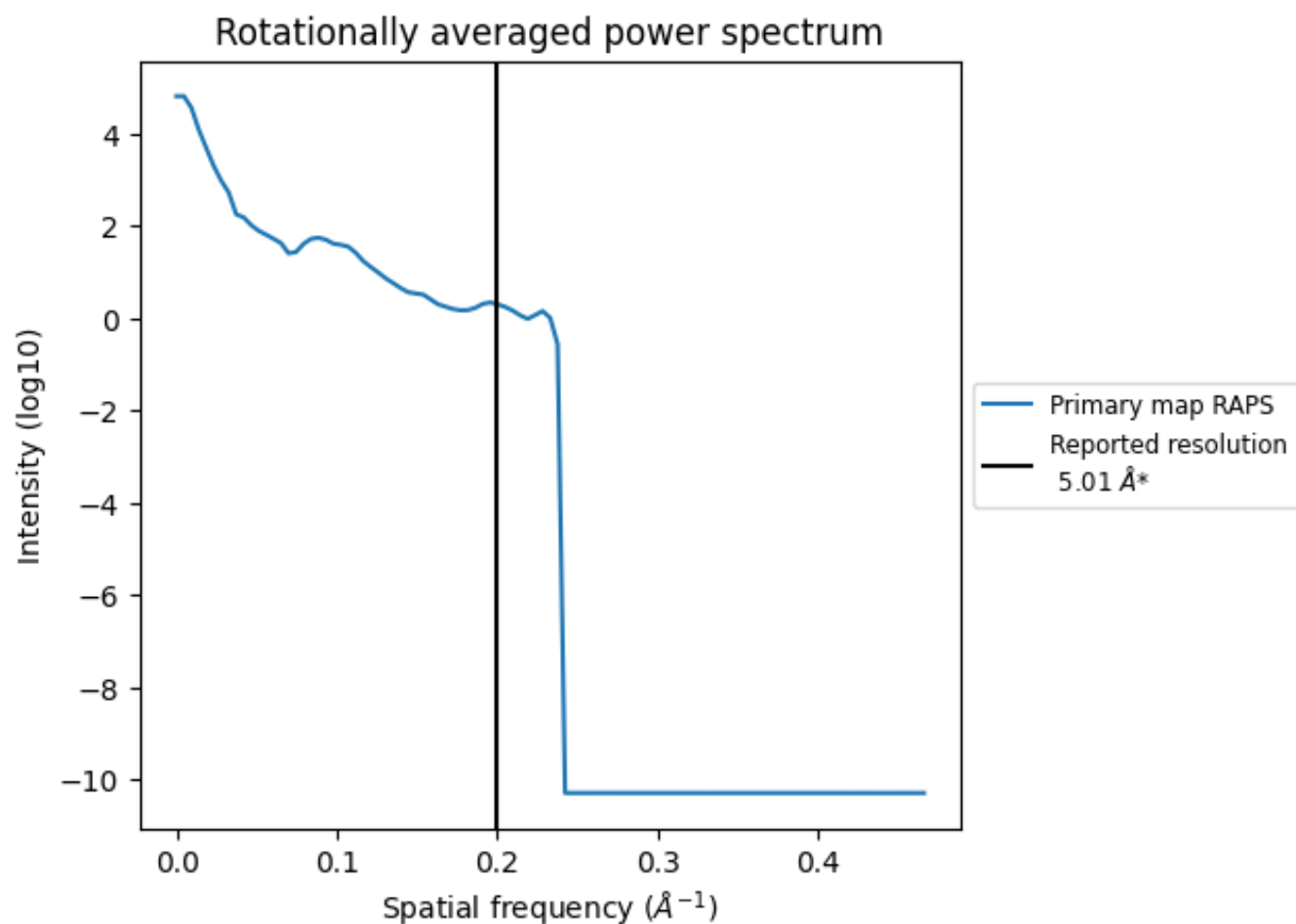


The volume at the recommended contour level is 99 nm<sup>3</sup>; this corresponds to an approximate mass of 89 kDa.

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 0.200 Å<sup>-1</sup>

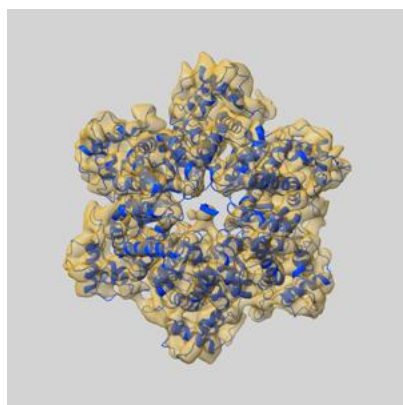
## 8 Fourier-Shell correlation ⓘ

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

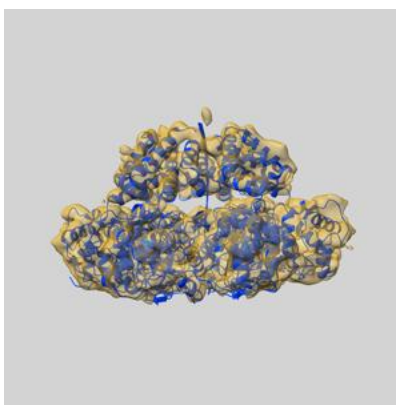
## 9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-22455 and PDB model 7JSI. Per-residue inclusion information can be found in [section 3](#) on [page 4](#).

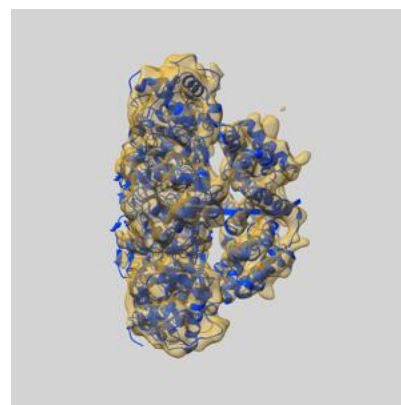
### 9.1 Map-model overlay [i](#)



X



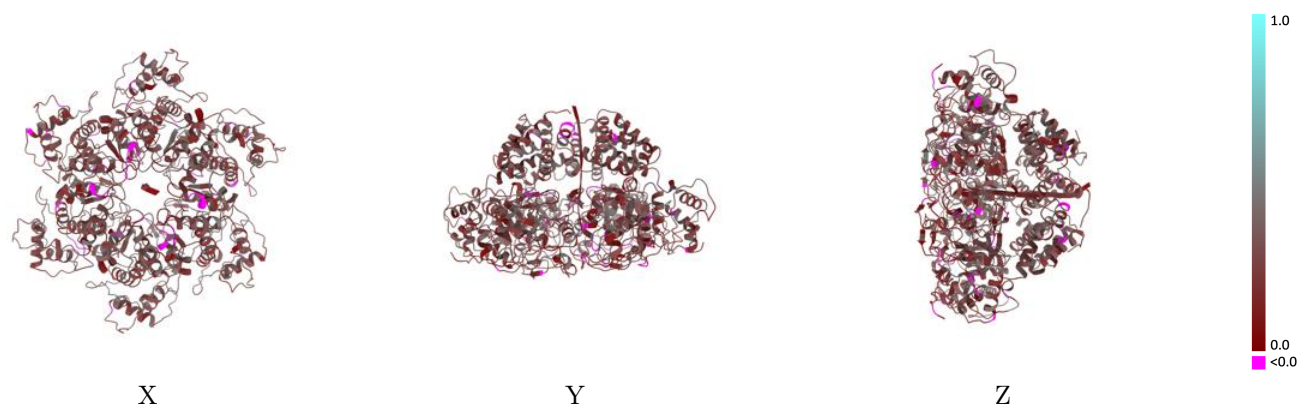
Y



Z

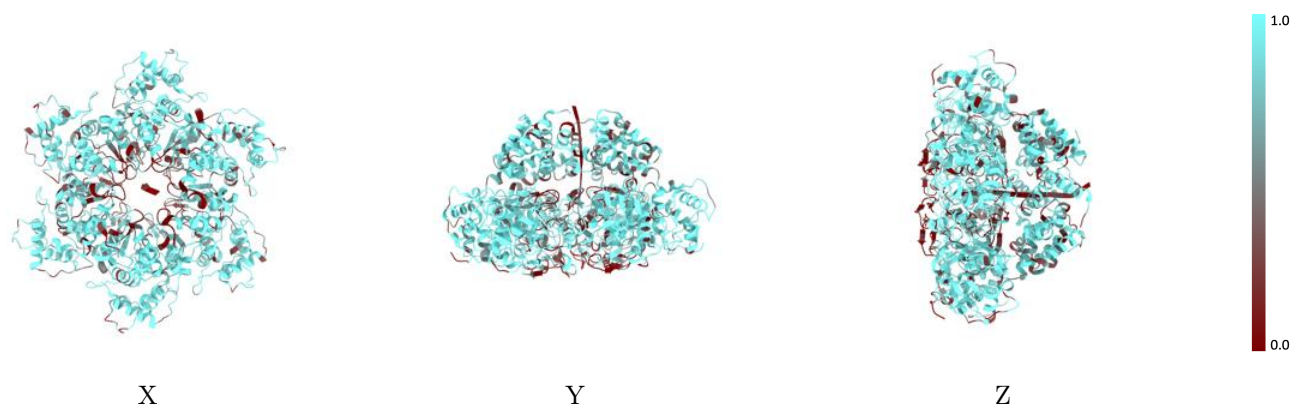
The images above show the 3D surface view of the map at the recommended contour level 0.34 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



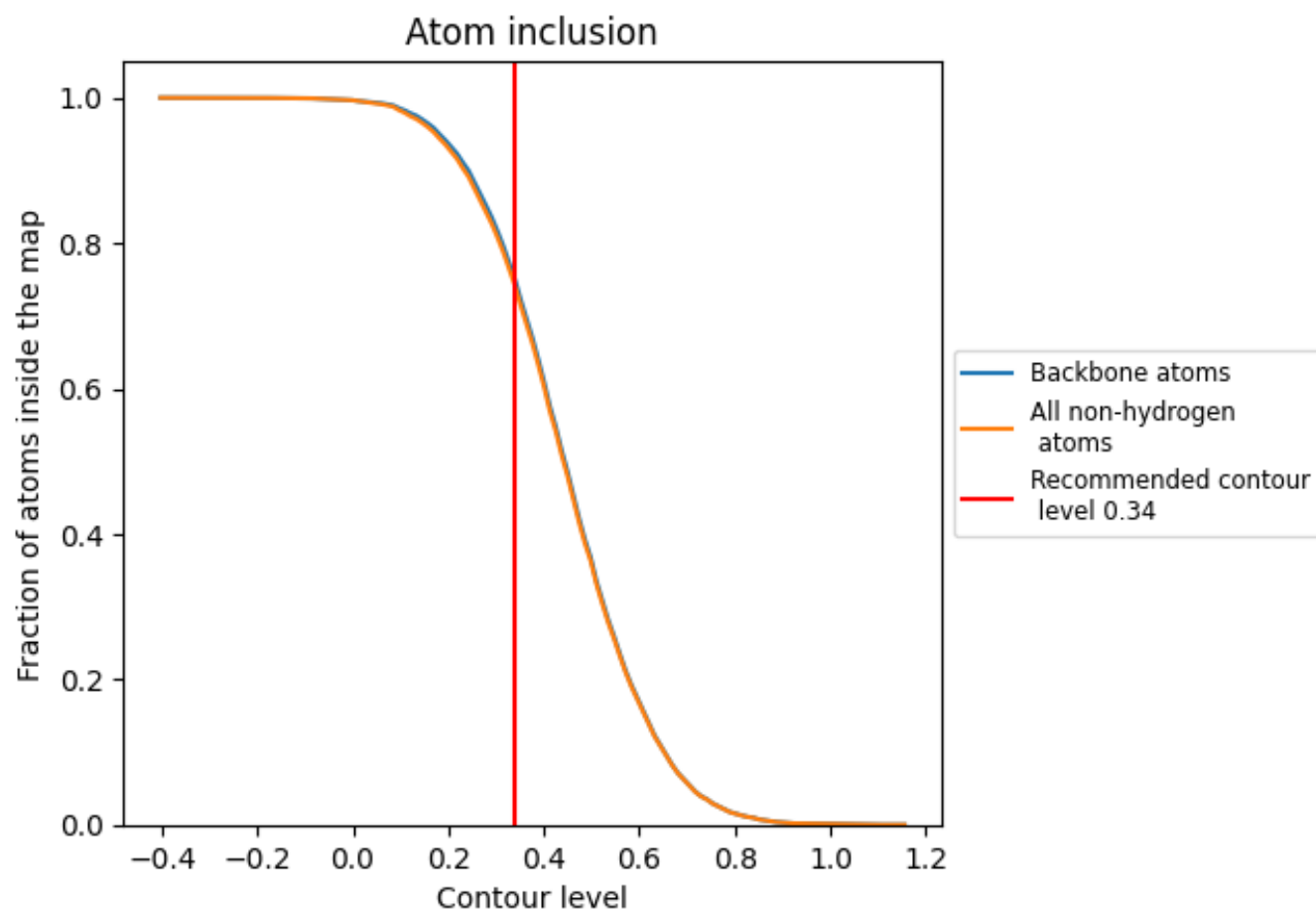
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.34).

## 9.4 Atom inclusion [i](#)



At the recommended contour level, 75% of all backbone atoms, 74% 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 (0.34) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div>0.7410</div>	<div><div></div>0.2600</div>
A	<div><div></div>0.7138</div>	<div><div></div>0.2700</div>
B	<div><div></div>0.7690</div>	<div><div></div>0.2650</div>
C	<div><div></div>0.8043</div>	<div><div></div>0.2750</div>
D	<div><div></div>0.7654</div>	<div><div></div>0.2610</div>
E	<div><div></div>0.7092</div>	<div><div></div>0.2510</div>
F	<div><div></div>0.7554</div>	<div><div></div>0.2510</div>
H	<div><div></div>0.2500</div>	<div><div></div>0.1850</div>

1.0

0.0

<0.0