wwPDB EM Validation Summary Report

Apr 15, 2026 – 02:24 PM JST

PDB ID	:	21BH / pdb_000021bh
EMDB ID	:	EMD-67552
Title	:	Cryo-EM structure of type VII CRISPR-Cas complex at the target engagement state
Authors	:	Zhang, H.; Zhang, S.
Deposited on	:	2025-12-06
Resolution	:	3.10 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:

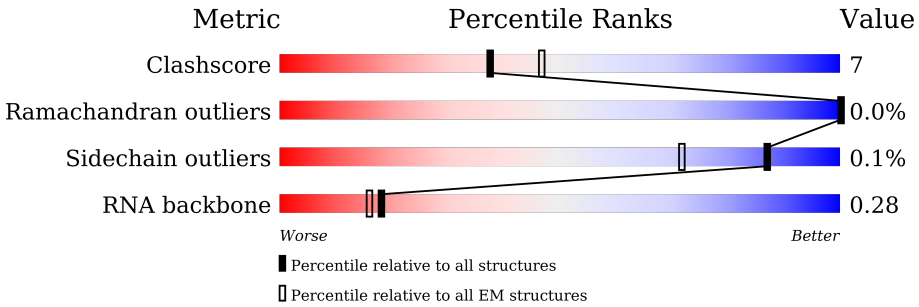
EMDB validation analysis	:	0.0.1.dev132
MolProbity	:	4-5-2 with Phenix2.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh &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

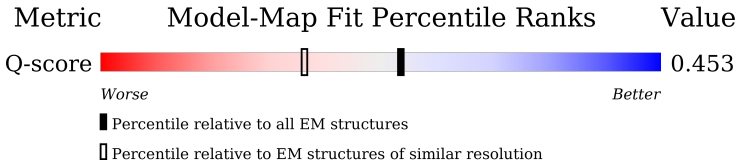
1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.10 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	14724 (2.60 - 3.60)

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 <=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	B	200	82% 17%
1	C	200	88% 12%
1	D	200	89% 10%

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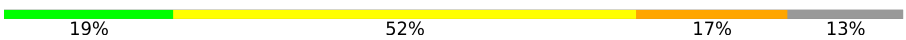
Mol	Chain	Length	Quality of chain
1	E	200	
1	F	200	
1	J	200	
1	O	200	
2	G	240	
3	H	609	
3	I	609	
3	K	609	
3	L	609	
4	M	54	
5	N	60	

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 25416 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	B	199	Total	C	N	O	S	0	0
			1525	950	271	296	8		
1	C	198	Total	C	N	O	S	0	0
			1530	954	275	293	8		
1	D	199	Total	C	N	O	S	0	0
			1529	953	272	296	8		
1	E	199	Total	C	N	O	S	0	0
			1527	953	270	296	8		
1	F	196	Total	C	N	O	S	0	0
			1489	926	262	293	8		
1	J	198	Total	C	N	O	S	0	0
			1534	956	275	295	8		
1	O	186	Total	C	N	O	S	0	0
			1416	889	245	274	8		

- Molecule 2 is a protein called a protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	G	239	Total	C	N	O	S	0	0
			1791	1149	291	347	4		

- Molecule 3 is a protein called a protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	H	577	Total	C	N	O	S	0	0
			4594	2922	763	894	15		
3	I	155	Total	C	N	O	S	0	0
			1296	833	215	241	7		
3	K	531	Total	C	N	O	S	0	0
			4156	2636	689	816	15		
3	L	157	Total	C	N	O	S	0	0
			1246	794	206	241	5		

- Molecule 4 is a RNA chain called RNA (54-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
4	M	47	Total 986	C 443	N 163	O 333	P 47	0	0

- Molecule 5 is a RNA chain called RNA (60-MER).

Mol	Chain	Residues	Atoms					AltConf	Trace
5	N	37	Total 788	C 353	N 144	O 254	P 37	0	0

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

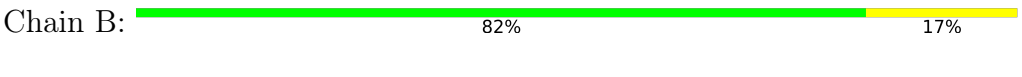
Mol	Chain	Residues	Atoms		AltConf
6	B	1	Total 1	Zn 1	0
6	C	1	Total 1	Zn 1	0
6	D	1	Total 1	Zn 1	0
6	E	1	Total 1	Zn 1	0
6	F	1	Total 1	Zn 1	0
6	G	1	Total 1	Zn 1	0
6	H	1	Total 1	Zn 1	0
6	J	1	Total 1	Zn 1	0
6	O	1	Total 1	Zn 1	0

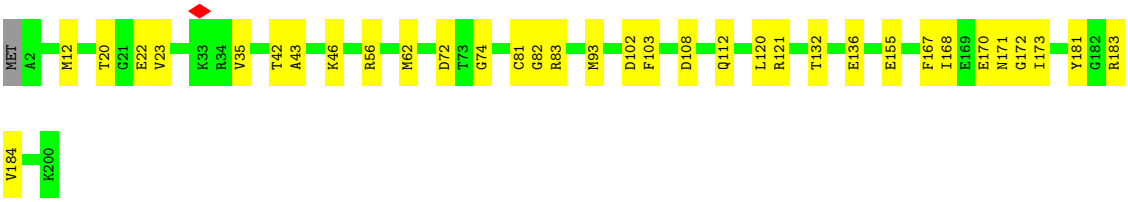
3 Residue-property plots (i)

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: a protein

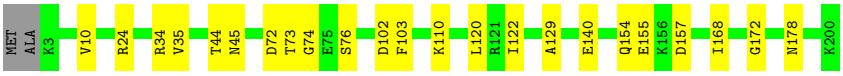
Chain B:





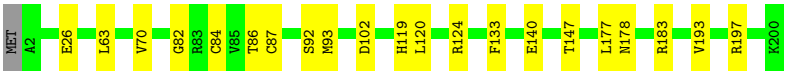
• Molecule 1: a protein

Chain C:



• Molecule 1: a protein

Chain D:

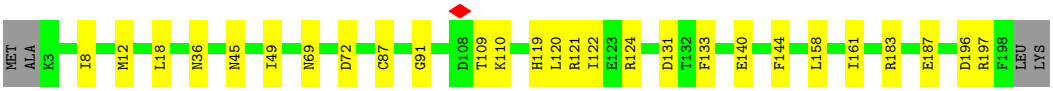


• Molecule 1: a protein

Chain E:

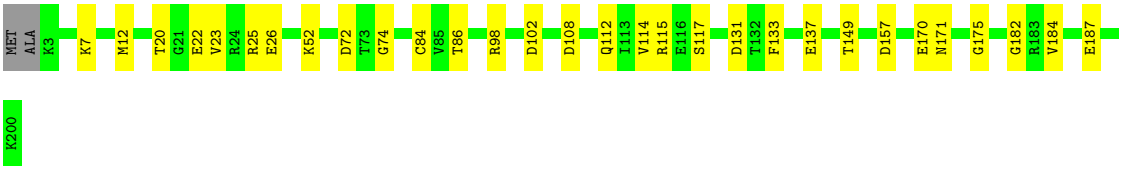
• Molecule 1: a protein

Chain F:



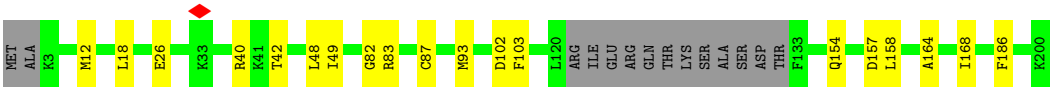
• Molecule 1: a protein

Chain J:



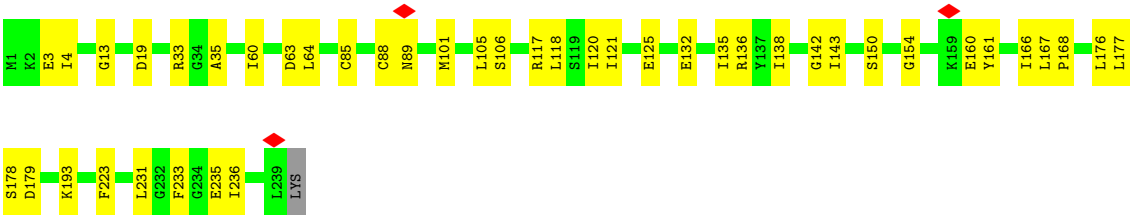
• Molecule 1: a protein

Chain O:



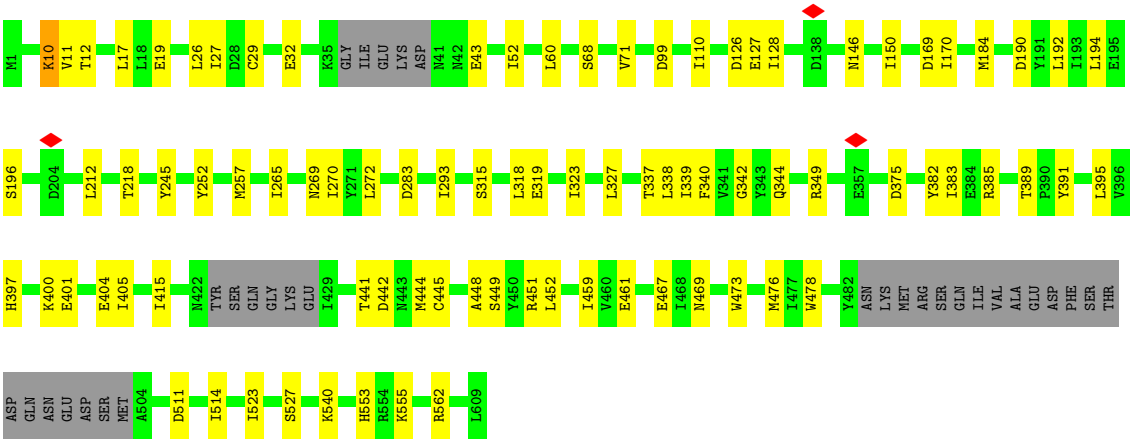
• Molecule 2: a protein

Chain G:



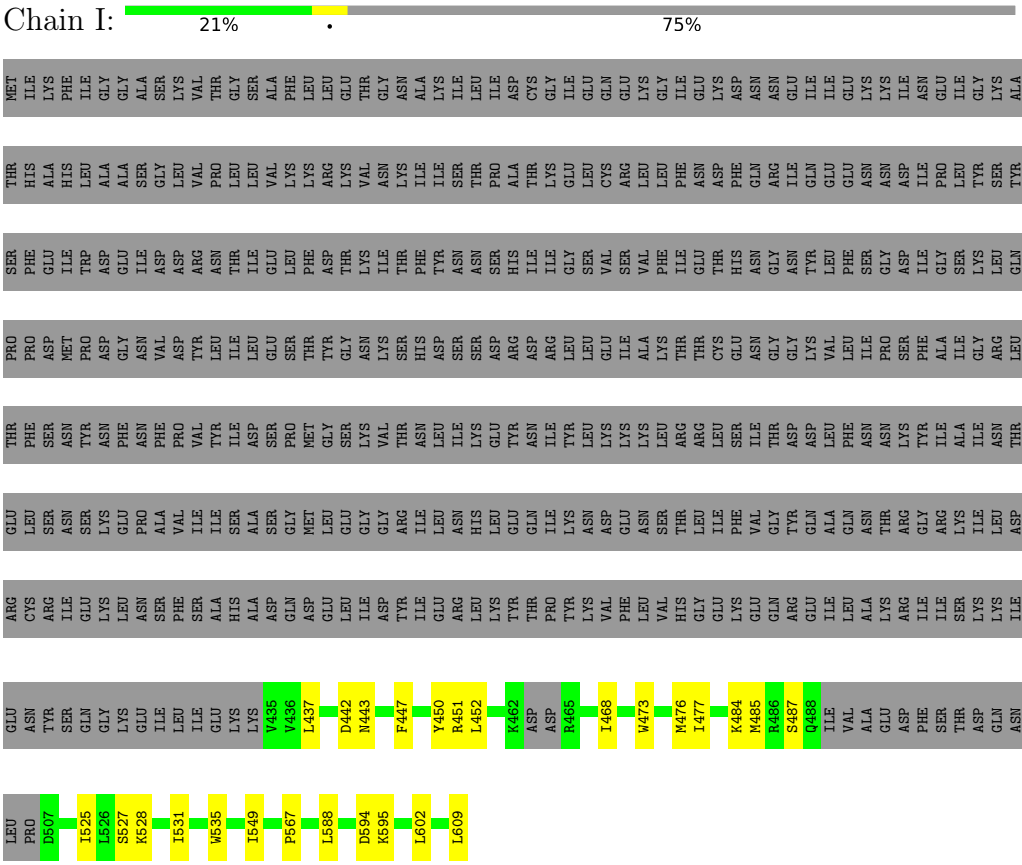
• Molecule 3: a protein

Chain H:



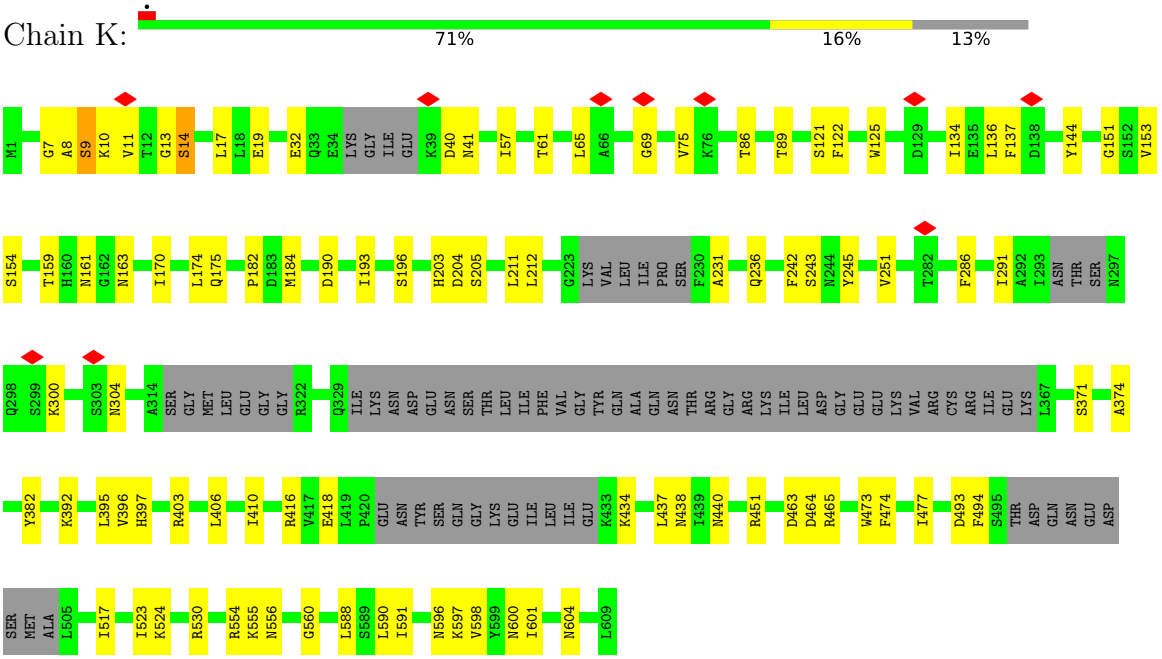
• Molecule 3: a protein

Chain I:



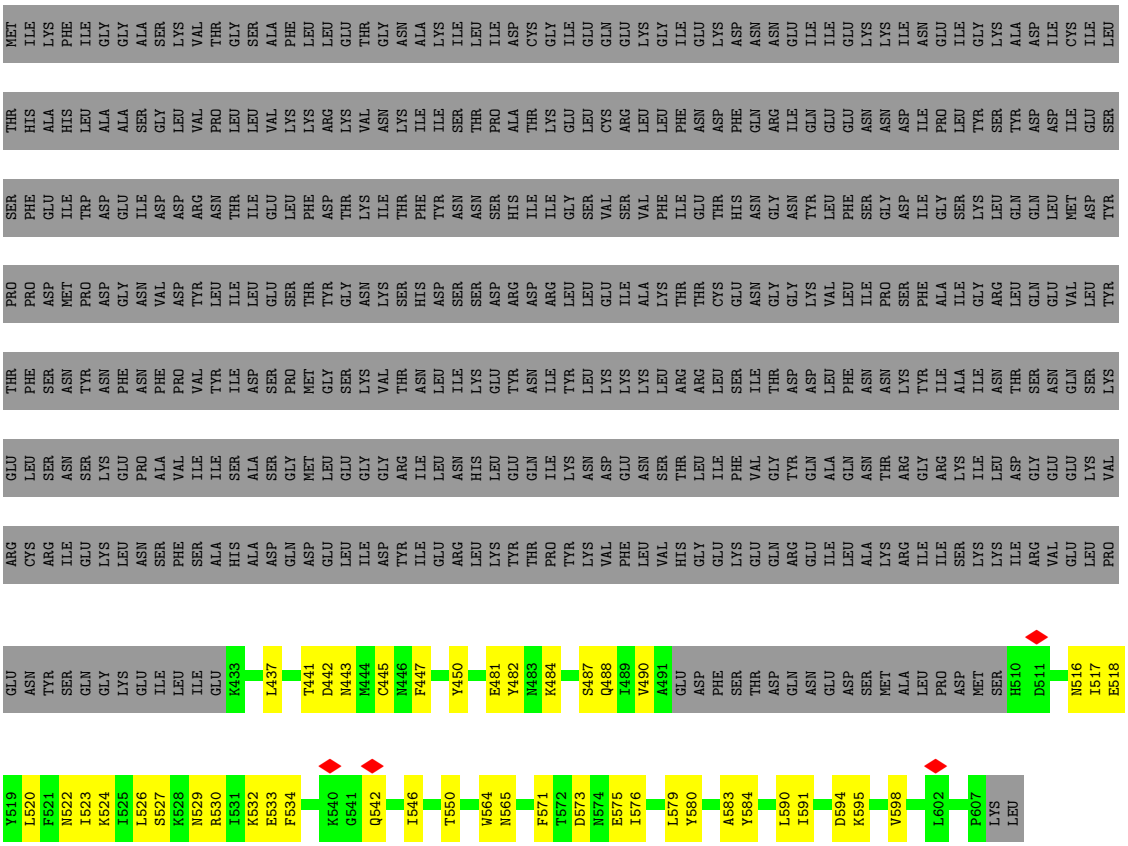
• Molecule 3: a protein

Chain K:



• Molecule 3: a protein

Chain L:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	65006	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 ($e^-/Å^2$)	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	0.712	Depositor
Minimum map value	-0.262	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.016	Depositor
Recommended contour level	0.11	Depositor
Map size (Å)	395.19998, 395.19998, 395.19998	wwPDB
Map dimensions	416, 416, 416	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.9499999, 0.9499999, 0.9499999	Depositor

5 Model quality

5.1 Standard geometry

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	B	0.15	0/1544	0.33	0/2076
1	C	0.15	0/1549	0.34	0/2079
1	D	0.16	0/1548	0.35	0/2080
1	E	0.15	0/1546	0.34	0/2077
1	F	0.16	0/1507	0.36	0/2028
1	J	0.16	0/1553	0.32	0/2084
1	O	0.12	0/1433	0.30	0/1924
2	G	0.16	0/1825	0.37	1/2469 (0.0%)
3	H	0.16	0/4678	0.36	1/6333 (0.0%)
3	I	0.14	0/1325	0.32	0/1789
3	K	0.18	0/4225	0.34	0/5722
3	L	0.17	0/1274	0.40	0/1734
4	M	0.23	0/1099	0.44	0/1706
5	N	0.23	0/881	0.36	0/1370
All	All	0.17	0/25987	0.35	2/35471 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	H	10	LYS	CB-CA-C	-7.08	99.87	110.06
2	G	168	PRO	CA-N-CD	-6.43	103.00	112.00

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	B	1525	0	1531	23	0
1	C	1530	0	1555	20	0
1	D	1529	0	1542	16	0
1	E	1527	0	1542	24	0
1	F	1489	0	1484	25	0
1	J	1534	0	1559	20	0
1	O	1416	0	1432	14	0
2	G	1791	0	1675	29	0
3	H	4594	0	4448	72	0
3	I	1296	0	1244	16	0
3	K	4156	0	3958	73	0
3	L	1246	0	1115	37	0
4	M	986	0	500	34	0
5	N	788	0	401	20	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	0	0
6	H	1	0	0	0	0
6	J	1	0	0	0	0
6	O	1	0	0	0	0
All	All	25416	0	23986	370	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 7.

The worst 5 of 370 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)
3:K:7:GLY:HA2	3:K:14:SER:HA	1.47	0.97
1:C:76:SER:OG	1:J:131:ASP:OD1	1.92	0.86
3:K:8:ALA:HB2	3:K:396:VAL:HG12	1.56	0.85
1:C:102:ASP:OD1	1:C:103:PHE:N	2.10	0.84

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:102:ASP:OD2	1:E:147:THR:OG1	1.95	0.83

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	B	197/200 (98%)	194 (98%)	3 (2%)	0	100	100
1	C	196/200 (98%)	193 (98%)	3 (2%)	0	100	100
1	D	197/200 (98%)	194 (98%)	3 (2%)	0	100	100
1	E	197/200 (98%)	193 (98%)	4 (2%)	0	100	100
1	F	194/200 (97%)	182 (94%)	12 (6%)	0	100	100
1	J	196/200 (98%)	194 (99%)	2 (1%)	0	100	100
1	O	182/200 (91%)	178 (98%)	4 (2%)	0	100	100
2	G	237/240 (99%)	225 (95%)	12 (5%)	0	100	100
3	H	569/609 (93%)	550 (97%)	19 (3%)	0	100	100
3	I	149/609 (24%)	146 (98%)	3 (2%)	0	100	100
3	K	515/609 (85%)	493 (96%)	21 (4%)	1 (0%)	43	73
3	L	153/609 (25%)	148 (97%)	5 (3%)	0	100	100
All	All	2982/4076 (73%)	2890 (97%)	91 (3%)	1 (0%)	100	100

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
3	K	9	SER

5.3.2 Protein sidechains (i)

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	B	164/170 (96%)	164 (100%)	0	100 100
1	C	166/170 (98%)	166 (100%)	0	100 100
1	D	165/170 (97%)	165 (100%)	0	100 100
1	E	165/170 (97%)	165 (100%)	0	100 100
1	F	160/170 (94%)	160 (100%)	0	100 100
1	J	167/170 (98%)	167 (100%)	0	100 100
1	O	153/170 (90%)	153 (100%)	0	100 100
2	G	178/209 (85%)	178 (100%)	0	100 100
3	H	500/556 (90%)	500 (100%)	0	100 100
3	I	142/556 (26%)	142 (100%)	0	100 100
3	K	446/556 (80%)	444 (100%)	2 (0%)	84 86
3	L	126/556 (23%)	126 (100%)	0	100 100
All	All	2532/3623 (70%)	2530 (100%)	2 (0%)	87 90

All (2) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	K	10	LYS
3	K	14	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 18 such sidechains are listed below:

Mol	Chain	Res	Type
3	L	438	ASN
3	L	558	ASN
3	L	516	ASN
3	H	326	HIS
3	K	606	ASN

5.3.3 RNA

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
4	M	46/54 (85%)	14 (30%)	1 (2%)
5	N	36/60 (60%)	14 (38%)	0
All	All	82/114 (71%)	28 (34%)	1 (1%)

5 of 28 RNA backbone outliers are listed below:

Mol	Chain	Res	Type
4	M	-7	U
4	M	-6	A
4	M	-1	U
4	M	11	U
4	M	12	G

All (1) RNA pucker outliers are listed below:

Mol	Chain	Res	Type
4	M	34	C

5.4 Non-standard residues in protein, DNA, RNA chains

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 9 ligands modelled in this entry, 9 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.

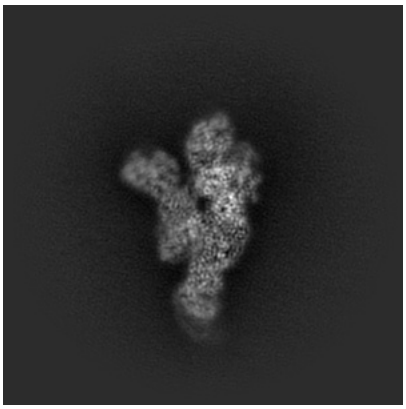
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-67552. These allow visual inspection of the internal detail of the map and identification of artifacts.

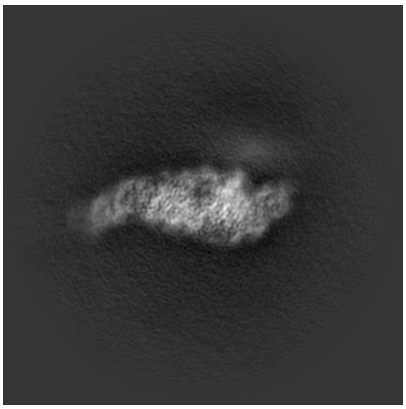
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

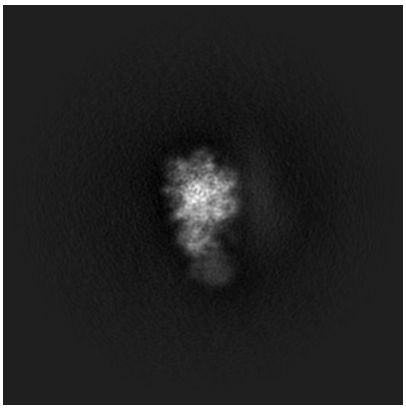
6.1.1 Primary map



X

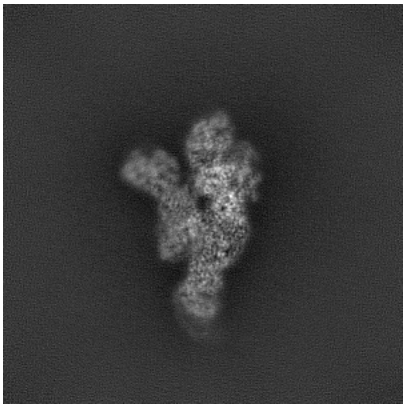


Y

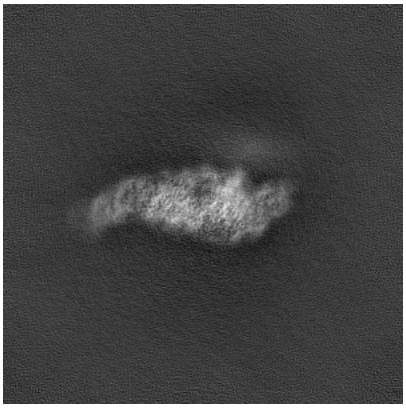


Z

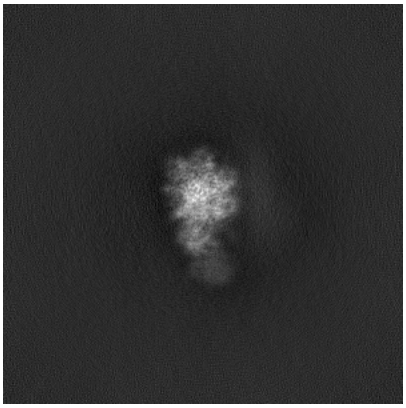
6.1.2 Raw map



X



Y

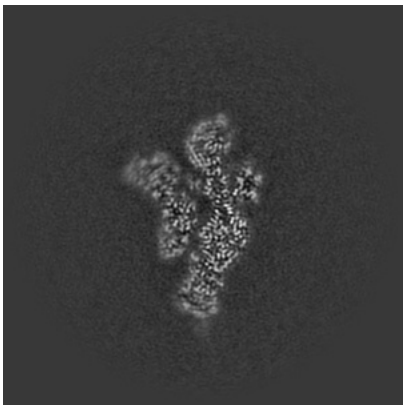


Z

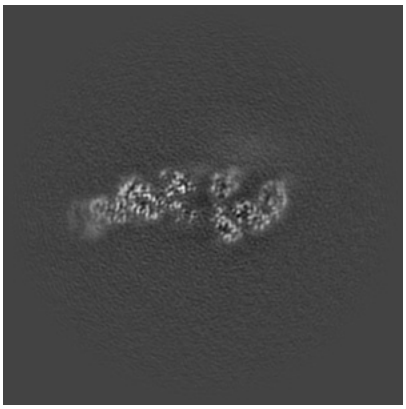
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

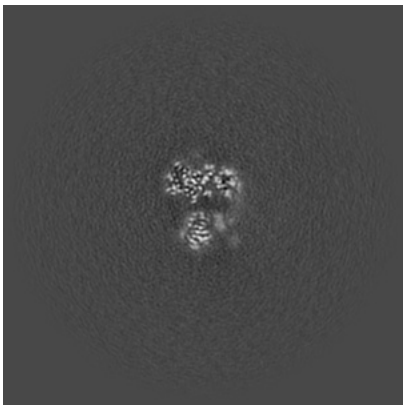
6.2.1 Primary map



X Index: 208

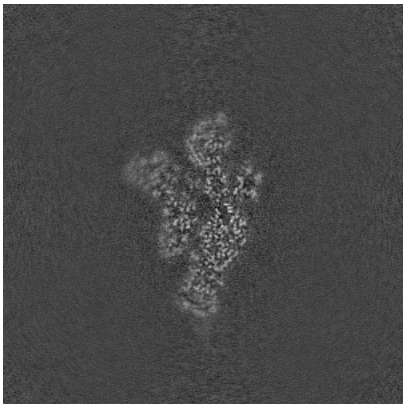


Y Index: 208

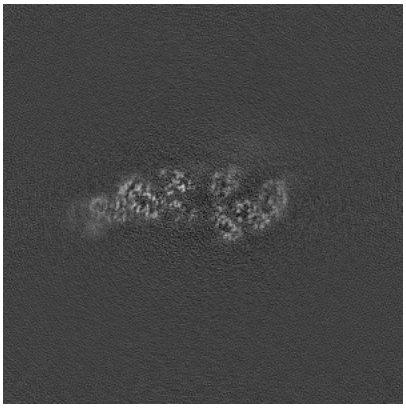


Z Index: 208

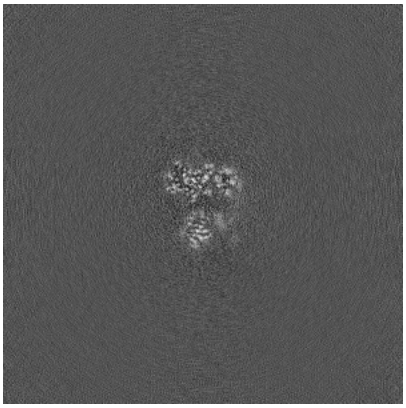
6.2.2 Raw map



X Index: 208



Y Index: 208

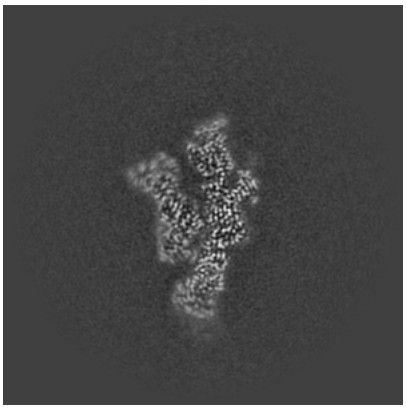


Z Index: 208

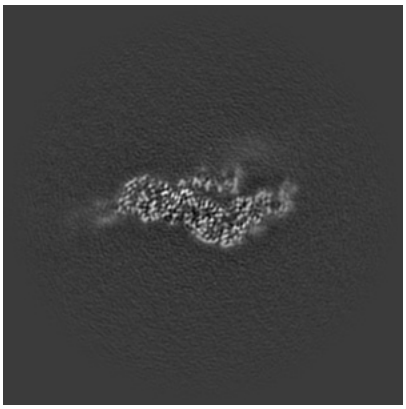
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

6.3.1 Primary map



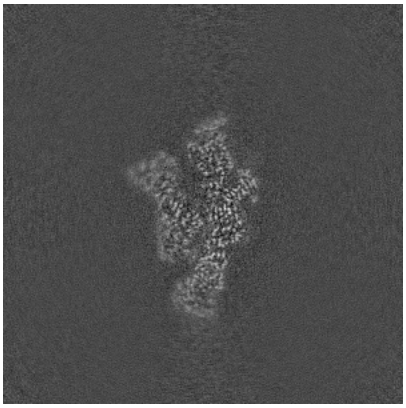
X Index: 202



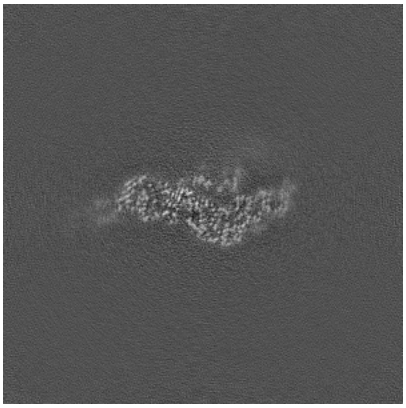
Y Index: 223

Z Index: 231

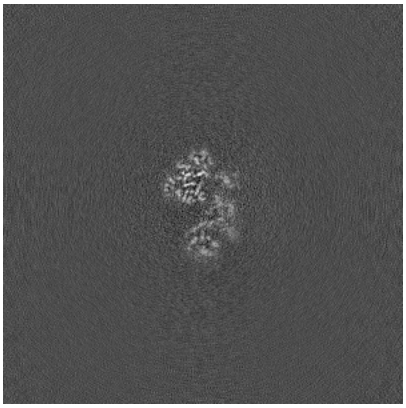
6.3.2 Raw map



X Index: 202



Y Index: 222

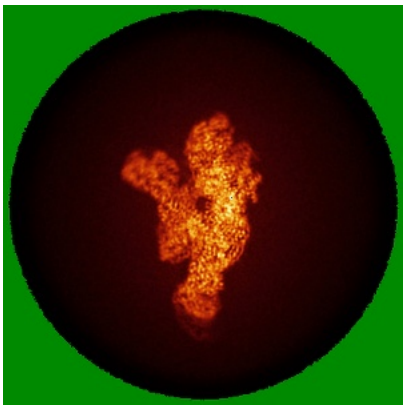


Z Index: 219

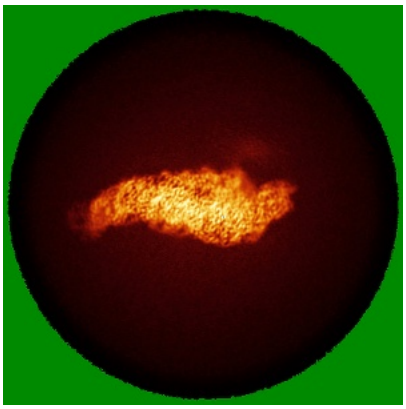
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

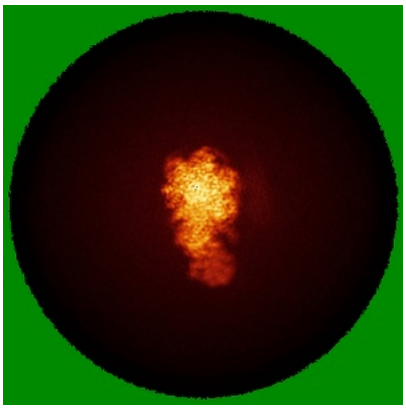
6.4.1 Primary map



X

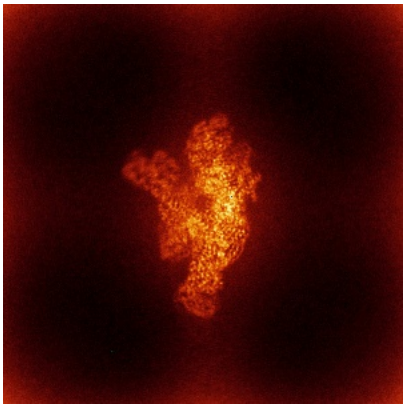


Y

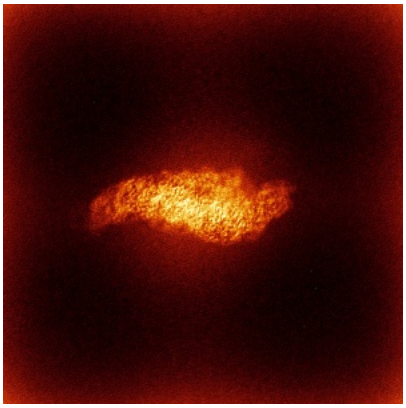


Z

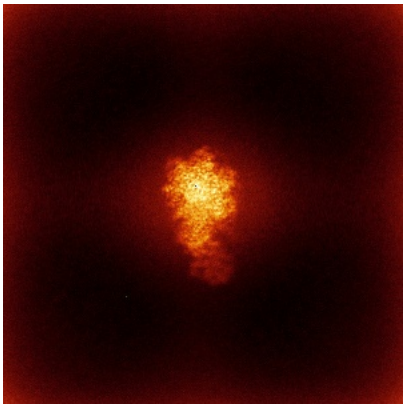
6.4.2 Raw map



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.

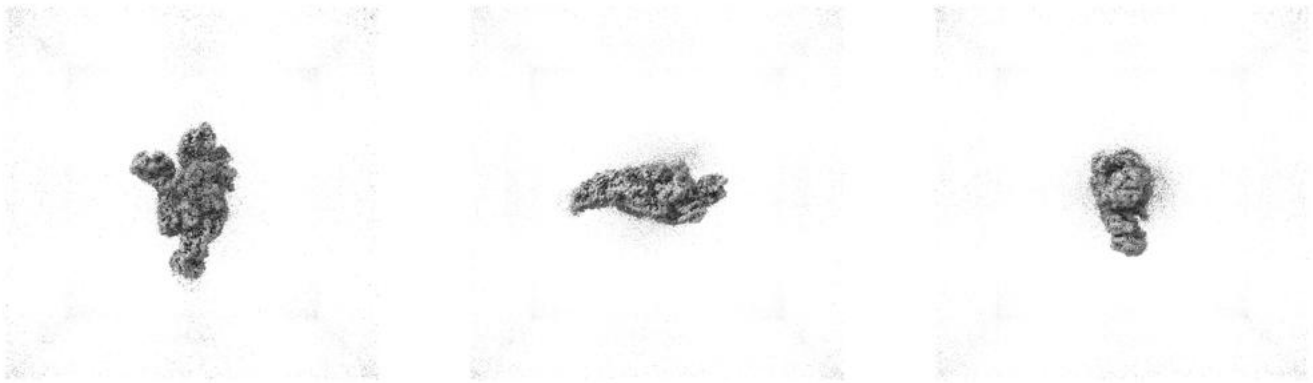
6.5 Orthogonal surface views (i)

6.5.1 Primary map

X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.11. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



X Y Z

These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

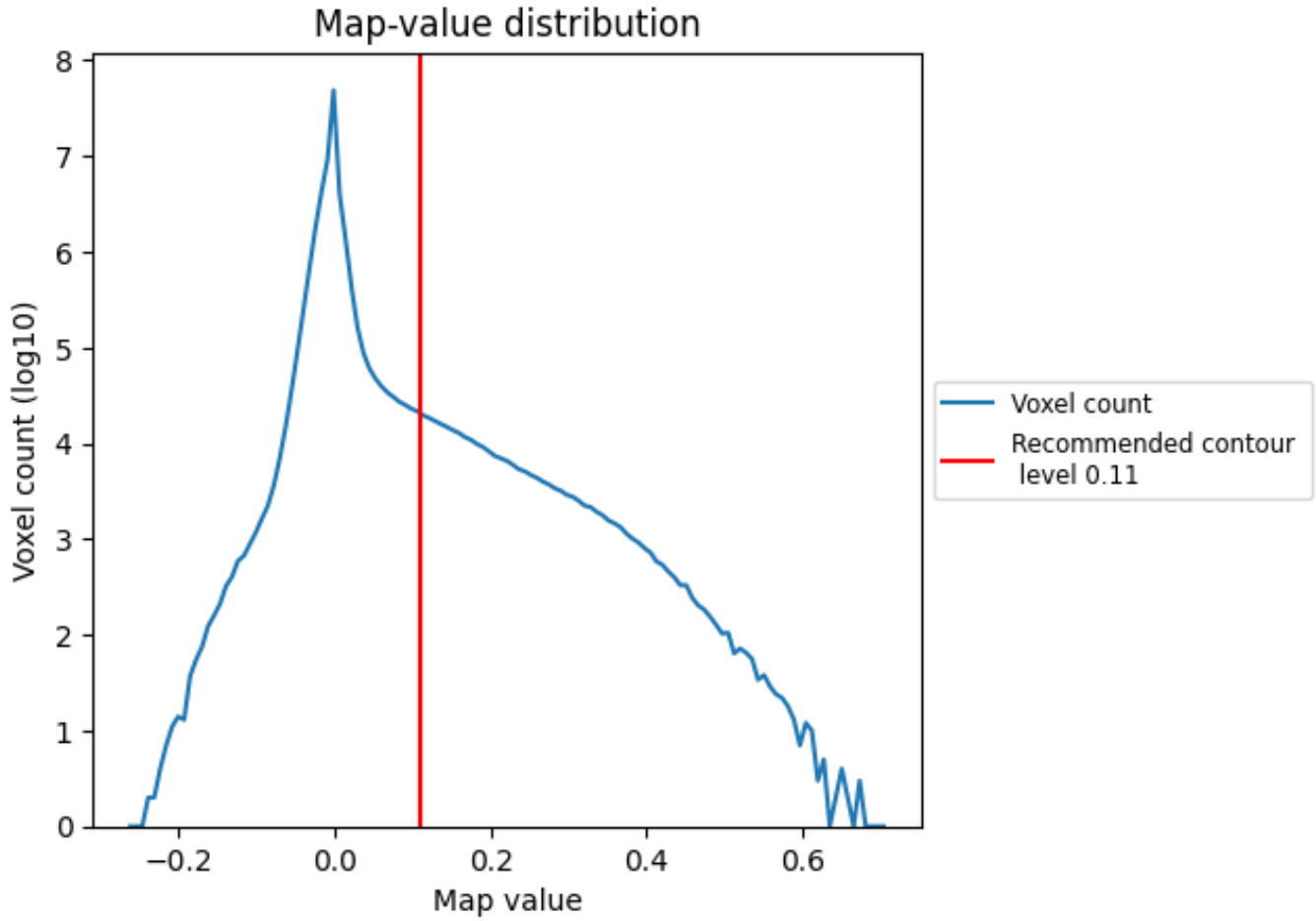
6.6 Mask visualisation (i)

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

7 Map analysis

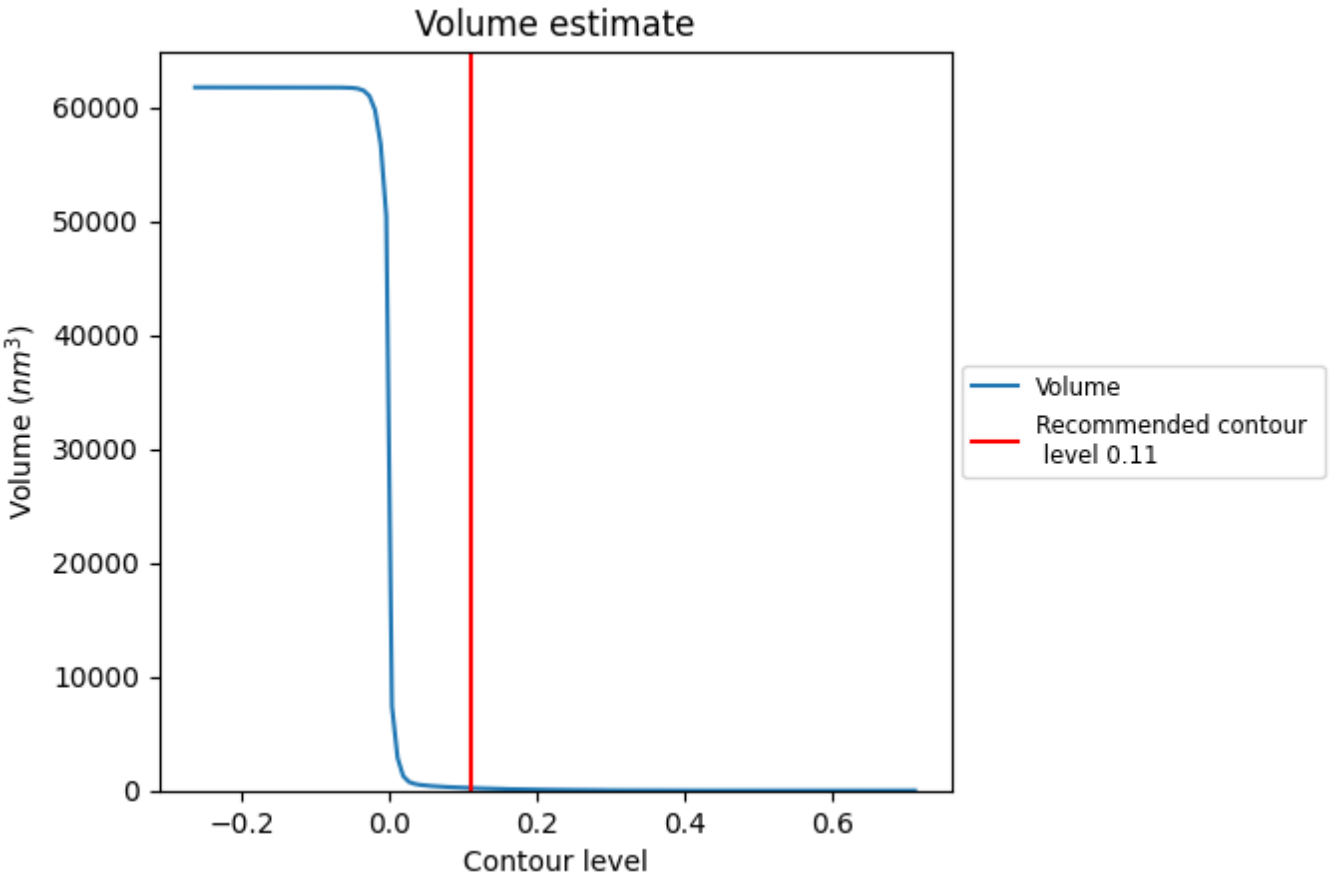
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

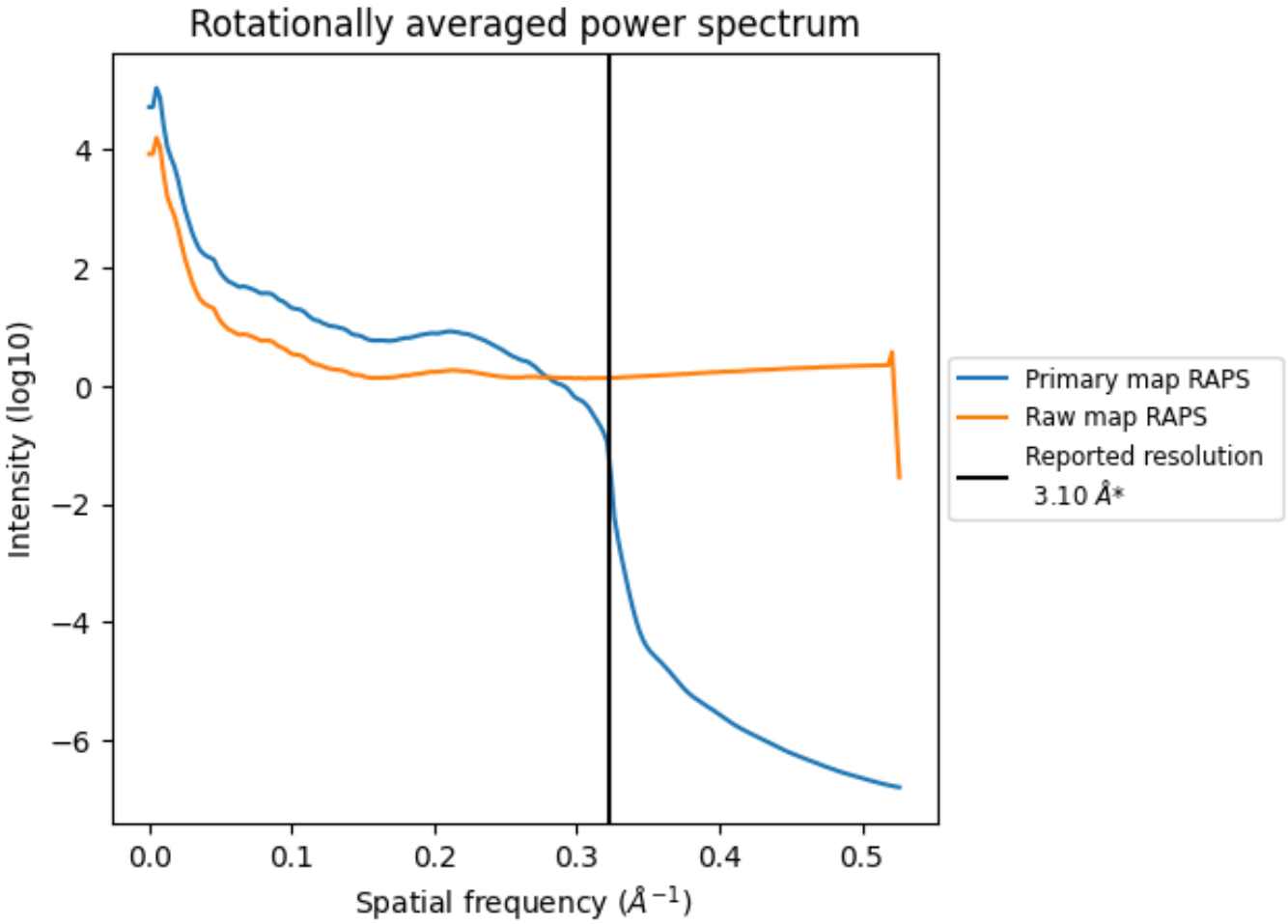
7.2 Volume estimate (i)



The volume at the recommended contour level is 225 nm^3; this corresponds to an approximate mass of 203 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 (i)

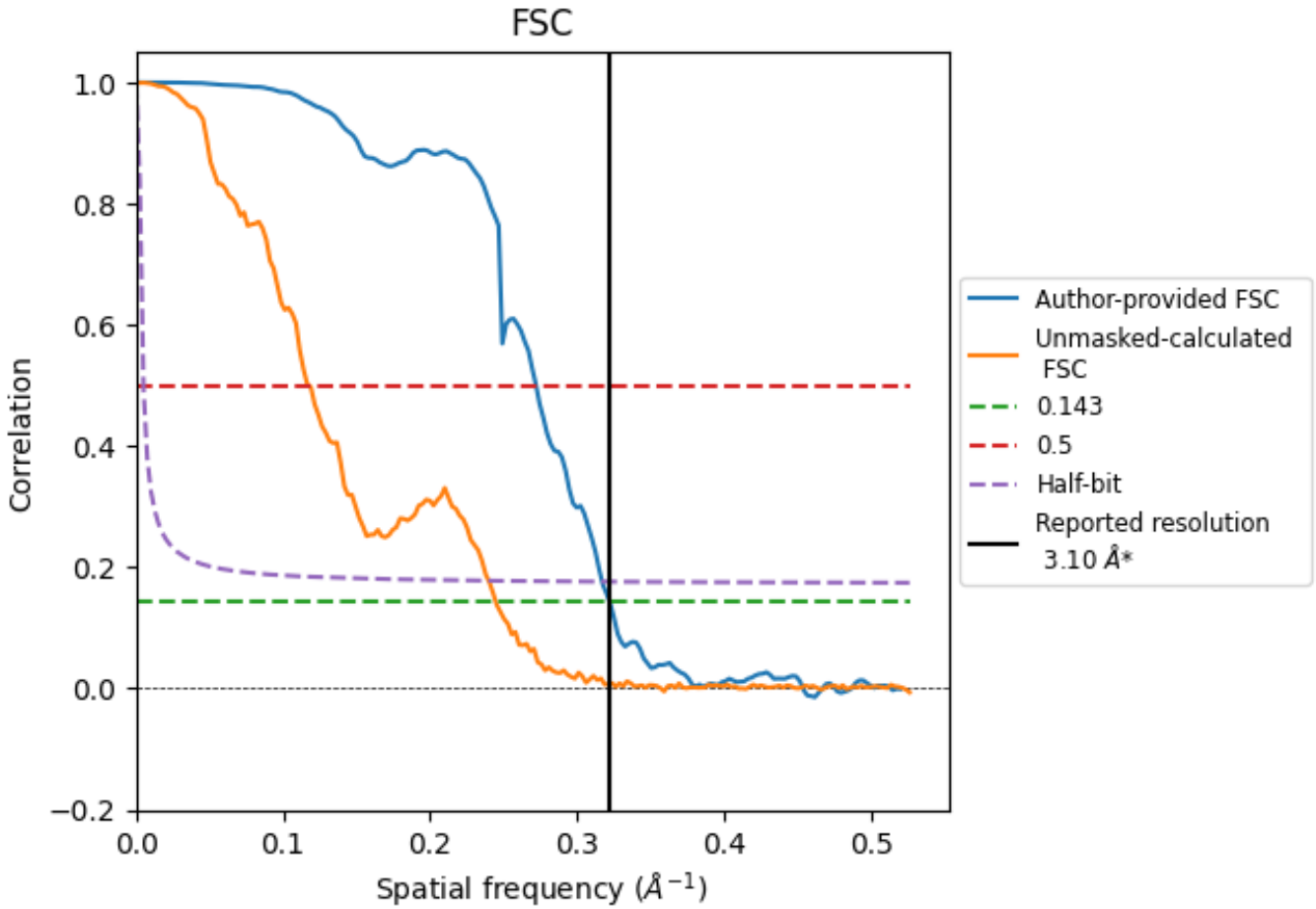


*Reported resolution corresponds to spatial frequency of 0.323 Å^{-1}

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.323 Å^{-1}

8.2 Resolution estimates (i)

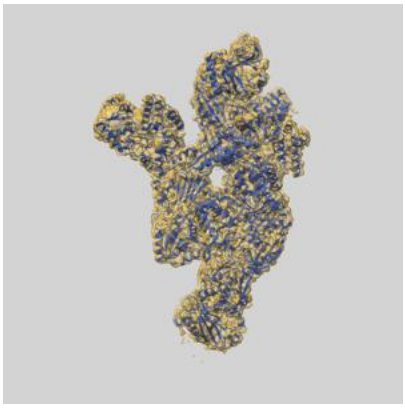
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.10	3.67	3.15
Unmasked-calculated*	4.08	8.50	4.17

*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 4.08 differs from the reported value 3.1 by more than 10 %

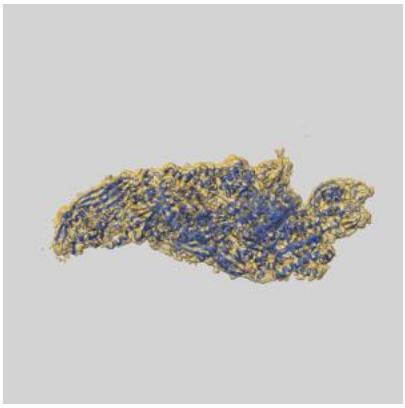
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-67552 and PDB model 21BH. Per-residue inclusion information can be found in section 3 on page 6.

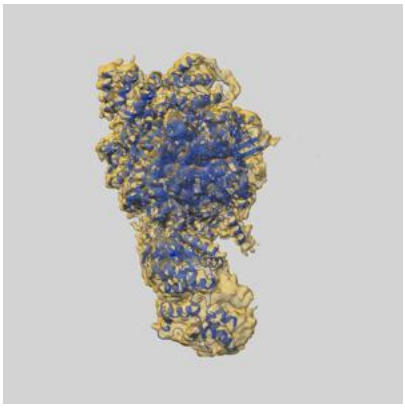
9.1 Map-model overlay



X



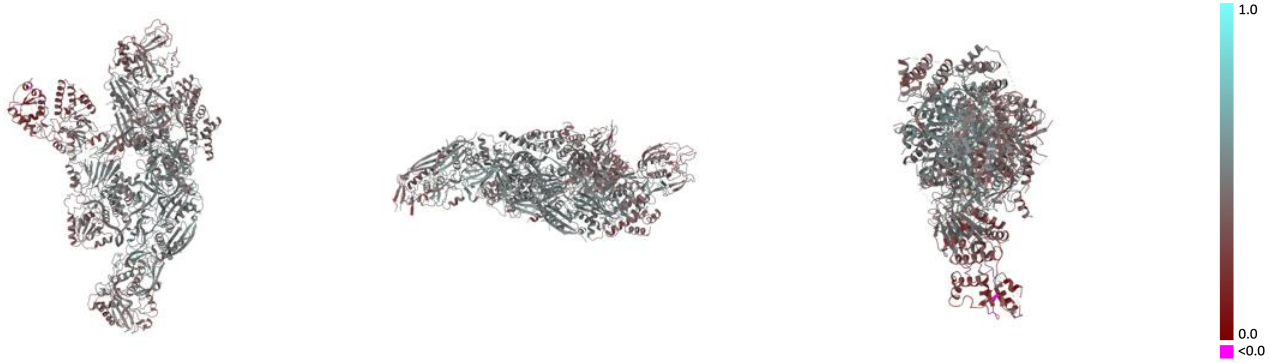
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.11 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.

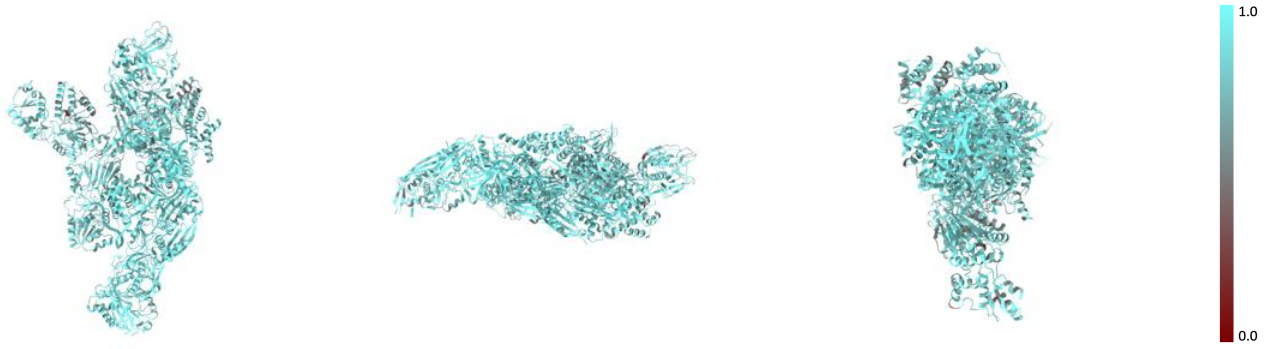
9.2 Q-score mapped to coordinate model



X Y Z

The images above show the model with each residue coloured according 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.

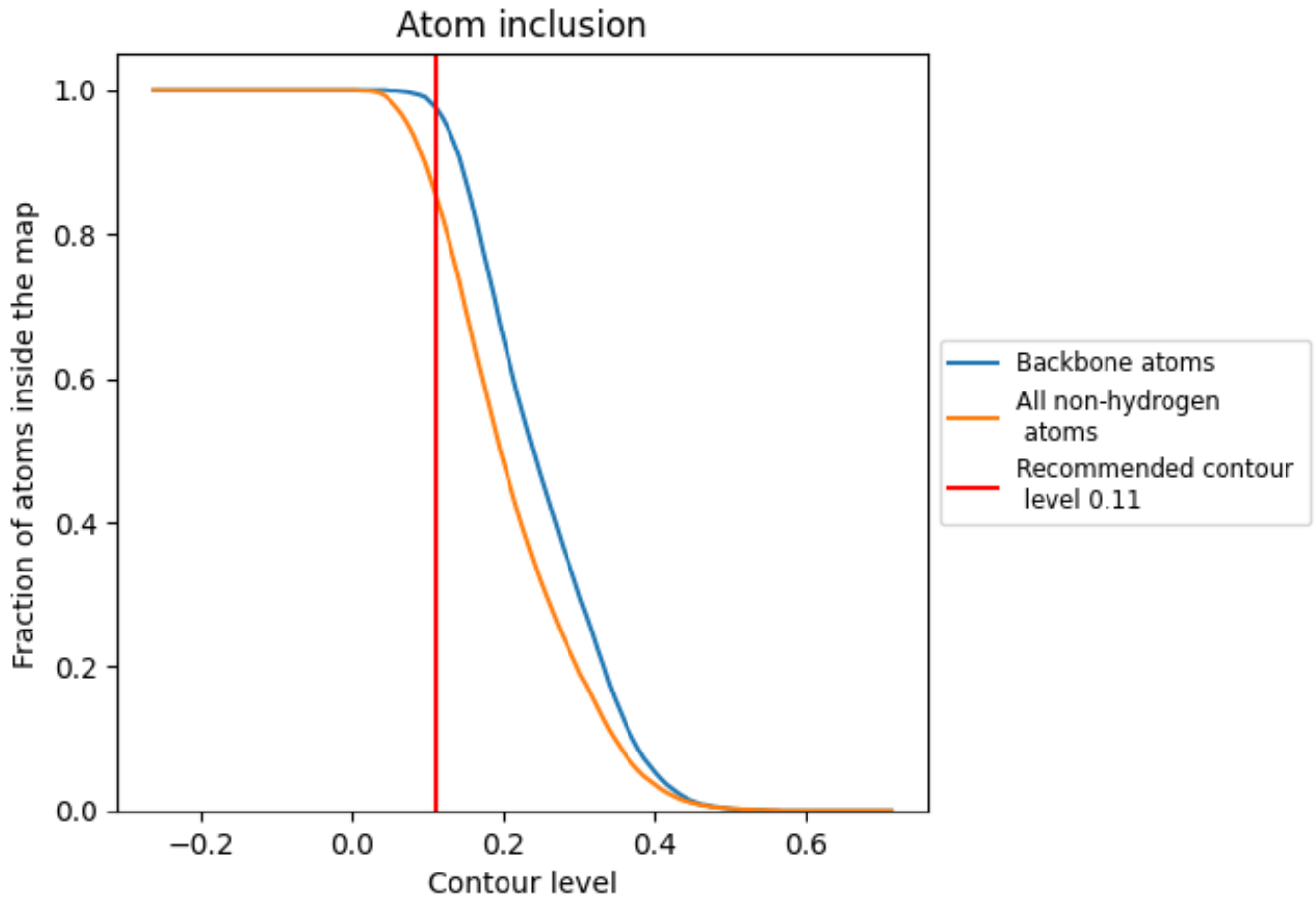
9.3 Atom inclusion mapped to coordinate model



X Y Z

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 (0.11).

9.4 Atom inclusion (i)



At the recommended contour level, 98% of all backbone atoms, 85% 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 (0.11) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.8530	0.4530
B	0.8920	0.4710
C	0.9170	0.5130
D	0.8980	0.5240
E	0.8760	0.5140
F	0.8530	0.4820
G	0.8260	0.3990
H	0.8330	0.4500
I	0.8590	0.4420
J	0.9100	0.5240
K	0.7870	0.3750
L	0.7600	0.4120
M	0.9720	0.5050
N	0.9610	0.4960
O	0.8360	0.4120