wwPDB EM Validation Summary Report

Oct 19, 2026 – 07:22 pm BST

PDB ID	:	32ZR / pdb_000032zr
EMDB ID	:	EMD-59288
Title	:	B. subtilis RNA Polymerase delta omega variant with SigA
Authors	:	Loll, B.; Hilal, T.
Deposited on	:	2026-08-03
Resolution	:	2.80 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:

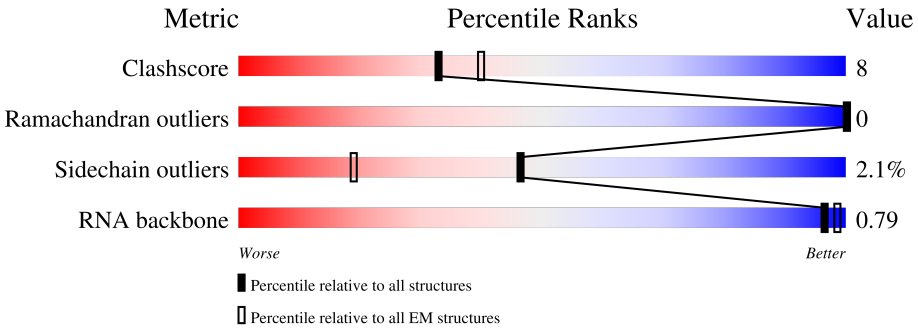
EMDB validation analysis	:	0.0.1.dev133
MolProbity	:	4-5-2 with Phenix2.2
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.51

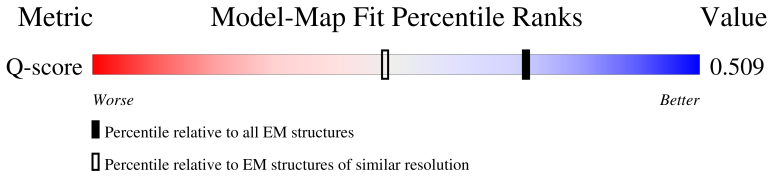
1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.80 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.





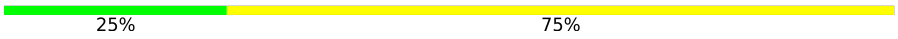
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
RNA backbone	8273	3508	-
Q-score	-	25397	11806 (2.30 - 3.30)

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	E	69	77% 22% ·
2	U	314	17% 77% 14% · 8%
2	V	314	56% 17% 27%

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Mol	Chain	Length	Quality of chain
3	X	1193	
4	Y	1209	
5	R	5	
6	G	371	
7	K	60	
8	L	63	
9	D	173	

2 Entry composition (i)

There are 11 unique types of molecules in this entry. The entry contains 28710 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 DNA-directed RNA polymerase subunit epsilon.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	E	69	Total	C	N	O	S	0	0
			580	373	92	114	1		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
E	42	LEU	PHE	conflict	UNP A0A9W5LJC9

- Molecule 2 is a protein called DNA-directed RNA polymerase subunit alpha.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	U	290	Total	C	N	O	S	3	0
			2279	1425	395	455	4		
2	V	229	Total	C	N	O	S	0	0
			1761	1110	295	355	1		

- Molecule 3 is a protein called DNA-directed RNA polymerase subunit beta.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	X	1157	Total	C	N	O	S	1	0
			9142	5735	1599	1774	34		

- Molecule 4 is a protein called DNA-directed RNA polymerase subunit beta'.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	Y	1181	Total	C	N	O	S	7	0
			9355	5898	1673	1740	44		

There are 10 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
Y	1200	HIS	-	expression tag	UNP P37871
Y	1201	HIS	-	expression tag	UNP P37871

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Chain	Residue	Modelled	Actual	Comment	Reference
Y	1202	HIS	-	expression tag	UNP P37871
Y	1203	HIS	-	expression tag	UNP P37871
Y	1204	HIS	-	expression tag	UNP P37871
Y	1205	HIS	-	expression tag	UNP P37871
Y	1206	HIS	-	expression tag	UNP P37871
Y	1207	HIS	-	expression tag	UNP P37871
Y	1208	HIS	-	expression tag	UNP P37871
Y	1209	HIS	-	expression tag	UNP P37871

- Molecule 5 is a RNA chain called RNA.

Mol	Chain	Residues	Atoms					AltConf	Trace
5	R	5	Total	C	N	O	P	0	0
			105	47	17	36	5		

- Molecule 6 is a protein called RNA polymerase sigma factor SigA.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	G	278	Total	C	N	O	S	0	0
			2263	1418	408	431	6		

- Molecule 7 is a DNA chain called SigA controlled promoter Pveg.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	K	60	Total	C	N	O	P	0	0
			1223	587	226	350	60		

There are 23 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	1	DA	-	expression tag	GB 2731940075
K	2	DT	-	expression tag	GB 2731940075
K	3	DC	-	expression tag	GB 2731940075
K	4	DC	-	expression tag	GB 2731940075
K	5	DA	-	expression tag	GB 2731940075
K	6	DC	-	expression tag	GB 2731940075
K	7	DC	-	expression tag	GB 2731940075
K	8	DT	-	expression tag	GB 2731940075
K	9	DC	-	expression tag	GB 2731940075
K	10	DA	-	expression tag	GB 2731940075
K	11	DG	-	expression tag	GB 2731940075
K	12	DA	-	expression tag	GB 2731940075

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Chain	Residue	Modelled	Actual	Comment	Reference
K	13	DT	-	expression tag	GB 2731940075
K	14	DG	-	expression tag	GB 2731940075
K	15	DT	-	expression tag	GB 2731940075
K	16	DA	-	expression tag	GB 2731940075
K	17	DA	-	expression tag	GB 2731940075
K	18	DA	-	expression tag	GB 2731940075
K	19	DT	-	expression tag	GB 2731940075
K	20	DA	-	expression tag	GB 2731940075
K	21	DA	-	expression tag	GB 2731940075
K	22	DC	-	expression tag	GB 2731940075
K	23	DA	-	expression tag	GB 2731940075

- Molecule 8 is a DNA chain called SigA controlled promoter Pveg.

Mol	Chain	Residues	Atoms					AltConf	Trace
8	L	60	Total	C	N	O	P	0	0
			1244	596	226	362	60		

- Molecule 9 is a protein called DNA-directed RNA polymerase subunit delta.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	D	91	Total	C	N	O	S	0	0
			755	484	121	149	1		

- Molecule 10 is MAGNESIUM ION (CCD ID: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		AltConf
10	Y	1	Total	Mg	0
			1	1	

- Molecule 11 is ZINC ION (CCD ID: ZN) (formula: Zn).

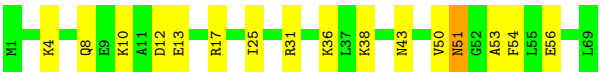
Mol	Chain	Residues	Atoms		AltConf
11	Y	2	Total	Zn	0
			2	2	

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: DNA-directed RNA polymerase subunit epsilon

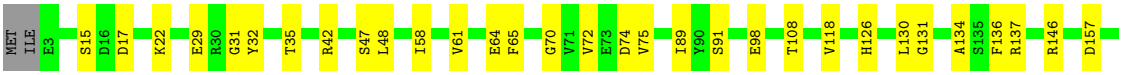
Chain E:

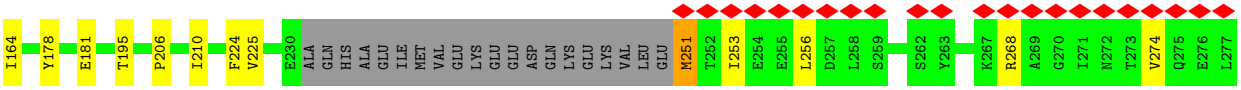


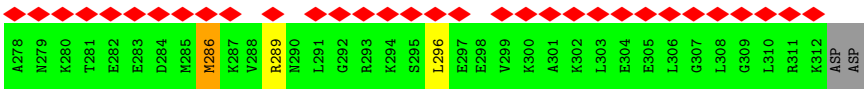
• Molecule 2: DNA-directed RNA polymerase subunit alpha

Chain U:







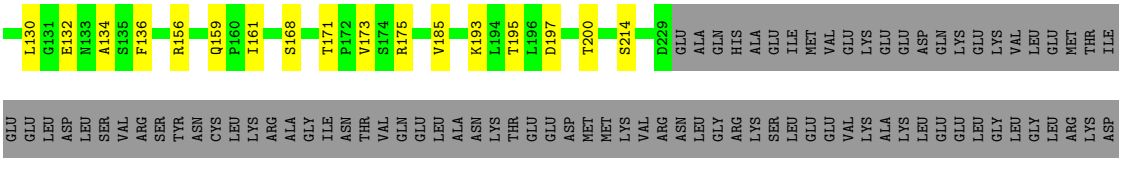


• Molecule 2: DNA-directed RNA polymerase subunit alpha

Chain V:



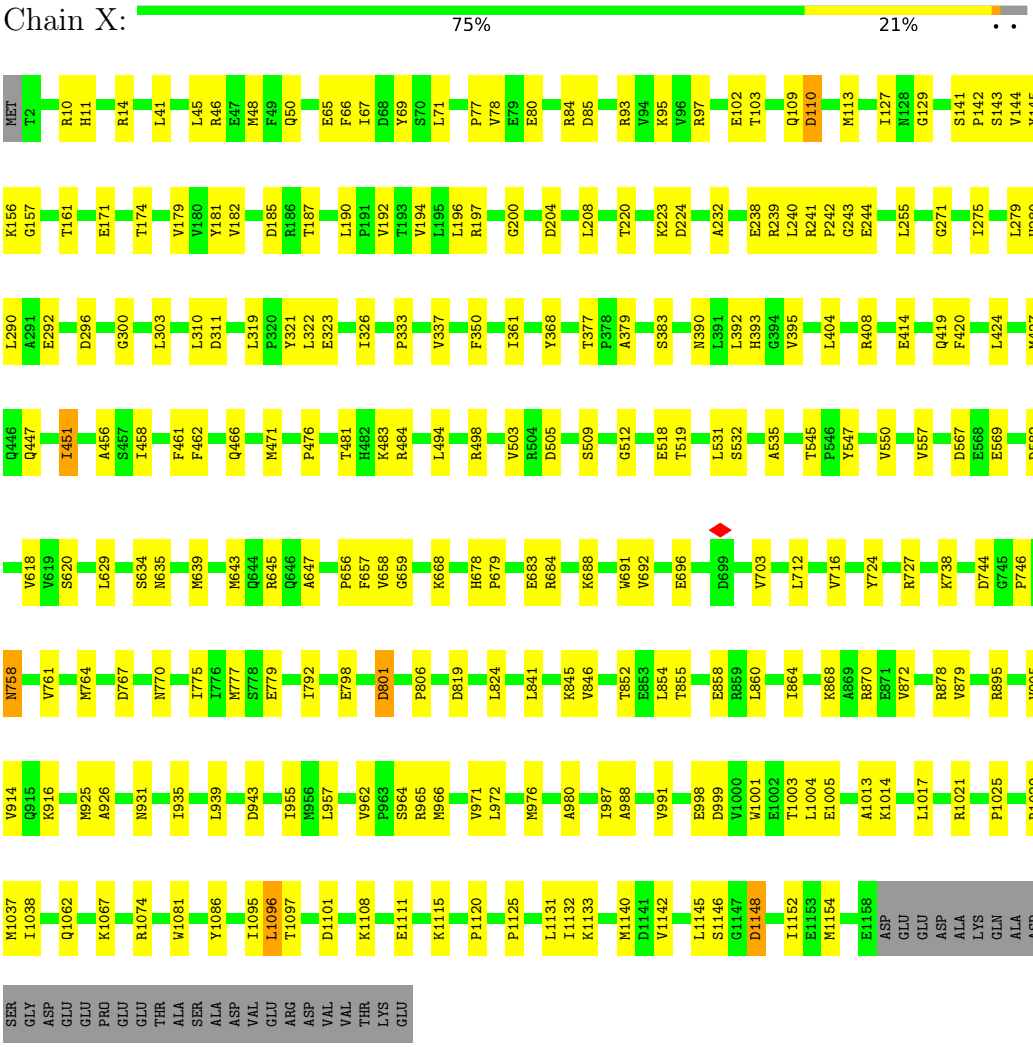






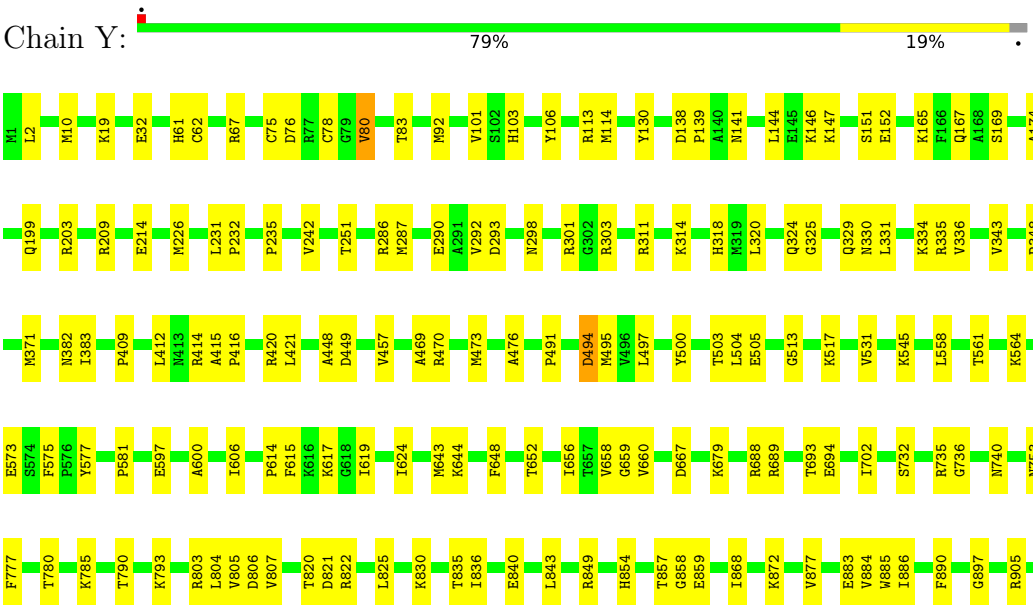
• Molecule 3: DNA-directed RNA polymerase subunit beta

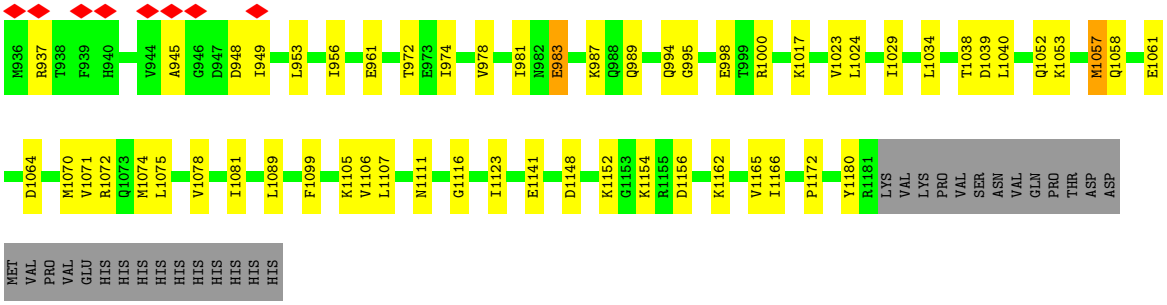
Chain X:



• Molecule 4: DNA-directed RNA polymerase subunit beta'

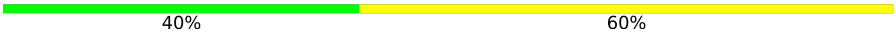
Chain Y:





• Molecule 5: RNA

Chain R:

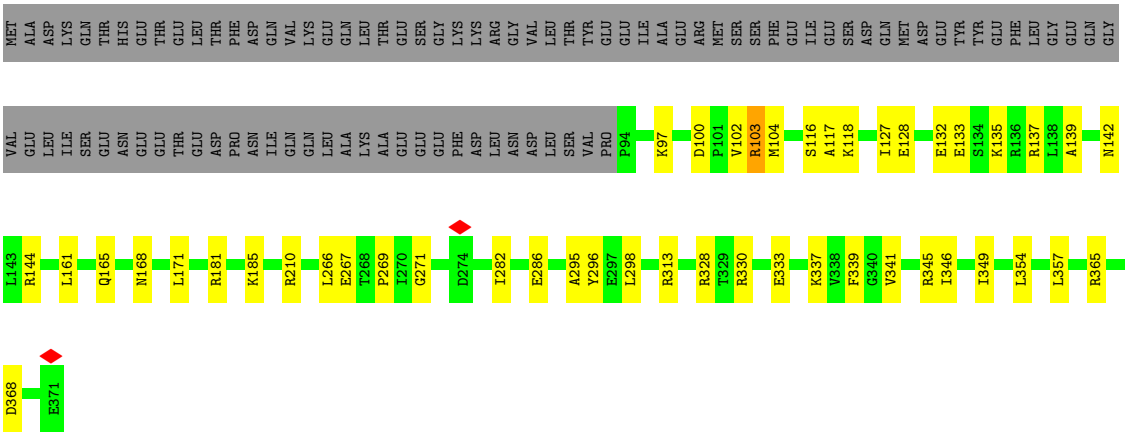




• Molecule 6: RNA polymerase sigma factor SigA

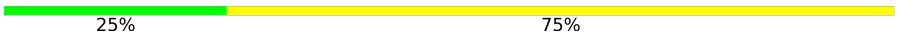
Chain G:

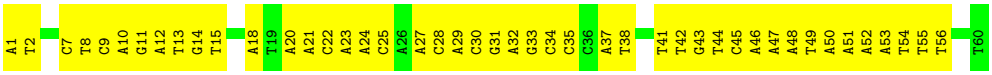




• Molecule 7: SigA controlled promoter Pveg

Chain K:

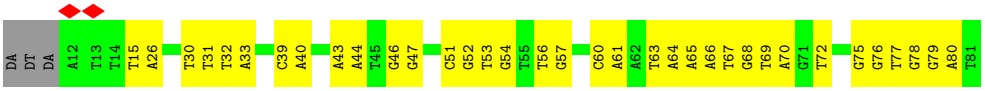




• Molecule 8: SigA controlled promoter Pveg

Chain L:

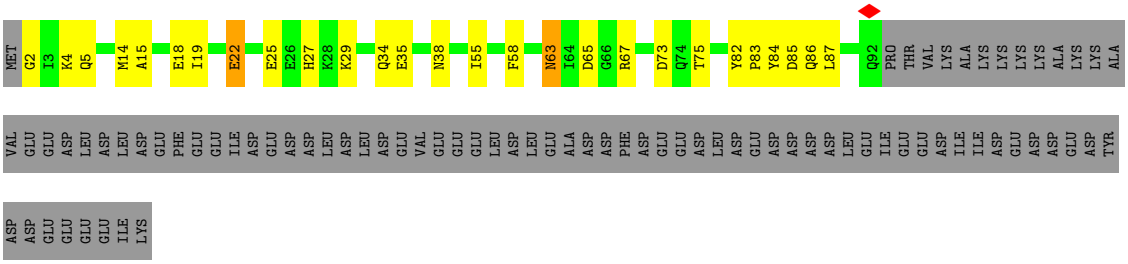




• Molecule 9: DNA-directed RNA polymerase subunit delta

Chain D:




MET G2 I3 K4 Q5 M14 A15 E18 I19 E22 E25 E26 H27 K28 K29 Q34 E35 N38 I55 F58 N63 I64 D65 G66 R67 D73 Q74 T75 Y82 P83 Y84 D85 Q86 L87 Q92 PRO THR VAL LYS ALA LYS LYS LYS LYS ALA LYS LYS ALA
VAL GLU GLU ASP LEU ASP LEU ASP GLU PHE GLU GLU ILE ASP GLU ASP ASP LEU ASP LEU ASP GLU VAL GLU GLU GLU LEU ASP LEU GLU ALA ASP ASP PHE ASP GLU GLU ASP LEU ASP GLU ASP ASP ASP ASP LEU GLU ILE GLU GLU ASP ILE ILE ASP GLU ASP ASP GLU ASP TYR
ASP ASP GLU GLU GLU GLU ILE LYS

4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	82187	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/Å^2$)	42	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	96000	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	2.283	Depositor
Minimum map value	0.000	Depositor
Average map value	0.011	Depositor
Map value standard deviation	0.065	Depositor
Recommended contour level	0.22	Depositor
Map size (Å)	314.496, 314.496, 314.496	wwPDB
Map dimensions	384, 384, 384	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.819, 0.819, 0.819	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, MG

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	E	0.10	0/589	0.23	0/789
2	U	0.09	0/2305	0.22	0/3116
2	V	0.11	0/1787	0.27	0/2430
3	X	0.12	0/9301	0.27	0/12573
4	Y	0.10	0/9505	0.24	0/12817
5	R	0.22	0/116	0.68	0/178
6	G	0.10	0/2294	0.24	0/3083
7	K	0.68	0/1373	0.87	0/2113
8	L	0.66	0/1397	0.90	0/2158
9	D	0.12	0/771	0.35	0/1041
All	All	0.23	0/29438	0.38	0/40298

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

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	E	580	0	585	12	0
2	U	2279	0	2319	35	0
2	V	1761	0	1779	37	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	X	9142	0	9093	174	0
4	Y	9355	0	9601	148	0
5	R	105	0	53	3	0
6	G	2263	0	2309	43	0
7	K	1223	0	677	40	0
8	L	1244	0	683	32	0
9	D	755	0	728	23	0
10	Y	1	0	0	0	0
11	Y	2	0	0	0	0
All	All	28710	0	27827	475	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 8.

The worst 5 of 475 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:U:42:ARG:HE	2:V:35:THR:HG22	1.38	0.87
3:X:48:MET:HE1	3:X:420:PHE:HB3	1.65	0.79
3:X:67:ILE:HD12	3:X:95:LYS:HG2	1.70	0.73
4:Y:114:MET:HG2	4:Y:226:MET:HE3	1.70	0.73
4:Y:857:THR:OG1	4:Y:859:GLU:OE1	2.07	0.72

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	E	67/69 (97%)	66 (98%)	1 (2%)	0	100	100
2	U	289/314 (92%)	285 (99%)	4 (1%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
2	V	227/314 (72%)	218 (96%)	9 (4%)	0	100	100
3	X	1156/1193 (97%)	1122 (97%)	34 (3%)	0	100	100
4	Y	1186/1209 (98%)	1162 (98%)	24 (2%)	0	100	100
6	G	276/371 (74%)	274 (99%)	2 (1%)	0	100	100
9	D	89/173 (51%)	88 (99%)	1 (1%)	0	100	100
All	All	3290/3643 (90%)	3215 (98%)	75 (2%)	0	100	100

There are no Ramachandran outliers to report.

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	E	65/65 (100%)	64 (98%)	1 (2%)	57	84
2	U	254/273 (93%)	249 (98%)	5 (2%)	48	80
2	V	197/273 (72%)	190 (96%)	7 (4%)	31	66
3	X	998/1026 (97%)	969 (97%)	29 (3%)	37	73
4	Y	1016/1037 (98%)	1001 (98%)	15 (2%)	57	84
6	G	244/329 (74%)	242 (99%)	2 (1%)	73	90
9	D	81/159 (51%)	79 (98%)	2 (2%)	42	76
All	All	2855/3162 (90%)	2794 (98%)	61 (2%)	46	79

5 of 61 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
3	X	727	ARG
4	Y	1105	LYS
3	X	879	VAL
4	Y	1078	VAL
9	D	22	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 27 such sidechains are listed below:

Mol	Chain	Res	Type
4	Y	394	GLN
4	Y	542	ASN
9	D	74	GLN
4	Y	478	ASN
4	Y	632	HIS

5.3.3 RNA (i)

Mol	Chain	Analysed	Backbone Outliers	Pucker Outliers
5	R	4/5 (80%)	0	0

There are no RNA backbone outliers to report.

There are no RNA pucker outliers to report.

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 3 ligands modelled in this entry, 3 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.

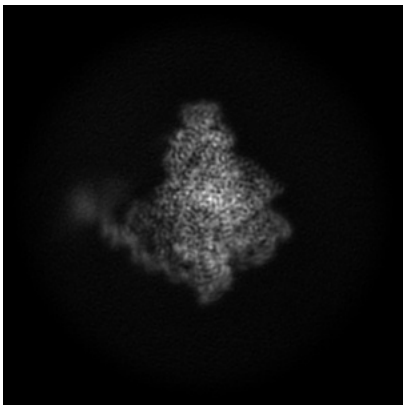
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-59288. These allow visual inspection of the internal detail of the map and identification of artifacts.

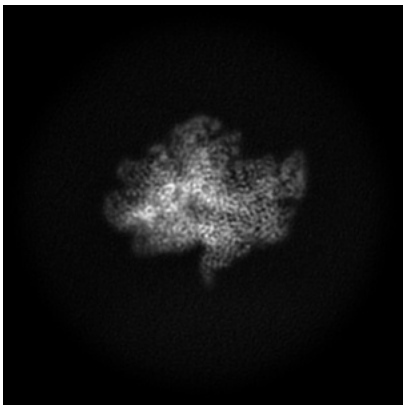
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

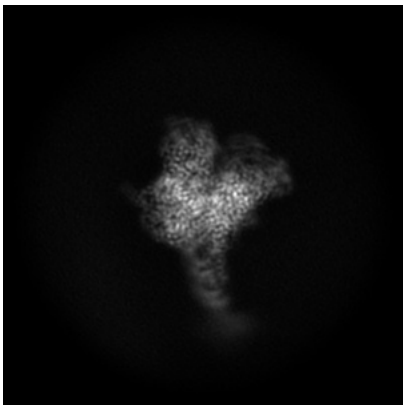
6.1.1 Primary map



X

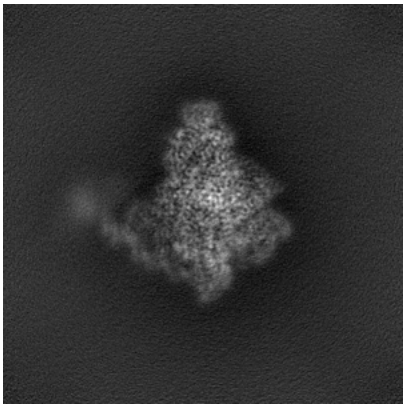


Y

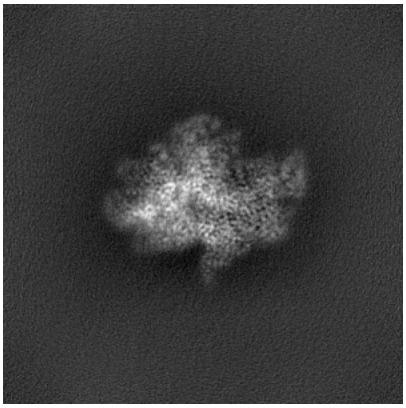


Z

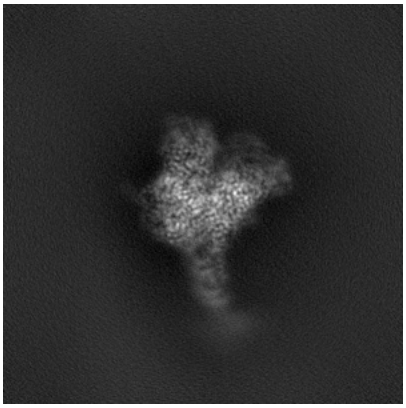
6.1.2 Raw map



X



Y

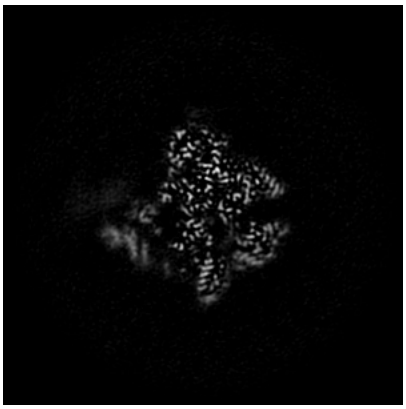


Z

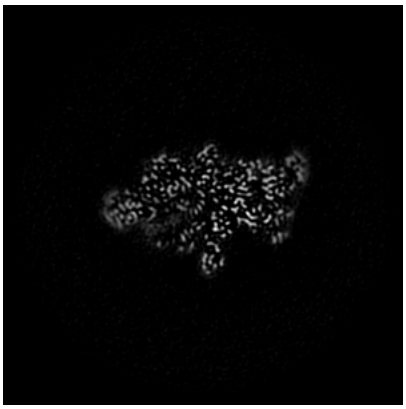
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

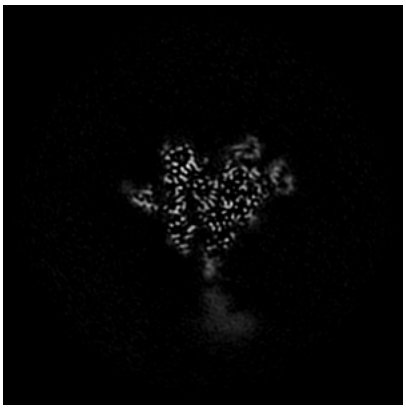
6.2.1 Primary map



X Index: 192

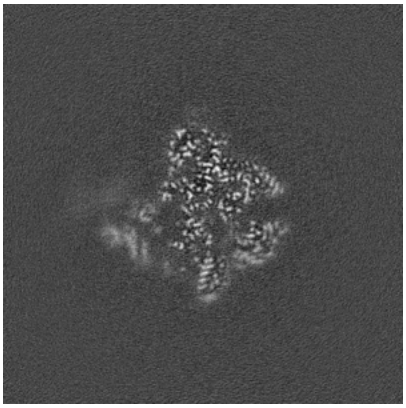


Y Index: 192

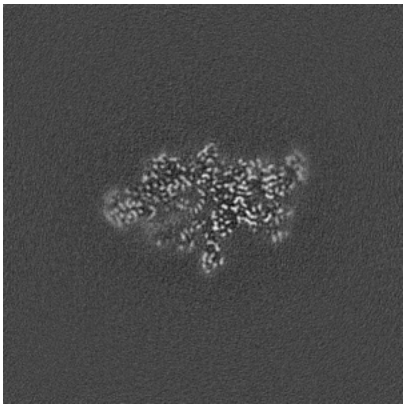


Z Index: 192

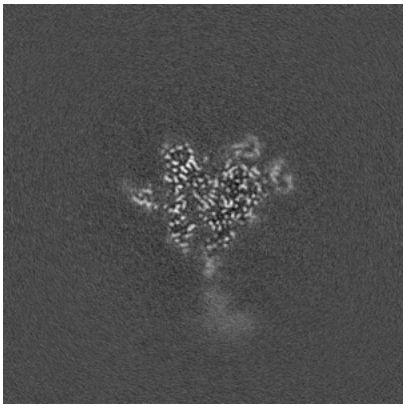
6.2.2 Raw map



X Index: 192



Y Index: 192

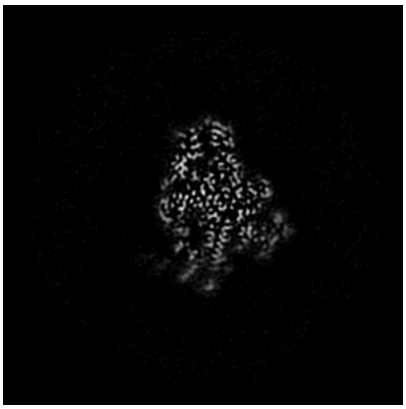


Z Index: 192

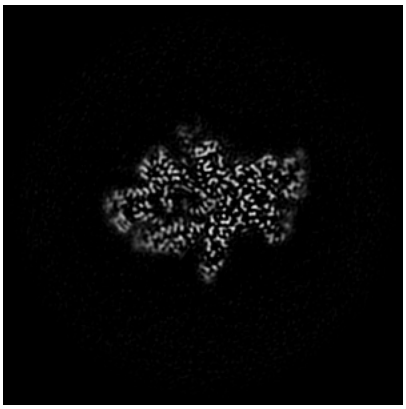
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

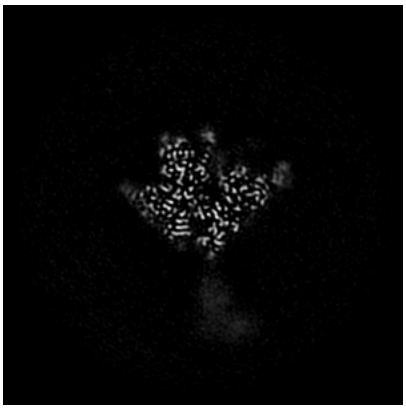
6.3.1 Primary map



X Index: 168

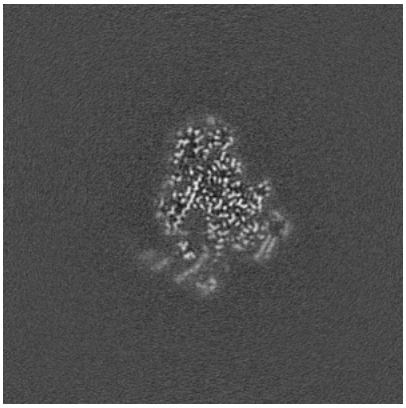


Y Index: 199

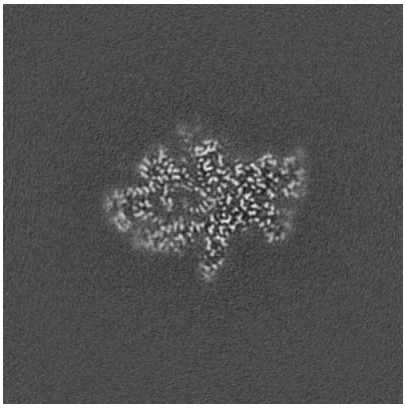


Z Index: 200

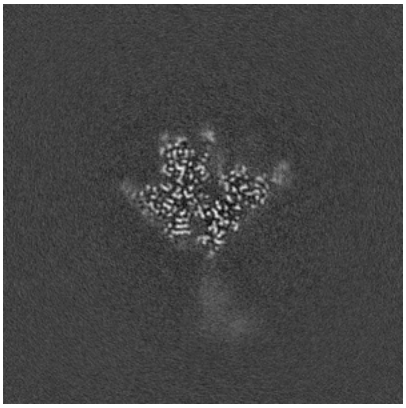
6.3.2 Raw map



X Index: 171



Y Index: 199

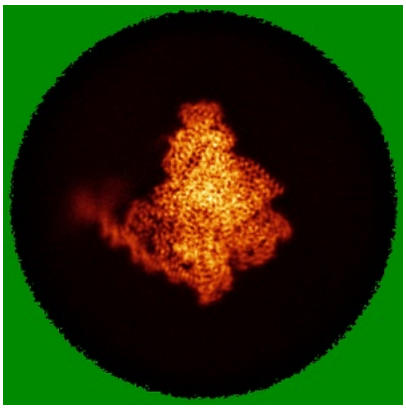


Z Index: 200

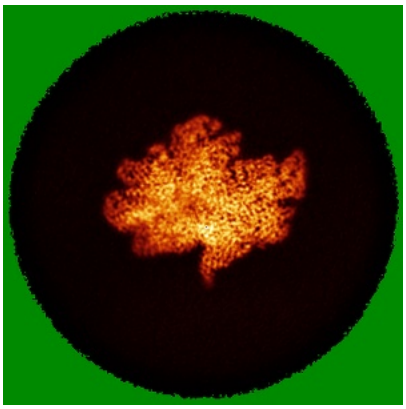
The images above show the largest variance slices of the map in three orthogonal directions.

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

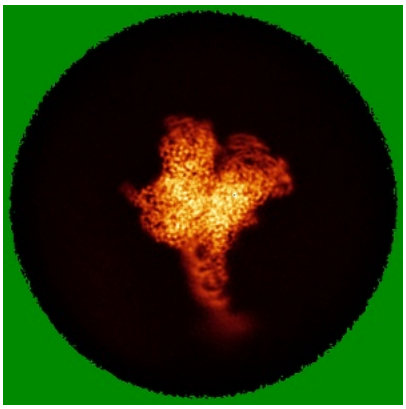
6.4.1 Primary map



X

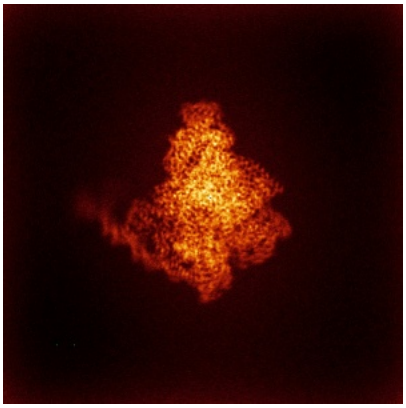


Y

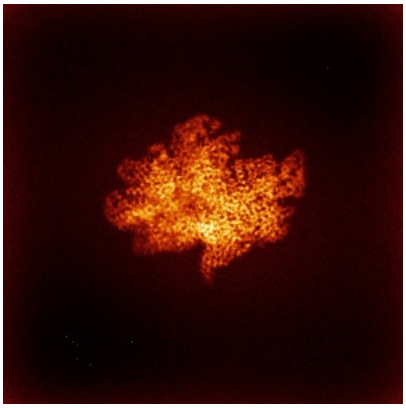


Z

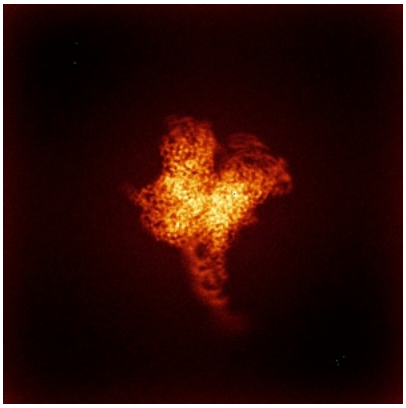
6.4.2 Raw map



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)

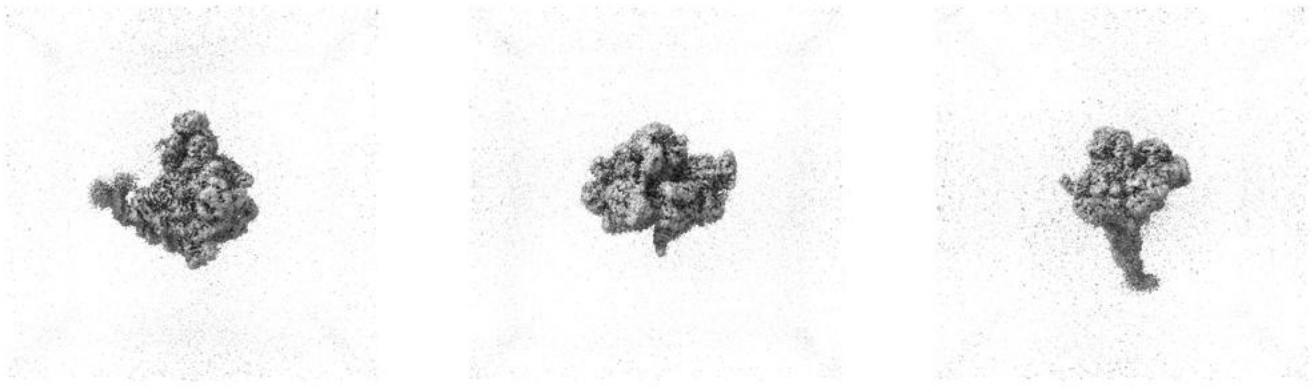
6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.22. 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.

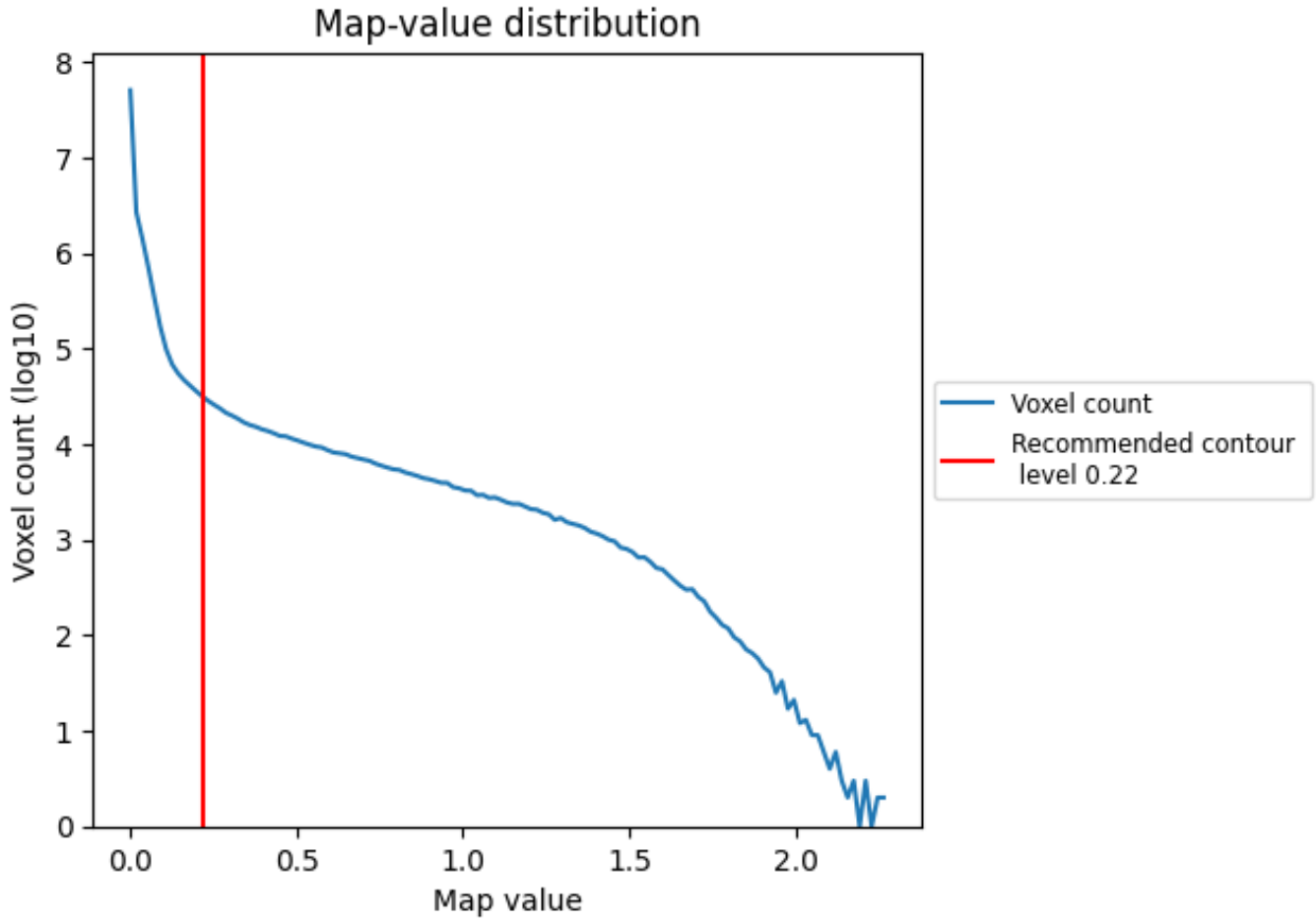
6.6 Mask visualisation (i)

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

7 Map analysis

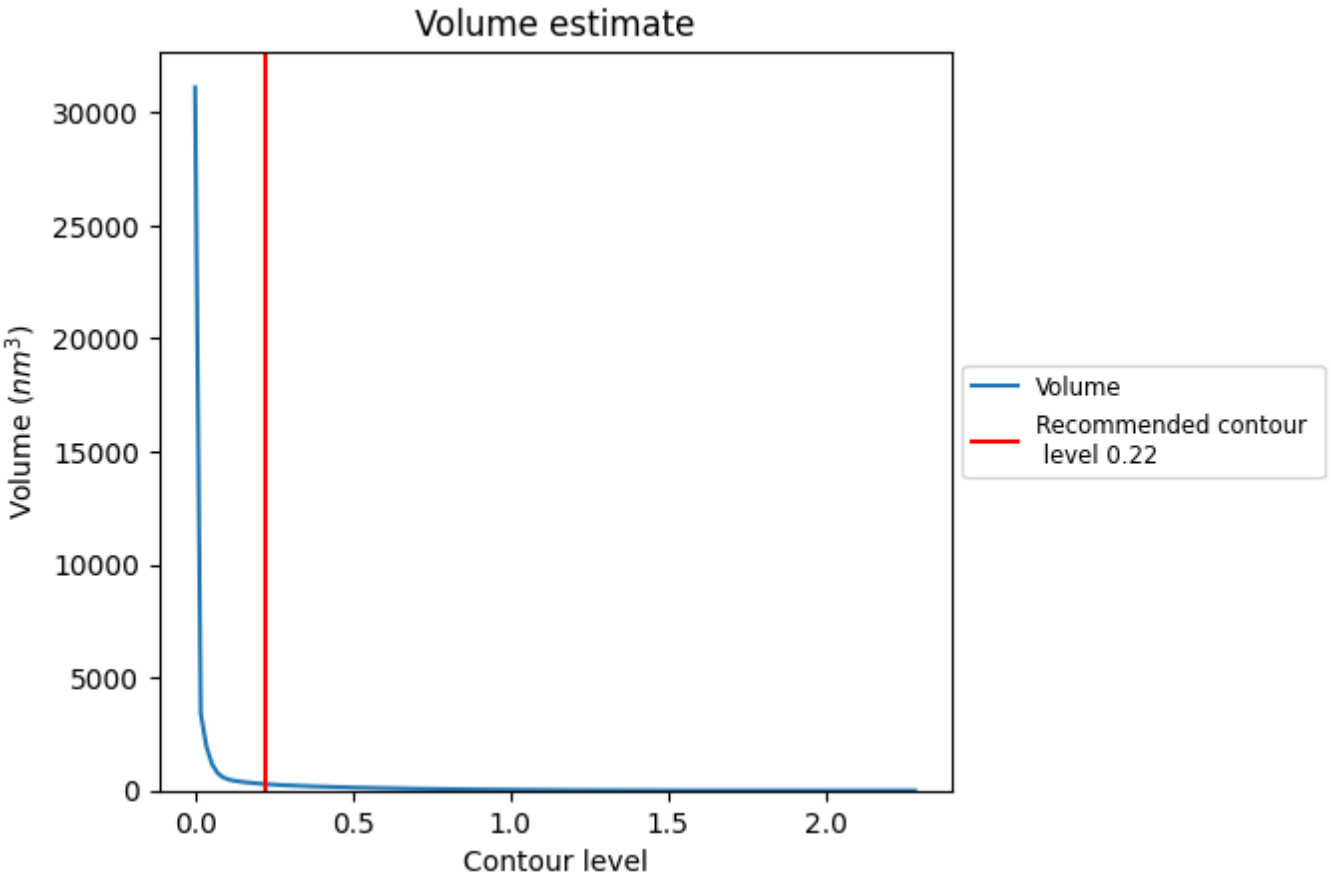
This section contains the results of statistical analysis of the map.

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.

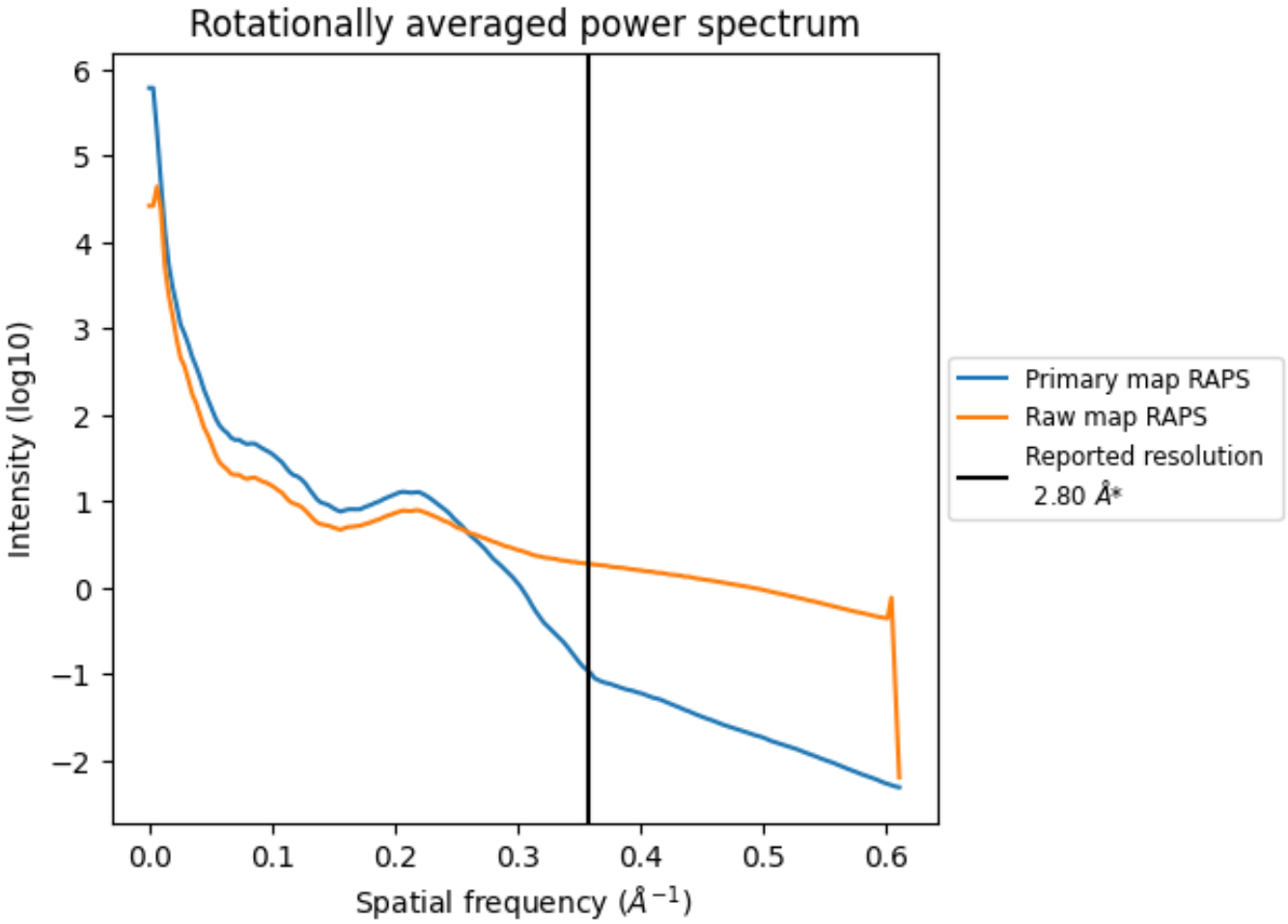
7.2 Volume estimate (i)



The volume at the recommended contour level is 290 nm^3; this corresponds to an approximate mass of 262 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)

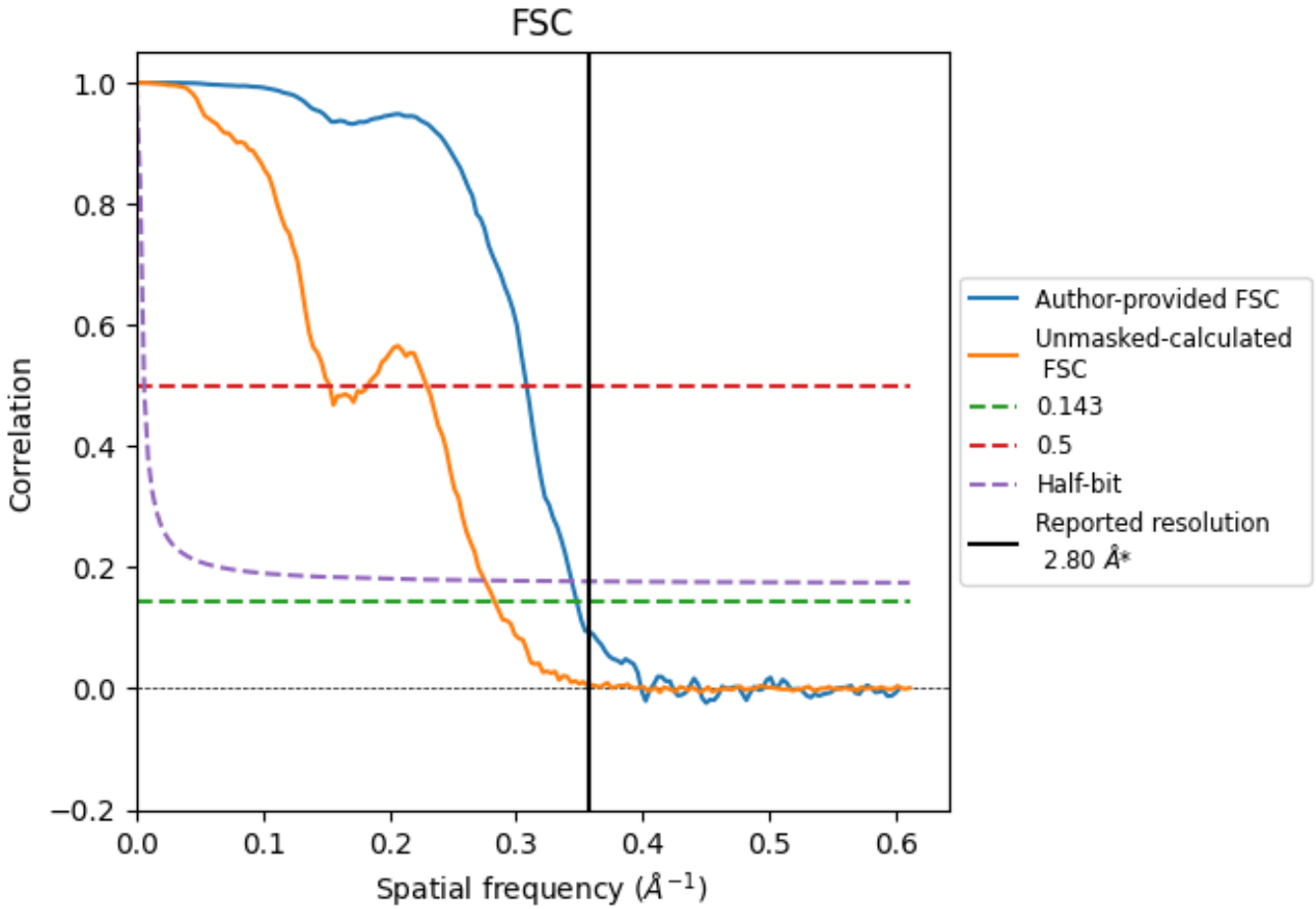


*Reported resolution corresponds to spatial frequency of 0.357 Å^{-1}

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 0.357 Å^{-1}

8.2 Resolution estimates (i)

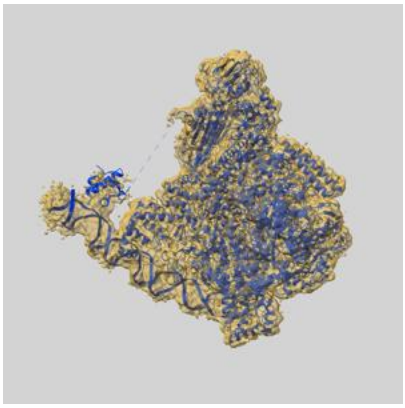
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.80	-	-
Author-provided FSC curve	2.87	3.24	2.91
Unmasked-calculated*	3.53	6.55	3.64

*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 3.53 differs from the reported value 2.8 by more than 10 %

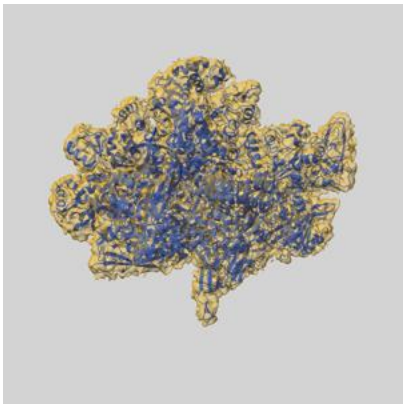
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-59288 and PDB model 32ZR. Per-residue inclusion information can be found in section 3 on page 7.

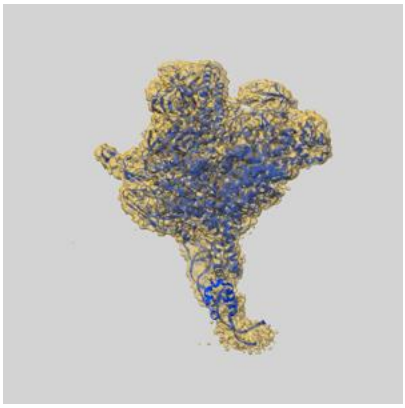
9.1 Map-model overlay



X



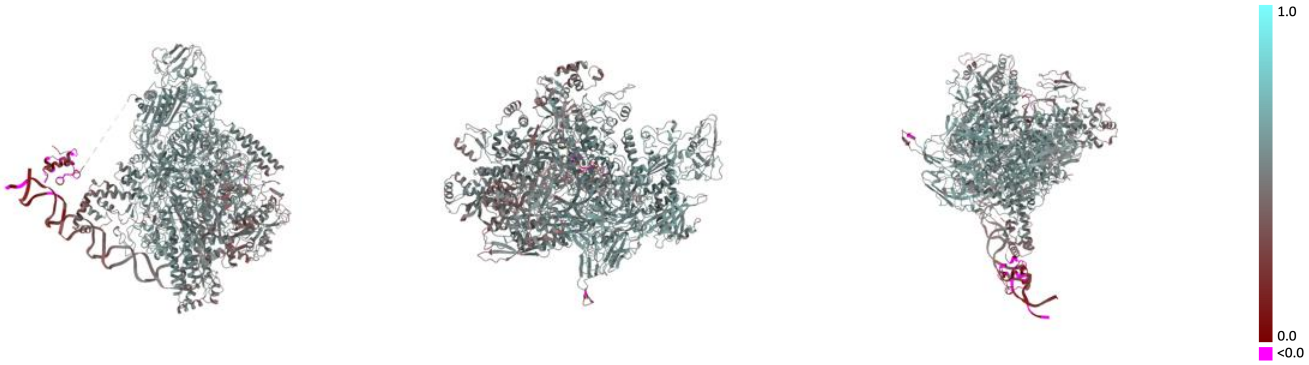
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.22 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.

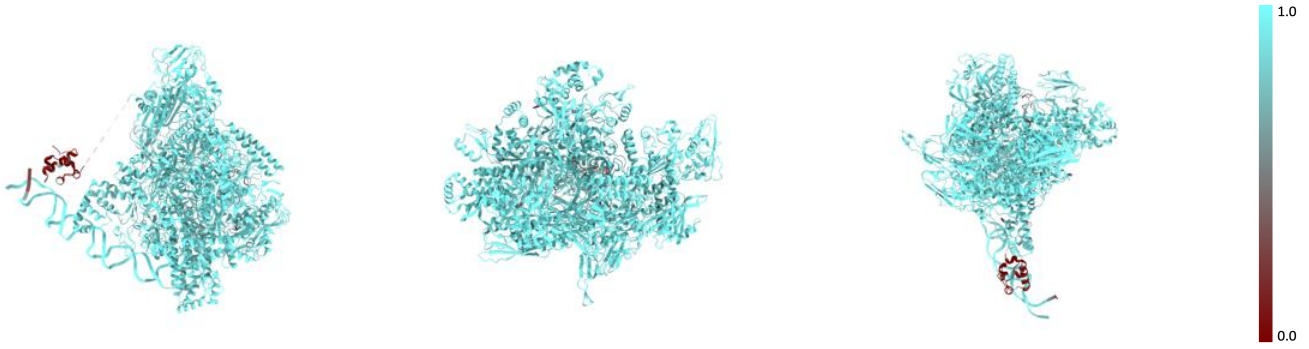
9.2 Q-score mapped to coordinate model (i)



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.

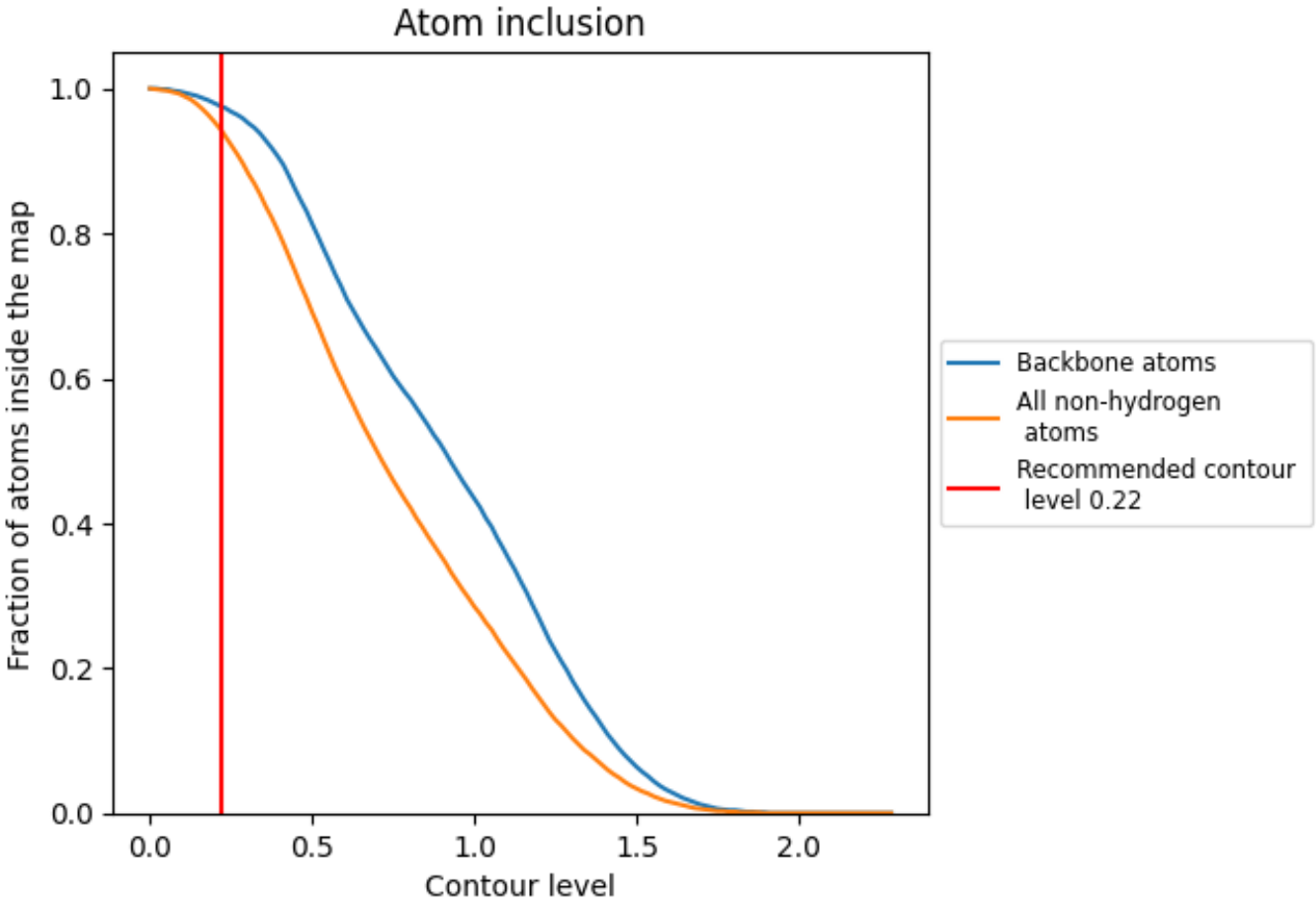
9.3 Atom inclusion mapped to coordinate model (i)



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 (0.22).

9.4 Atom inclusion (i)



At the recommended contour level, 98% of all backbone atoms, 94% 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 (0.22) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9420	0.5090
D	0.9490	0.4410
E	0.9670	0.5580
G	0.9350	0.4820
K	0.9410	0.3790
L	0.9120	0.3580
R	0.9330	0.5370
U	0.7920	0.4500
V	0.9590	0.5250
X	0.9630	0.5330
Y	0.9580	0.5410

