



## Full wwPDB EM Validation Report ⓘ

Oct 15, 2024 – 12:42 AM JST

PDB ID : 8HQZ  
EMDB ID : EMD-34955  
Title : Baseplate of DT57C bacteriophage in the full state  
Authors : Ayala, R.; Moiseenko, A.V.; Chen, T.H.; Kulikov, E.E.; Golomidova, A.K.;  
Orekhov, P.S.; Street, M.A.; Sokolova, O.S.; Letarov, A.V.; Wolf, M.  
Deposited on : 2022-12-14  
Resolution : 3.80 Å(reported)

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.dev113  
MolProbity : 4.02b-467  
Percentile statistics : 20231227.v01 (using entries in the PDB archive December 27th 2023)  
MapQ : 1.9.13  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.39

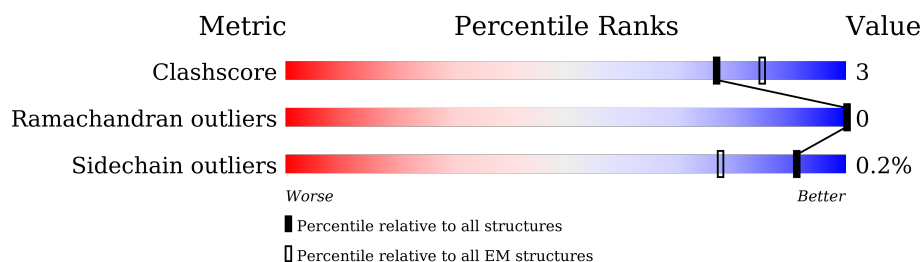
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.80 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.



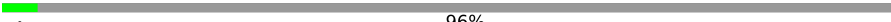
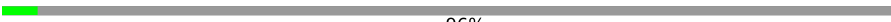



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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	949	 91% 7% .
2	H	204	 95% 5%
2	I	204	 94% 5%
3	L	300	 94% 5% .
4	Q	140	 91% 9% .
4	R	140	 91% 9% .
4	S	140	 94% 5% .
4	T	140	 91% 6% .

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Mol	Chain	Length	Quality of chain
5	d	1076	 96%
5	e	1076	 96%
5	f	1076	 96%
6	k	1227	 97%
7	p	468	 100%

## 2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 22127 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 Baseplate hub protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	930	Total	C	N	O	S	0	0
			7438	4707	1259	1462	10		

- Molecule 2 is a protein called Distal tail protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	H	204	Total	C	N	O	S	0	0
			1600	1023	266	308	3		
2	I	204	Total	C	N	O	S	0	0
			1600	1023	266	308	3		

- Molecule 3 is a protein called Minor tail protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	L	296	Total	C	N	O	S	0	0
			2400	1533	395	464	8		

- Molecule 4 is a protein called L-shaped tail fiber assembly.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	Q	139	Total	C	N	O	S	0	0
			1069	669	175	221	4		
4	R	139	Total	C	N	O	S	0	0
			1069	669	175	221	4		
4	S	139	Total	C	N	O	S	0	0
			1069	669	175	221	4		
4	T	137	Total	C	N	O	S	0	0
			1056	662	173	217	4		

- Molecule 5 is a protein called L-shaped tail fiber assembly.

Mol	Chain	Residues	Atoms				AltConf	Trace
5	d	46	Total	C	N	O	0	0
			338	213	53	72		
5	e	47	Total	C	N	O	0	0
			343	216	54	73		
5	f	47	Total	C	N	O	0	0
			336	209	53	74		

- Molecule 6 is a protein called Tape measure protein.

Mol	Chain	Residues	Atoms				AltConf	Trace
6	k	36	Total	C	N	O	0	0
			257	155	47	55		

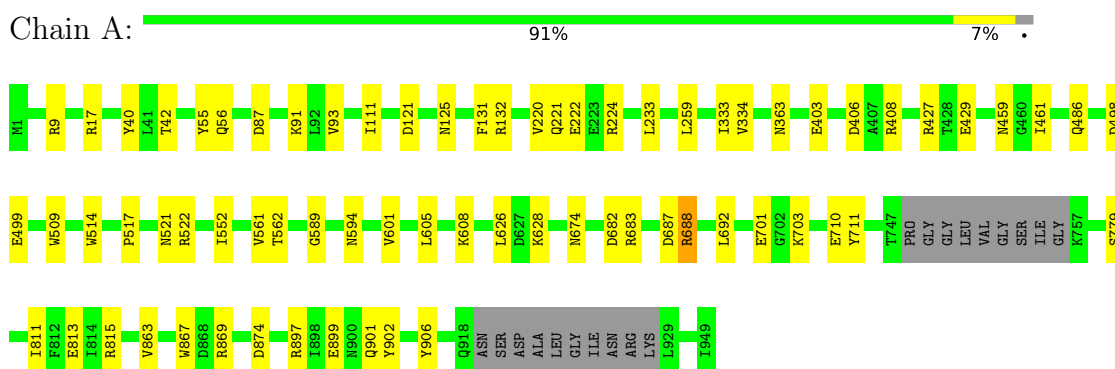
- Molecule 7 is a protein called Tail tube protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
7	p	466	Total	C	N	O	S	0	0
			3552	2222	577	744	9		

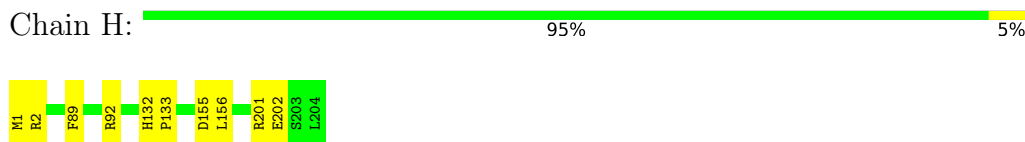
### 3 Residue-property plots

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

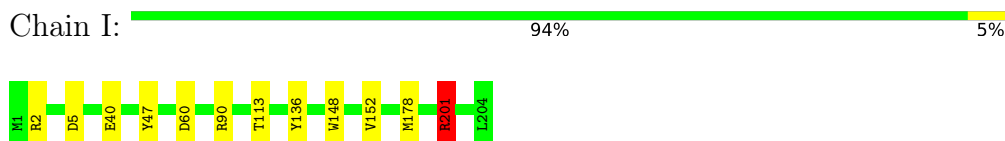
- Molecule 1: Baseplate hub protein



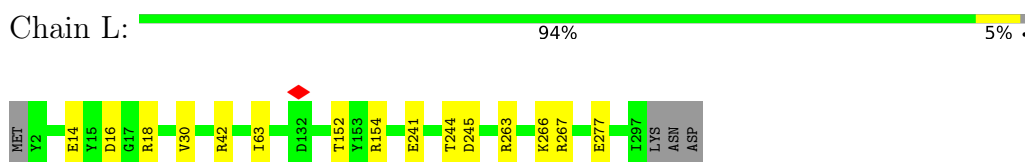
- Molecule 2: Distal tail protein



- Molecule 2: Distal tail protein



- Molecule 3: Minor tail protein



- Molecule 4: L-shaped tail fiber assembly

MET	S2	K15	T49 I50	V82	T86	R96	I111	L112	M113	T114	M128	D136	T139	Q140
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- Chain R:  91% 9%

Amino Acid	Percentage (%)
MET	~10
S2	~10
T3	~10
K15	~10
V16	~10
P17	~10
D26	~10
V27	~10
V82	~10
V85	~10
G89	~10
E129	~10
I134	~10
G137	~10
Q140	~10

- Chain S:  94% 5%

Diagram illustrating the structure of a protein, showing residues MET, S2, V16, P17, Q23, Q73, R102, L112, S120, E121, V122, and Q140. A red diamond marker is positioned above the S120/E121/V122 region.

- Chain T:  91% 6% .

MET  
 SER  
 THR  
 E4  
 E14  
 K15  
 M25  
 T42  
 I111  
 R115  
 G119  
 S120  
 E129  
 G137  
 Q140

- Chain d:  96%

ALA	VAL	ALA	ASP	GLU	ALA	MET
PHE	LEU	GLU	ARG	THR	GLY	ALA
ARG	PRO	THR	ALA	ALA	LEU	ALA
ALA	ASP	PRO	THR	THR	SER	LYS
THR	GLY	VAL	ILE	GLY	GLY	THR
LYS	ARG	GLU	VAL	VAL	ALA	ARG
GLY	GLU	LEU	ASP	ASP	SER	ALA
GLU	VAL	GLU	SER	SER	ALA	ALA
GLU	LEU	GLU	LYS	ALA	ASN	GLY
ASP	ARG	THR	LEU	ALA	SER	SER
ALA	VAL	GLU	ASP	THR	ALA	ALA
GLY	ASP	GLN	LYS	THR	ALA	ALA
VAL	TYR	LYS	VAL	ALA	SER	ALA
ILE	PRO	LEU	ASP	GLY	ALA	ALA
LYS	ASP	THR	ILE	GLU	GLN	ALA
ASP	THR	SER	SER	GLY	GLY	GLY
GLN	TRP	VAL	GLY	ALA	SER	SER
ILE	GLU	ASN	PHE	ALA	GLU	GLU
PRO	ALA	ASN	ILE	ALA	GLY	ILE
TYR	ILE	VAL	VAL	ALA	PHE	ALA
VAL	VAL	ARG	VAL	ARG	ALA	ALA
VAL	ALA	ASP	TYR	GLN	ASP	LYS
THR	GLY	ASP	LYS	SER	SER	GLU
VAL	LEU	ASP	THR	GLU	ALA	SER
ASN	ILE	GLY	LYS	LEU	LEU	LEU
GLY	PRO	GLY	ALA	ASN	GLU	ASN
ILE	SER	ASN	GLU	ALA	ALA	ASN
SER	VAL	VAL	GLU	LYS	ALA	ALA
PRO	SER	GLN	ASP	ILE	GLN	LYS
ASP	GLU	ILE	ALA	SER	ASN	ASP
ALA	ALA	THR	ASP	GLU	ALA	SER
ILE	GLU	PRO	VAL	THR	GLU	GLU
THR	TRP	PRO	VAL	ASN	GLN	ASN
GLY	GLN	GLY	ASN	ALA	SER	GLU
ASN	ALA	GLY	ARG	ALA	ARG	ALA
VAL	GLY	ASN	VAL	GLY	LEU	ALA
GLU	ALA	PRO	LEU	SER	LEU	ILE
ILE	SER	SER	ASP	ALA	ALA	SER
ASP	LEU	LEU	GLU	THR	GLU	ALA
THR	TYR	TRP	LYS	GLU	GLN	GLY
SER	PHE	LEU	VAL	ALA	ALA	ALA
LEU	SER	GLY	LEU	GLY	LYS	SER
GLN	THR	GLU	VAL	ASP	THR	GLU
GLY	GLY	VAL	TRP	LYS	ALA	GLU
THR	ASP	THR	ASN	ALA	ALA	SER
VAL	GLY	TRP	GLN	ILE	GLN	ALA
SER	PHE	THR	THR	ASP	GLN	SER
ILE	THR	PRO	ASN	ALA	GLN	ALA
THR	THR	TYR	SER	THR	GLN	ALA
GLN	PHE	ASP	LYS	THR	THR	SER
GLY	ARG	LYS	TYR	GLY	ALA	ALA
GLY	LEU	ASP	GLY	ASP	ALA	SER
THR	PRO	SER	TRP	THR	GLU	ALA
GLY	ASP	GLY	TYR	ARG	GLU	ALA
ALA	MET	VAL	LYS	ALA	ALA	GLU
THR	MET	GLY	VAL	LYS	LYS	SER
THR	GLN	TYR	VAL	ALA	THR	GLU
LYS	GLY	PRO	GLY	ALA	GLY	ARG
THR	GLN	GLN	THR	ALA	ALA	CYS









## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	42697	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 ( $e^-/\text{\AA}^2$ )	67	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	4000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	29.178	Depositor
Minimum map value	-11.233	Depositor
Average map value	0.001	Depositor
Map value standard deviation	1.012	Depositor
Recommended contour level	4	Depositor
Map size (Å)	560.0, 560.0, 560.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.4, 1.4, 1.4	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	0.49	0/7597	0.58	0/10315
2	H	0.51	0/1638	0.56	0/2229
2	I	0.53	0/1638	0.56	0/2229
3	L	0.51	0/2455	0.58	0/3334
4	Q	0.45	0/1084	0.61	0/1469
4	R	0.47	0/1084	0.58	0/1469
4	S	0.48	0/1084	0.59	0/1469
4	T	0.44	0/1071	0.60	0/1451
5	d	0.45	0/340	0.56	0/464
5	e	0.38	0/345	0.52	0/471
5	f	0.41	0/338	0.61	0/462
6	k	0.39	0/258	0.58	0/346
7	p	0.44	0/3613	0.57	0/4931
All	All	0.48	0/22545	0.58	0/30639

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
2	I	0	2
3	L	0	1
All	All	0	4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (4) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	522	ARG	Sidechain
2	I	201	ARG	Sidechain
2	I	90	ARG	Sidechain
3	L	42	ARG	Sidechain

## 5.2 Too-close contacts

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	7438	0	7221	46	0
2	H	1600	0	1579	5	0
2	I	1600	0	1579	8	0
3	L	2400	0	2324	8	0
4	Q	1069	0	1054	6	0
4	R	1069	0	1054	8	0
4	S	1069	0	1054	5	0
4	T	1056	0	1042	5	0
5	d	338	0	351	0	0
5	e	343	0	356	0	0
5	f	336	0	340	0	0
6	k	257	0	247	0	0
7	p	3552	0	3459	0	0
All	All	22127	0	21660	88	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 3.

All (88) 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:A:589:GLY:O	1:A:628:LYS:NZ	2.31	0.64
1:A:682:ASP:OD1	1:A:683:ARG:N	2.35	0.60
1:A:687:ASP:OD1	1:A:688:ARG:NE	2.26	0.60
4:T:14:GLU:OE1	4:T:137:GLY:N	2.26	0.60
1:A:562:THR:O	1:A:674:ASN:ND2	2.31	0.59
1:A:486:GLN:NE2	1:A:561:VAL:O	2.31	0.58
3:L:263:ARG:HH11	3:L:263:ARG:HG2	1.70	0.57

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:L:267:ARG:NH2	3:L:277:GLU:OE1	2.36	0.56
1:A:403:GLU:OE1	1:A:408:ARG:NH2	2.29	0.56
1:A:220:VAL:HG12	1:A:222:GLU:H	1.70	0.56
1:A:869:ARG:NH1	1:A:874:ASP:OD1	2.38	0.55
1:A:9:ARG:HH11	1:A:9:ARG:HG2	1.72	0.54
3:L:241:GLU:O	3:L:266:LYS:NZ	2.41	0.53
1:A:93:VAL:HG12	1:A:93:VAL:O	2.10	0.52
1:A:701:GLU:OE1	1:A:703:LYS:N	2.37	0.52
4:Q:15:LYS:HE3	4:Q:139:THR:O	2.09	0.52
1:A:40:TYR:HB3	1:A:55:TYR:CD1	2.46	0.51
2:H:2:ARG:HG2	2:H:2:ARG:HH11	1.75	0.51
1:A:867:TRP:CE2	1:A:902:TYR:HA	2.46	0.51
2:I:2:ARG:NH2	2:I:5:ASP:OD2	2.40	0.51
1:A:17:ARG:NE	2:I:40:GLU:OE2	2.38	0.51
1:A:459:ASN:OD1	1:A:461:ILE:N	2.41	0.51
1:A:514:TRP:CD2	1:A:521:ASN:HB3	2.46	0.50
1:A:514:TRP:CG	1:A:521:ASN:HB3	2.46	0.50
1:A:498:ASP:OD1	1:A:499:GLU:N	2.45	0.50
4:R:15:LYS:HE2	4:R:137:GLY:O	2.12	0.50
1:A:552:ILE:HG22	1:A:561:VAL:HG22	1.94	0.49
4:Q:113:MET:SD	4:Q:114:THR:N	2.86	0.49
1:A:333:ILE:HG23	1:A:334:VAL:HG23	1.93	0.49
1:A:710:GLU:O	1:A:711:TYR:CD1	2.66	0.48
4:T:25:MET:N	4:T:25:MET:SD	2.86	0.48
1:A:333:ILE:HG23	1:A:334:VAL:N	2.28	0.48
1:A:131:PHE:CG	1:A:132:ARG:N	2.81	0.48
1:A:514:TRP:CD2	1:A:521:ASN:CB	2.96	0.48
2:I:136:TYR:HB3	2:I:152:VAL:CG2	2.44	0.48
1:A:42:THR:HG22	1:A:56:GLN:O	2.13	0.47
2:H:201:ARG:NH1	2:H:202:GLU:O	2.41	0.47
4:T:42:THR:O	4:T:115:ARG:NH1	2.43	0.47
1:A:233:LEU:HD12	1:A:601:VAL:HG23	1.96	0.46
3:L:14:GLU:OE1	3:L:154:ARG:NE	2.48	0.46
4:Q:82:VAL:O	4:Q:86:THR:HG23	2.15	0.46
4:Q:96:ARG:NH2	4:Q:136:ASP:OD2	2.47	0.46
1:A:509:TRP:CH2	1:A:517:PRO:HD3	2.49	0.46
1:A:813:GLU:OE2	1:A:815:ARG:NE	2.40	0.46
1:A:692:LEU:HD12	1:A:711:TYR:CE2	2.51	0.46
4:R:82:VAL:HA	4:R:85:VAL:HG12	1.98	0.45
1:A:863:VAL:HG22	1:A:906:TYR:HB3	1.98	0.45
1:A:897:ARG:NE	1:A:899:GLU:OE2	2.46	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:S:122:VAL:O	4:S:122:VAL:HG13	2.15	0.45
2:I:113:THR:HG21	2:I:148:TRP:CE2	2.52	0.45
4:R:89:GLY:HA2	4:R:134:ILE:HD13	1.99	0.45
1:A:605:LEU:O	1:A:608:LYS:HG2	2.18	0.44
4:R:16:VAL:HG13	4:R:17:PRO:HD2	1.99	0.44
4:T:15:LYS:HE2	4:T:140:GLN:HG3	1.99	0.44
1:A:259:LEU:HD12	1:A:259:LEU:N	2.33	0.44
4:R:26:ASP:OD1	4:R:27:VAL:N	2.51	0.44
1:A:869:ARG:NH2	1:A:874:ASP:OD1	2.49	0.44
1:A:221:GLN:OE1	1:A:224:ARG:HD3	2.18	0.43
1:A:111:ILE:O	1:A:111:ILE:HG23	2.17	0.43
1:A:220:VAL:HG23	1:A:429:GLU:OE1	2.18	0.43
1:A:897:ARG:NH2	1:A:899:GLU:OE2	2.49	0.43
3:L:244:THR:HG22	3:L:245:ASP:N	2.34	0.43
4:Q:111:ILE:HD11	4:Q:128:MET:HB3	2.00	0.43
2:H:89:PHE:HB2	2:H:92:ARG:NH2	2.33	0.43
4:T:111:ILE:O	4:T:111:ILE:HG23	2.18	0.43
4:Q:49:THR:HG22	4:Q:50:ILE:N	2.34	0.43
1:A:901:GLN:CD	1:A:902:TYR:H	2.22	0.42
2:H:155:ASP:OD1	2:H:156:LEU:N	2.50	0.42
2:I:47:TYR:CE2	2:I:201:ARG:NH2	2.87	0.42
1:A:87:ASP:OD2	1:A:91:LYS:NZ	2.41	0.42
1:A:363:ASN:O	1:A:427:ARG:NH2	2.53	0.42
3:L:152:THR:HG23	4:R:3:THR:CG2	2.50	0.42
1:A:779:SER:HB3	1:A:811:ILE:HG12	2.01	0.42
2:I:47:TYR:CD2	2:I:201:ARG:CZ	3.03	0.41
4:S:16:VAL:HG13	4:S:17:PRO:HD2	2.02	0.41
4:S:23:GLN:NE2	4:S:73:GLN:O	2.50	0.41
2:H:132:HIS:CG	2:H:133:PRO:HD2	2.56	0.41
1:A:121:ASP:O	1:A:125:ASN:N	2.54	0.41
1:A:406:ASP:OD1	1:A:406:ASP:C	2.59	0.41
3:L:30:VAL:HG22	3:L:63:ILE:CD1	2.50	0.41
1:A:333:ILE:CG2	1:A:334:VAL:N	2.83	0.41
2:I:178:MET:SD	2:I:201:ARG:HD2	2.61	0.41
4:R:129:GLU:OE1	4:S:102:ARG:NH1	2.54	0.41
4:S:112:LEU:HD12	4:S:112:LEU:O	2.21	0.41
2:I:60:ASP:OD1	2:I:60:ASP:N	2.53	0.40
4:R:16:VAL:CG1	4:R:17:PRO:HD2	2.51	0.40
3:L:16:ASP:OD2	3:L:18:ARG:NH1	2.42	0.40
1:A:594:ASN:ND2	1:A:626:LEU:O	2.50	0.40

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	924/949 (97%)	891 (96%)	33 (4%)	0	100	100
2	H	202/204 (99%)	196 (97%)	6 (3%)	0	100	100
2	I	202/204 (99%)	195 (96%)	7 (4%)	0	100	100
3	L	294/300 (98%)	281 (96%)	13 (4%)	0	100	100
4	Q	137/140 (98%)	131 (96%)	6 (4%)	0	100	100
4	R	137/140 (98%)	134 (98%)	3 (2%)	0	100	100
4	S	137/140 (98%)	132 (96%)	5 (4%)	0	100	100
4	T	135/140 (96%)	132 (98%)	3 (2%)	0	100	100
5	d	44/1076 (4%)	42 (96%)	2 (4%)	0	100	100
5	e	45/1076 (4%)	43 (96%)	2 (4%)	0	100	100
5	f	45/1076 (4%)	39 (87%)	6 (13%)	0	100	100
6	k	34/1227 (3%)	34 (100%)	0	0	100	100
7	p	464/468 (99%)	446 (96%)	18 (4%)	0	100	100
All	All	2800/7140 (39%)	2696 (96%)	104 (4%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all PDB entries followed by that with respect to all EM entries.

The Analysed column shows the number of residues for which the sidechain conformation was analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	819/832 (98%)	818 (100%)	1 (0%)	92	95

*Continued on next page...*

Continued from previous page...

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	H	179/179 (100%)	178 (99%)	1 (1%)	84	88
2	I	179/179 (100%)	178 (99%)	1 (1%)	84	88
3	L	270/274 (98%)	270 (100%)	0	100	100
4	Q	120/121 (99%)	120 (100%)	0	100	100
4	R	120/121 (99%)	120 (100%)	0	100	100
4	S	120/121 (99%)	120 (100%)	0	100	100
4	T	118/121 (98%)	117 (99%)	1 (1%)	79	84
5	d	39/825 (5%)	39 (100%)	0	100	100
5	e	39/825 (5%)	39 (100%)	0	100	100
5	f	38/825 (5%)	38 (100%)	0	100	100
6	k	26/967 (3%)	26 (100%)	0	100	100
7	p	398/400 (100%)	398 (100%)	0	100	100
All	All	2465/5790 (43%)	2461 (100%)	4 (0%)	91	94

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	688	ARG
2	H	1	MET
2	I	201	ARG
4	T	129	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	14	ASN
1	A	446	HIS
1	A	521	ASN
1	A	597	GLN
1	A	631	GLN
1	A	721	GLN
2	I	25	ASN
6	k	152	ASN
7	p	20	HIS
7	p	31	GLN

### 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.

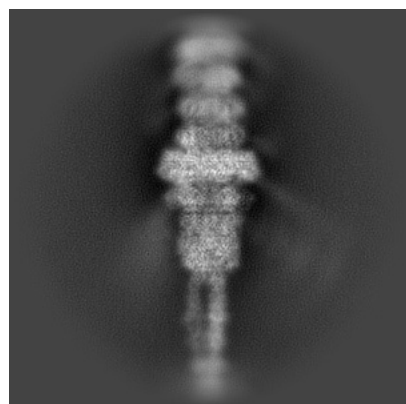
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-34955. These allow visual inspection of the internal detail of the map and identification of artifacts.

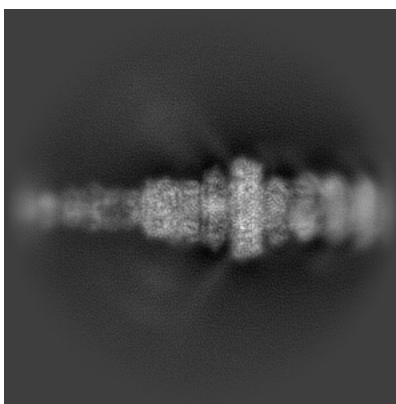
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

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

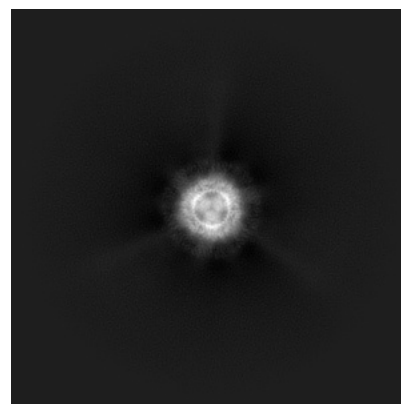
#### 6.1.1 Primary map



X

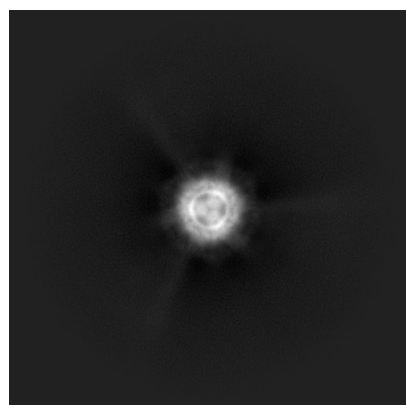


Y

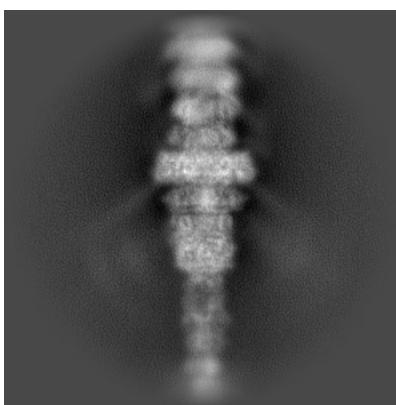


Z

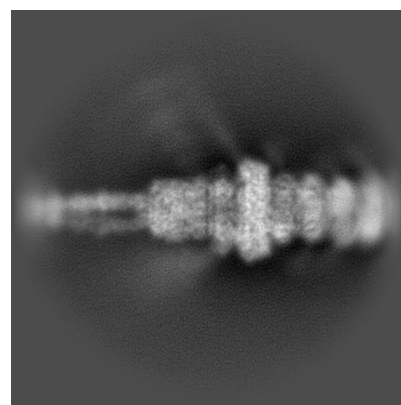
#### 6.1.2 Raw map



X



Y

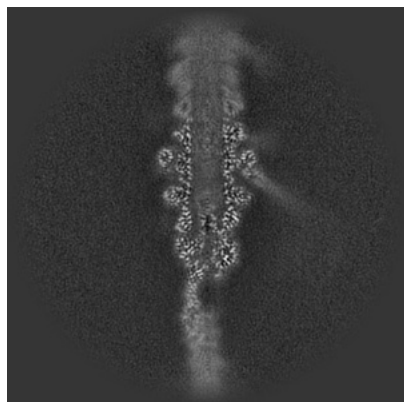


Z

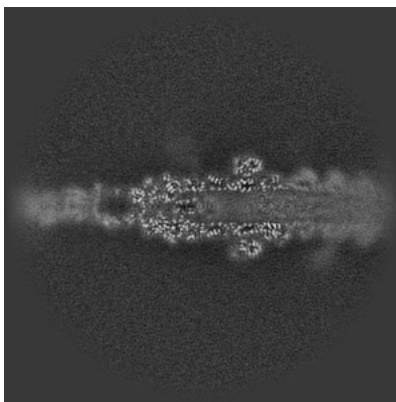
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

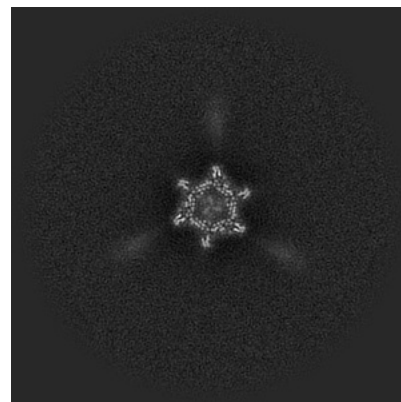
### 6.2.1 Primary map



X Index: 200

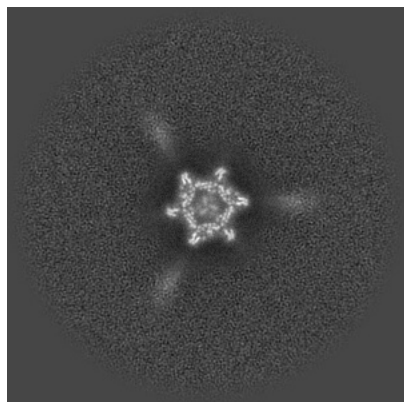


Y Index: 200

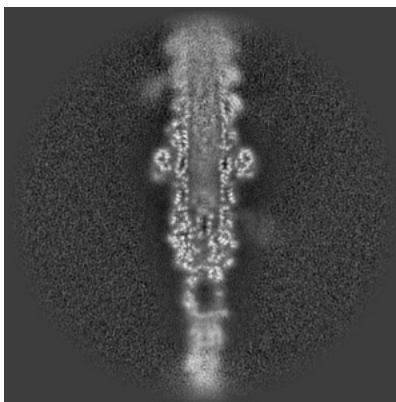


Z Index: 200

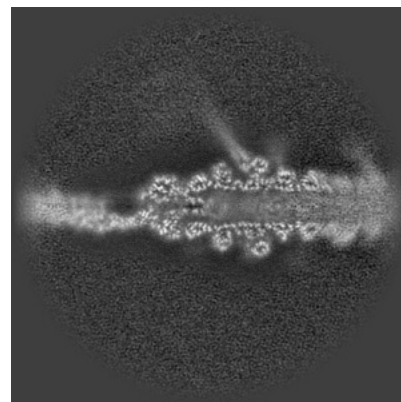
### 6.2.2 Raw map



X Index: 200



Y Index: 200

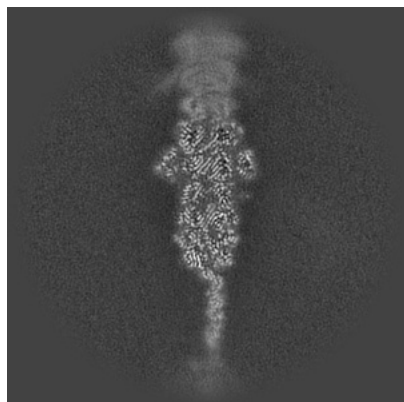


Z Index: 200

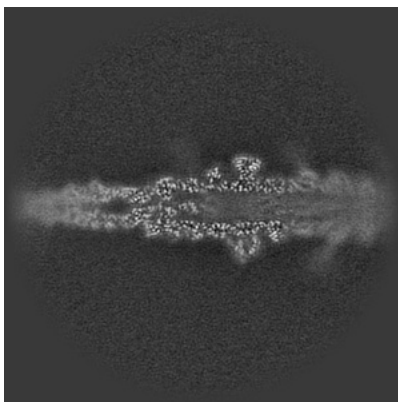
The images above show central slices of the map in three orthogonal directions.

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

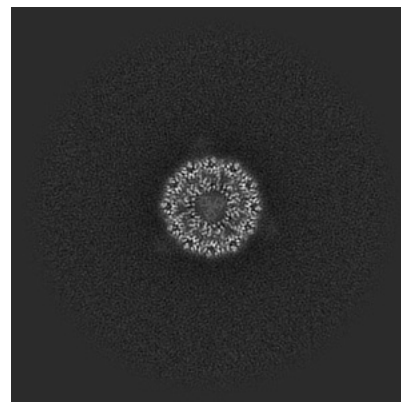
### 6.3.1 Primary map



X Index: 216

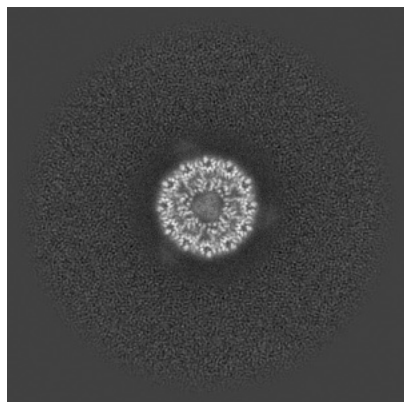


Y Index: 206

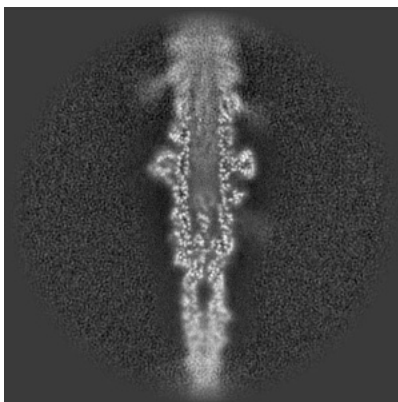


Z Index: 249

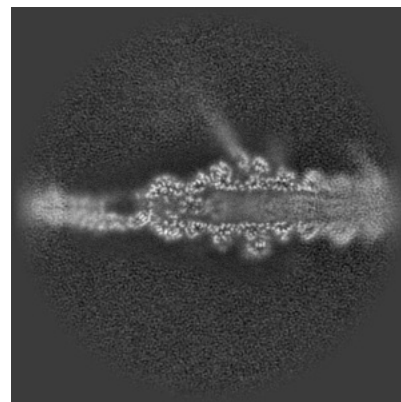
### 6.3.2 Raw map



X Index: 249



Y Index: 206

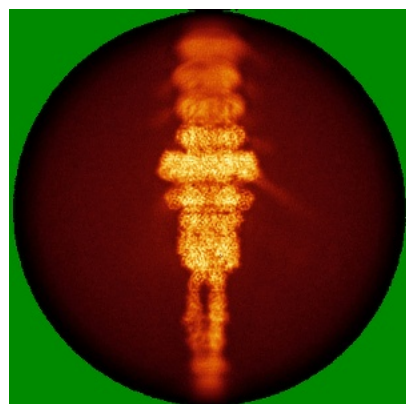


Z Index: 202

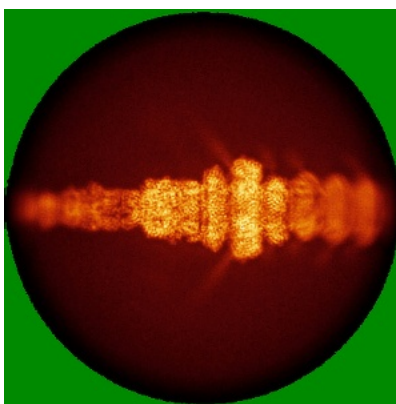
The images above show the largest variance slices of the map in three orthogonal directions.

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

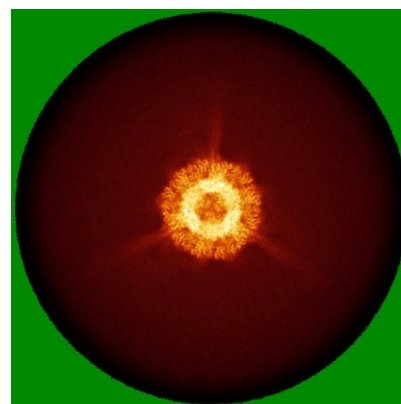
### 6.4.1 Primary map



X

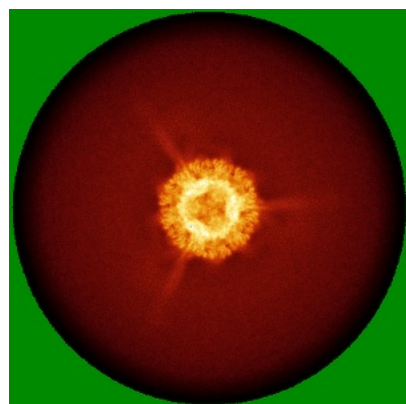


Y

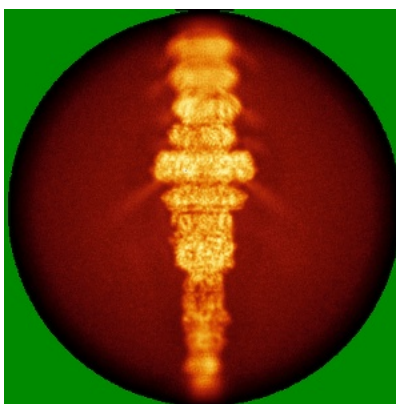


Z

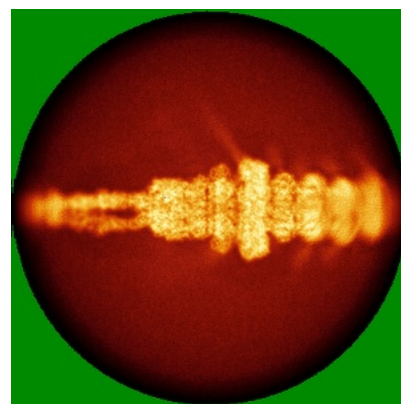
### 6.4.2 Raw map



X



Y

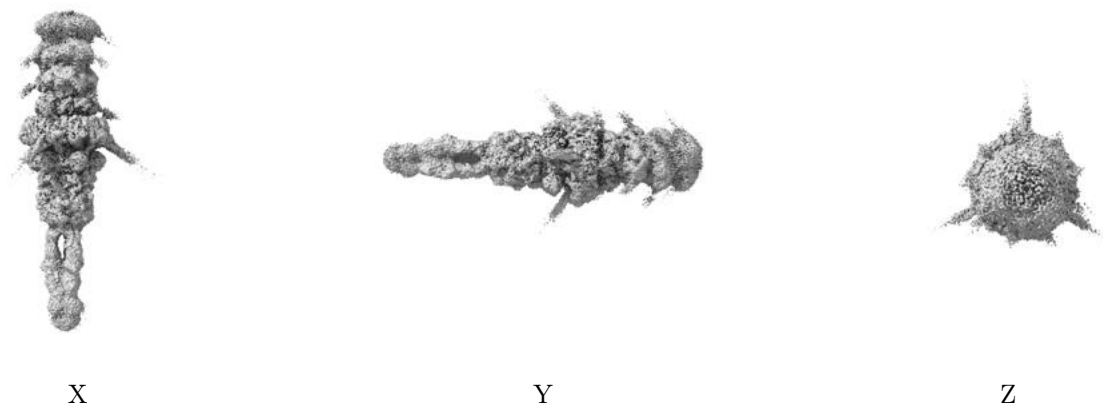


Z

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.

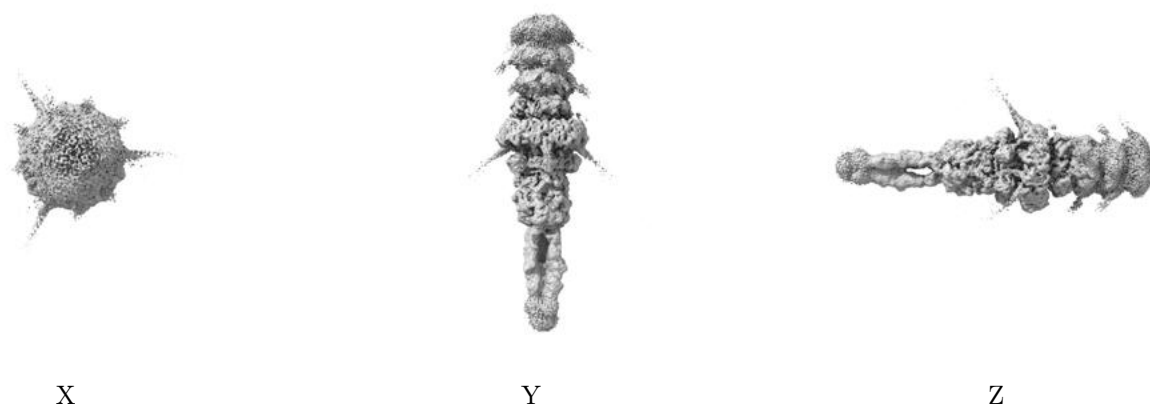
## 6.5 Orthogonal surface views [i](#)

### 6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 4.0. 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



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 shows the 3D surface view of the primary map at 50% transparency overlaid with the specified mask at 0% transparency

A mask typically either:

- Encompasses the whole structure
- Separates out a domain, a functional unit, a monomer or an area of interest from a larger structure

### 6.6.1 emd\_34955\_msk\_1.map [i](#)



X



Y

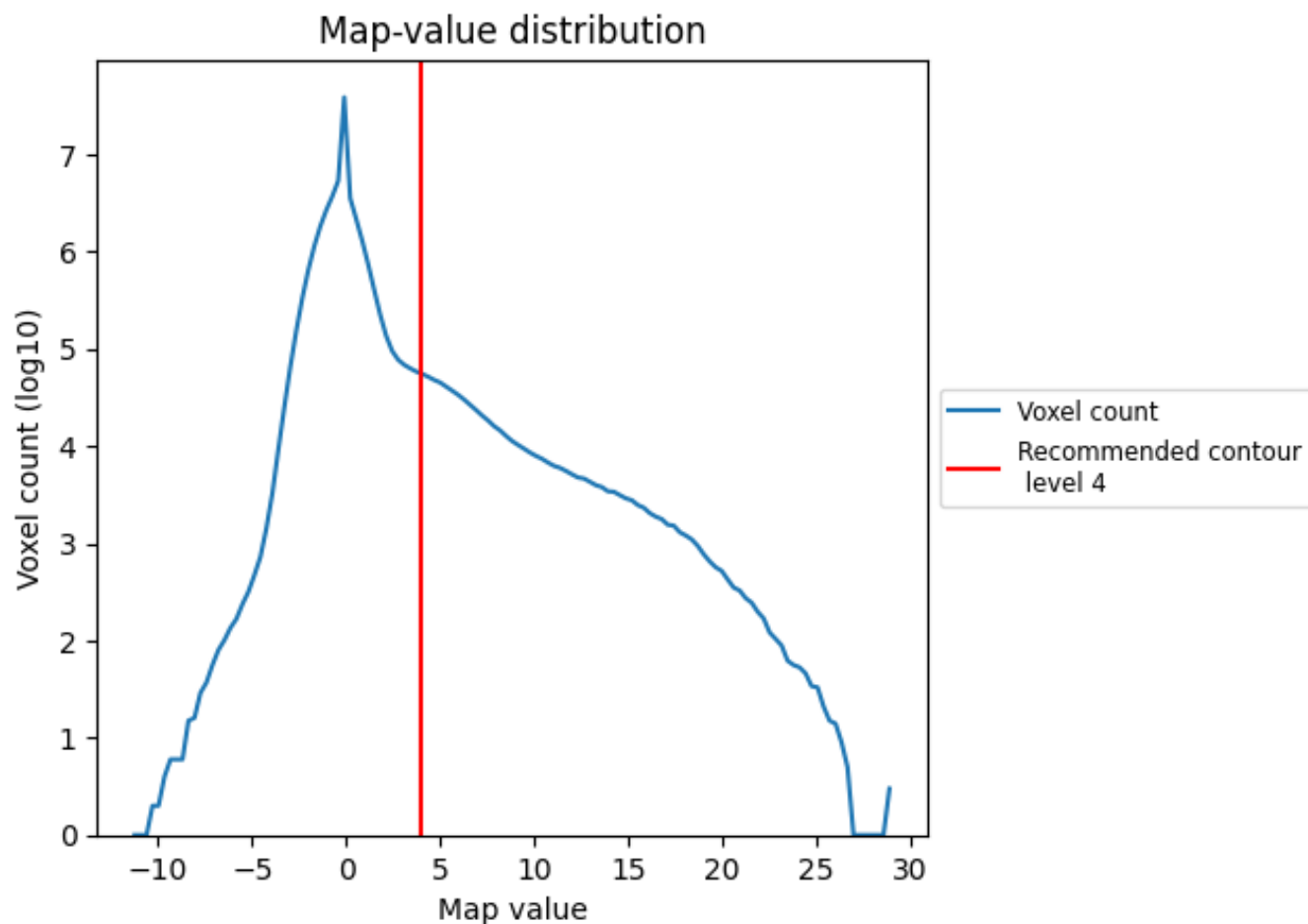


Z

## 7 Map analysis [i](#)

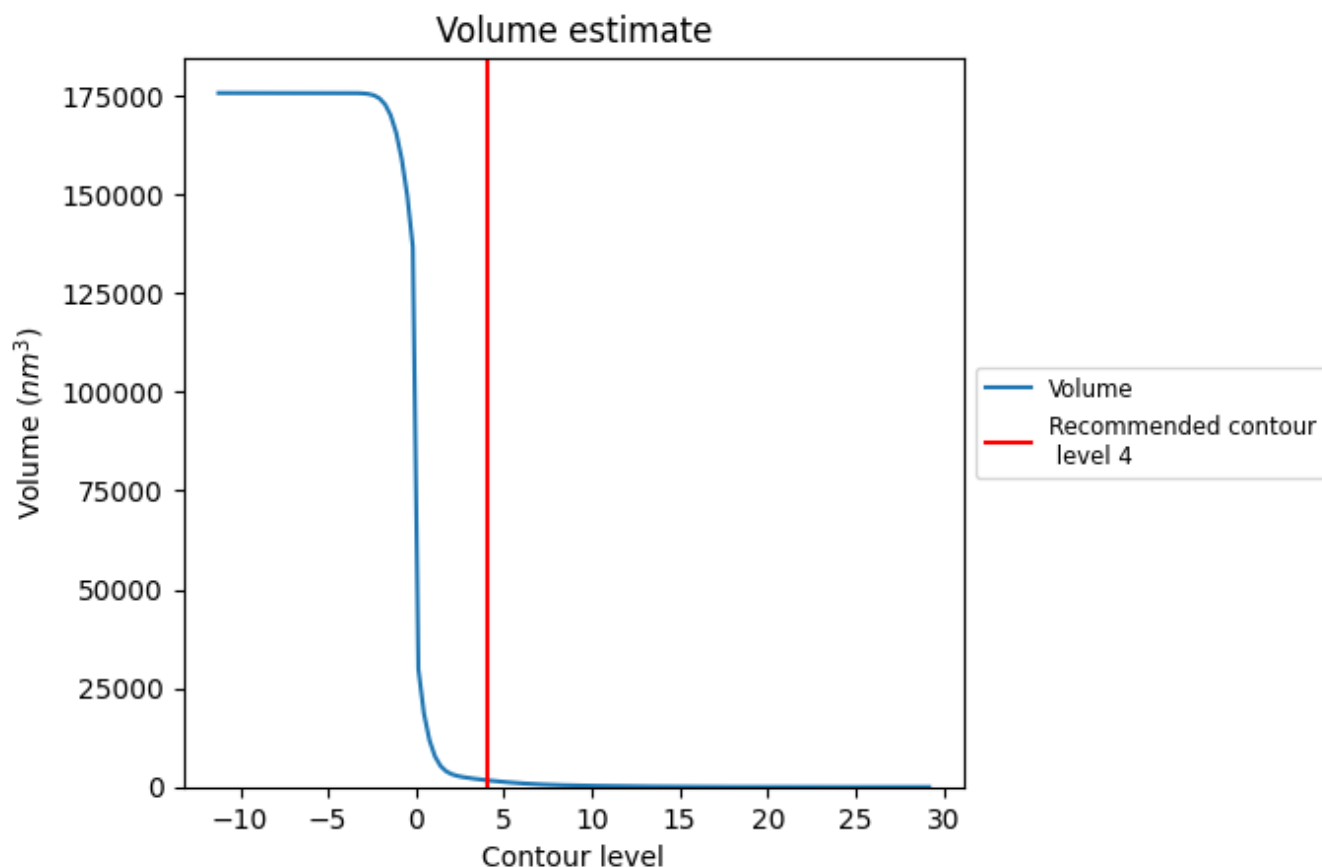
This section contains the results of statistical analysis of the map.

### 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.

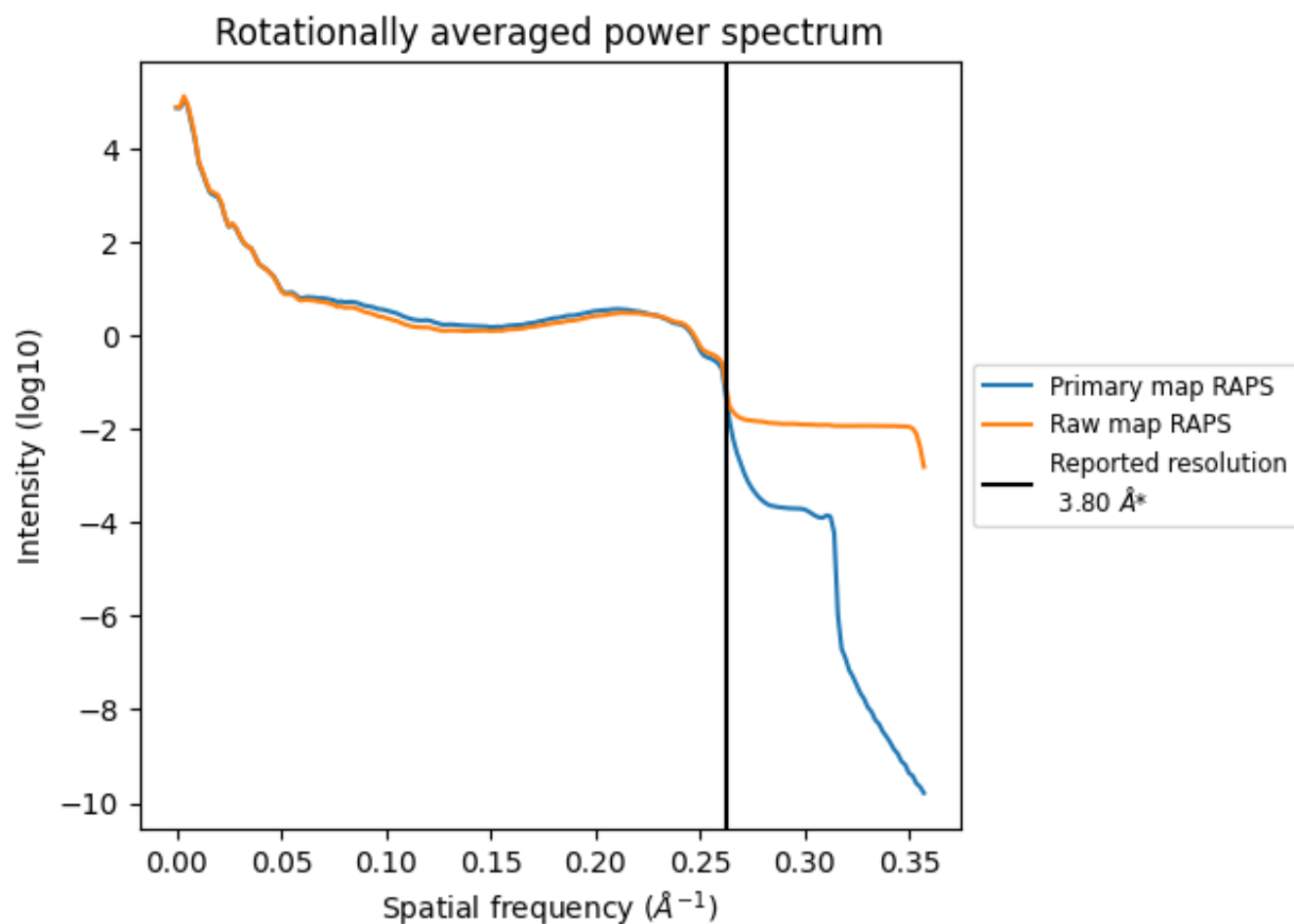
## 7.2 Volume estimate [i](#)



The volume at the recommended contour level is 1719  $\text{nm}^3$ ; this corresponds to an approximate mass of 1552 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 ⓘ

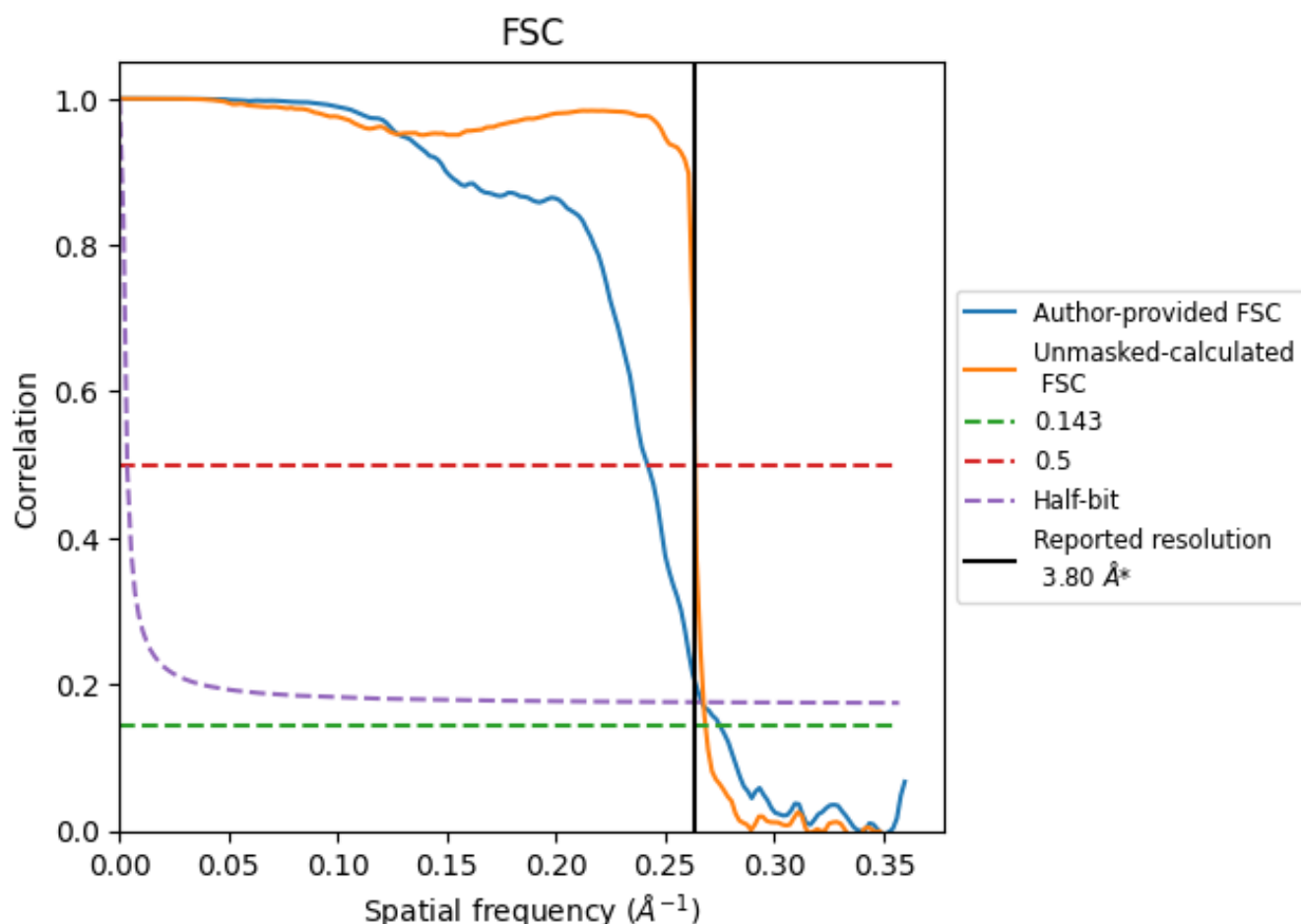


\*Reported resolution corresponds to spatial frequency of 0.263 Å<sup>-1</sup>

## 8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

### 8.1 FSC [i](#)



\*Reported resolution corresponds to spatial frequency of 0.263  $\text{\AA}^{-1}$

## 8.2 Resolution estimates [i](#)

Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.80	-	-
Author-provided FSC curve	3.64	4.14	3.75
Unmasked-calculated*	3.73	3.79	3.74

\*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps.

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

This section contains information regarding the fit between EMDB map EMD-34955 and PDB model 8HQZ. Per-residue inclusion information can be found in section 3 on page 6.

### 9.1 Map-model overlays

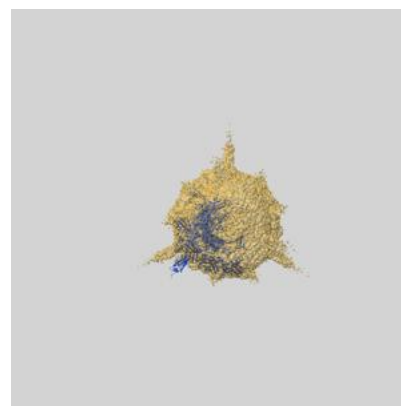
#### 9.1.1 Map-model overlay [i](#)



X

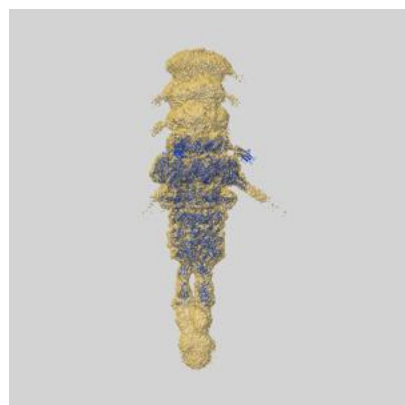


Y



Z

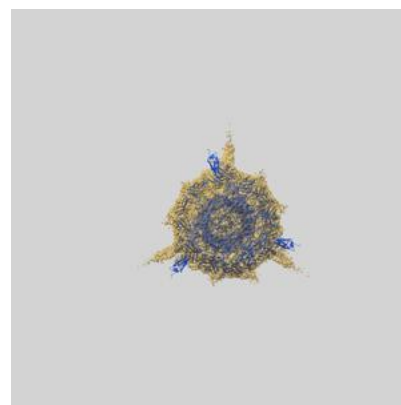
#### 9.1.2 Map-model assembly overlay [i](#)



X



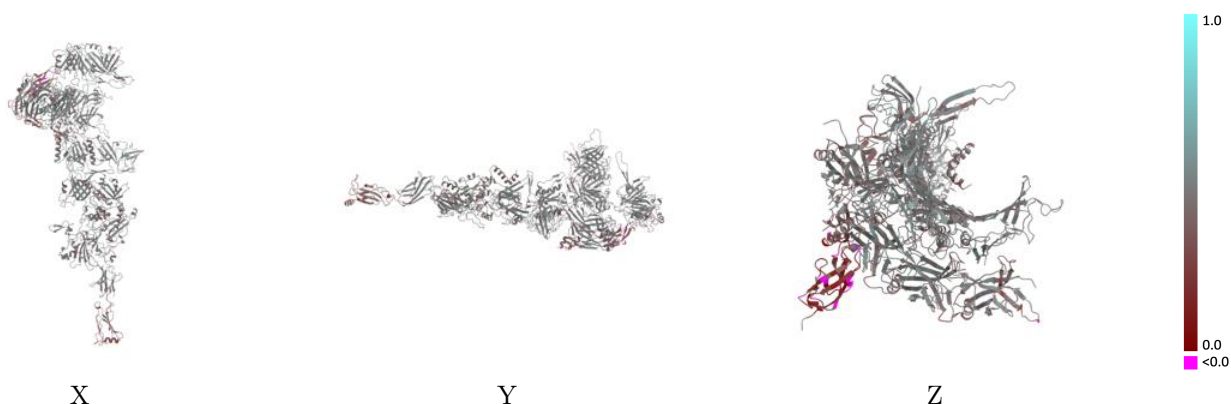
Y



Z

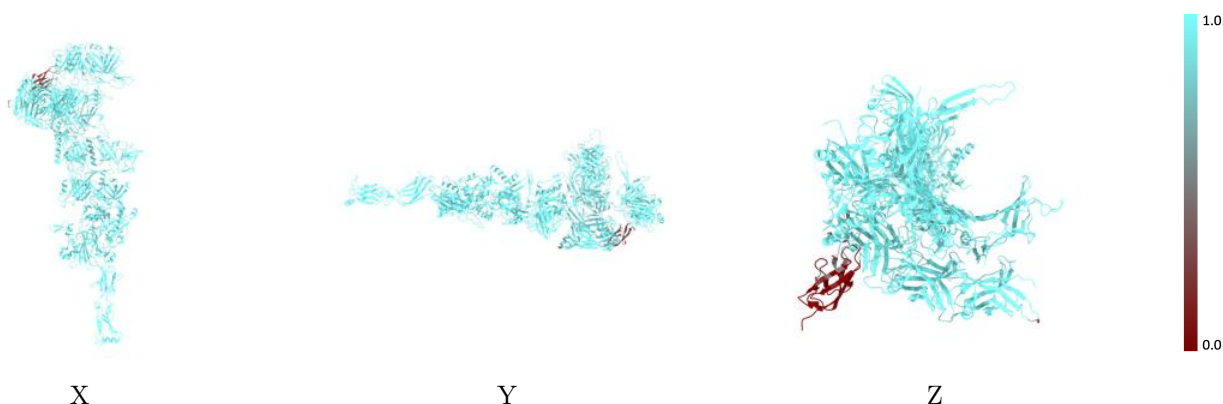
The images above show the 3D surface view of the map at the recommended contour level 4.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



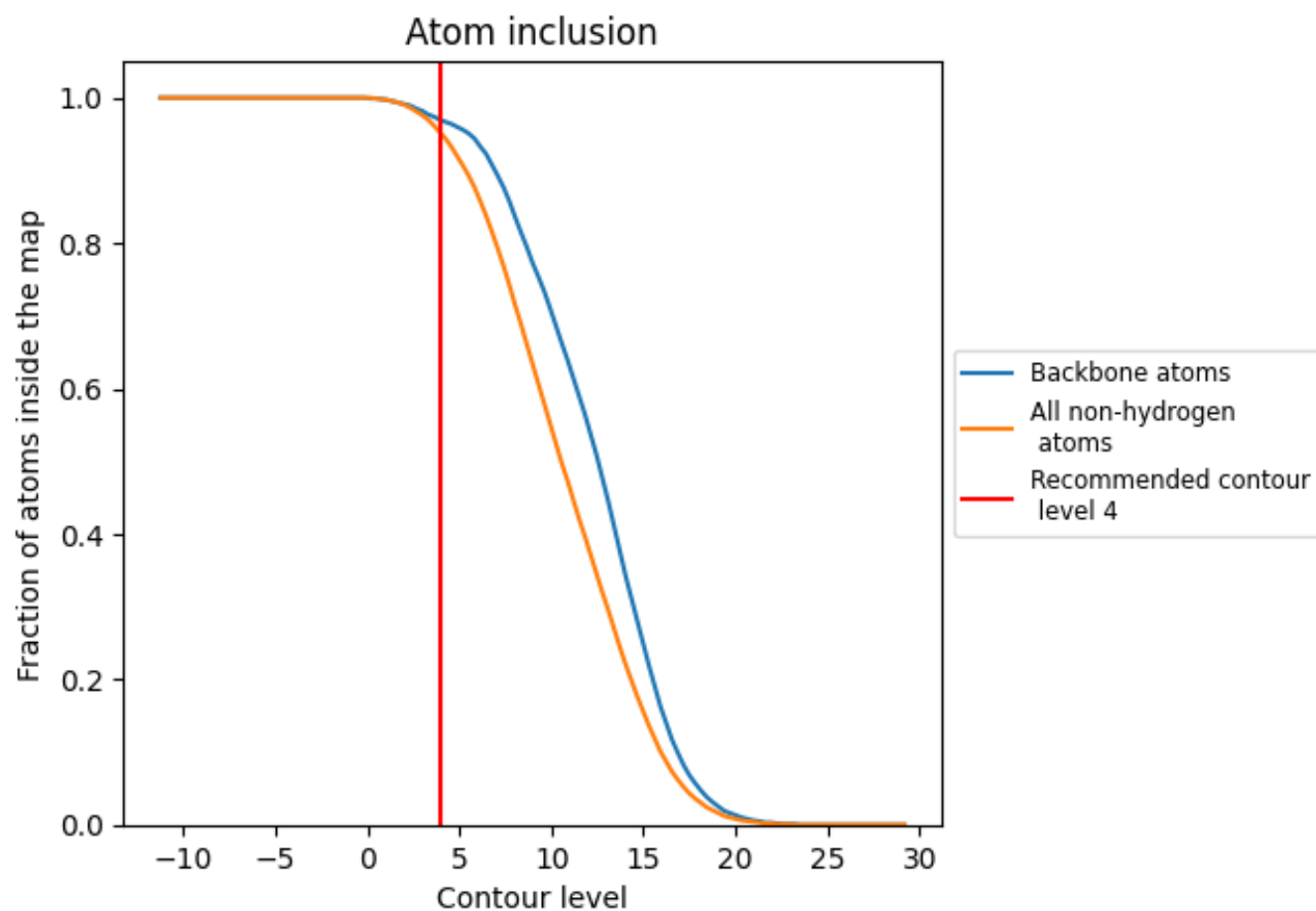
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (4).





























## 9.4 Atom inclusion [i](#)



At the recommended contour level, 97% of all backbone atoms, 95% 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 (4) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.9510	 0.4460
A	 0.9880	 0.4480
H	 0.9770	 0.4710
I	 0.9760	 0.4730
L	 0.9560	 0.4760
Q	 0.9750	 0.4440
R	 0.9780	 0.4550
S	 0.9650	 0.4430
T	 0.9670	 0.4290
d	 0.9730	 0.4360
e	 0.9530	 0.4060
f	 0.9760	 0.4130
k	 0.9490	 0.4010
p	 0.8210	 0.4120

