



Full wwPDB EM Validation Report ⓘ

Sep 22, 2026 – 03:29 PM EDT

PDB ID : 9YG8 / pdb_00009yg8
EMDB ID : EMD-72915
Title : CryoEM structure of F7 pyocin tail fiber in the pre-ejection state
Authors : He, Y.; Cai, X.Y.; Li, A.S.C.; Davidson, A.R.; Zhou, Z.H.
Deposited on : 2025-09-28
Resolution : 3.10 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

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:

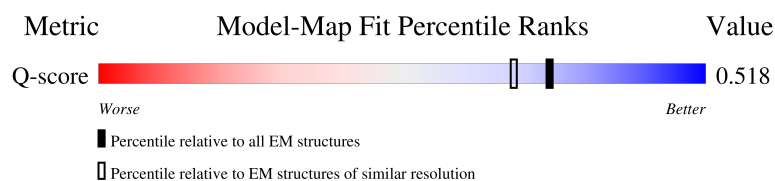
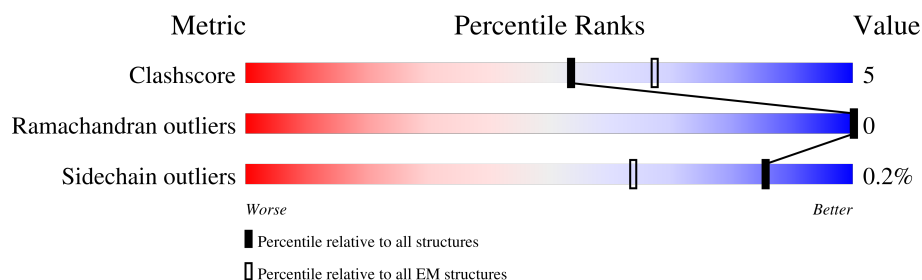
EMDB validation analysis : 0.0.1.dev133
MolProbity : 4-5-2 with Phenix2.0
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.50

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 3.10 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	-
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 $\leq 5\%$. 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	1222	
1	E	1222	
1	I	1222	
2	B	464	

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Mol	Chain	Length	Quality of chain
2	C	464	 86%14%
2	D	464	 87%12%
2	F	464	 88%12%
2	G	464	 85%15%
2	H	464	 91%9%
2	J	464	 87%12%
2	K	464	 84%16%
2	L	464	 91%9%

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 33666 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 Phage tail protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	163	Total	C	N	O	S	0	0
			1178	690	225	262	1		
1	E	163	Total	C	N	O	S	0	0
			1178	690	225	262	1		
1	I	163	Total	C	N	O	S	0	0
			1178	690	225	262	1		

There are 192 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	709	VAL	GLU	conflict	UNP A0A7M2ZXW7
A	863	VAL	ILE	conflict	UNP A0A7M2ZXW7
A	1049	ALA	VAL	conflict	UNP A0A7M2ZXW7
A	1155	THR	ALA	conflict	UNP A0A7M2ZXW7
A	1163	ASP	-	expression tag	UNP A0A7M2ZXW7
A	1164	GLU	-	expression tag	UNP A0A7M2ZXW7
A	1165	ILE	-	expression tag	UNP A0A7M2ZXW7
A	1166	ALA	-	expression tag	UNP A0A7M2ZXW7
A	1167	SER	-	expression tag	UNP A0A7M2ZXW7
A	1168	ASN	-	expression tag	UNP A0A7M2ZXW7
A	1169	ASN	-	expression tag	UNP A0A7M2ZXW7
A	1170	TYR	-	expression tag	UNP A0A7M2ZXW7
A	1171	LYS	-	expression tag	UNP A0A7M2ZXW7
A	1172	PRO	-	expression tag	UNP A0A7M2ZXW7
A	1173	ASP	-	expression tag	UNP A0A7M2ZXW7
A	1174	ARG	-	expression tag	UNP A0A7M2ZXW7
A	1175	THR	-	expression tag	UNP A0A7M2ZXW7
A	1176	GLY	-	expression tag	UNP A0A7M2ZXW7
A	1177	TRP	-	expression tag	UNP A0A7M2ZXW7
A	1178	ARG	-	expression tag	UNP A0A7M2ZXW7
A	1179	LEU	-	expression tag	UNP A0A7M2ZXW7
A	1180	ASP	-	expression tag	UNP A0A7M2ZXW7
A	1181	LYS	-	expression tag	UNP A0A7M2ZXW7
A	1182	ASN	-	expression tag	UNP A0A7M2ZXW7

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Chain	Residue	Modelled	Actual	Comment	Reference
A	1183	GLY	-	expression tag	UNP A0A7M2ZXW7
A	1184	ARG	-	expression tag	UNP A0A7M2ZXW7
A	1185	PHE	-	expression tag	UNP A0A7M2ZXW7
A	1186	GLU	-	expression tag	UNP A0A7M2ZXW7
A	1187	ILE	-	expression tag	UNP A0A7M2ZXW7
A	1188	ASN	-	expression tag	UNP A0A7M2ZXW7
A	1189	GLY	-	expression tag	UNP A0A7M2ZXW7
A	1190	VAL	-	expression tag	UNP A0A7M2ZXW7
A	1191	VAL	-	expression tag	UNP A0A7M2ZXW7
A	1192	GLU	-	expression tag	UNP A0A7M2ZXW7
A	1193	GLY	-	expression tag	UNP A0A7M2ZXW7
A	1194	GLN	-	expression tag	UNP A0A7M2ZXW7
A	1195	GLY	-	expression tag	UNP A0A7M2ZXW7
A	1196	ARG	-	expression tag	UNP A0A7M2ZXW7
A	1197	VAL	-	expression tag	UNP A0A7M2ZXW7
A	1198	THR	-	expression tag	UNP A0A7M2ZXW7
A	1199	ILE	-	expression tag	UNP A0A7M2ZXW7
A	1200	THR	-	expression tag	UNP A0A7M2ZXW7
A	1201	ASN	-	expression tag	UNP A0A7M2ZXW7
A	1202	ASN	-	expression tag	UNP A0A7M2ZXW7
A	1203	ALA	-	expression tag	UNP A0A7M2ZXW7
A	1204	ILE	-	expression tag	UNP A0A7M2ZXW7
A	1205	LYS	-	expression tag	UNP A0A7M2ZXW7
A	1206	VAL	-	expression tag	UNP A0A7M2ZXW7
A	1207	PHE	-	expression tag	UNP A0A7M2ZXW7
A	1208	ASP	-	expression tag	UNP A0A7M2ZXW7
A	1209	ASP	-	expression tag	UNP A0A7M2ZXW7
A	1210	LYS	-	expression tag	UNP A0A7M2ZXW7
A	1211	GLY	-	expression tag	UNP A0A7M2ZXW7
A	1212	ARG	-	expression tag	UNP A0A7M2ZXW7
A	1213	LEU	-	expression tag	UNP A0A7M2ZXW7
A	1214	ARG	-	expression tag	UNP A0A7M2ZXW7
A	1215	VAL	-	expression tag	UNP A0A7M2ZXW7
A	1216	HIS	-	expression tag	UNP A0A7M2ZXW7
A	1217	ILE	-	expression tag	UNP A0A7M2ZXW7
A	1218	GLY	-	expression tag	UNP A0A7M2ZXW7
A	1219	ASP	-	expression tag	UNP A0A7M2ZXW7
A	1220	LEU	-	expression tag	UNP A0A7M2ZXW7
A	1221	SER	-	expression tag	UNP A0A7M2ZXW7
A	1222	ALA	-	expression tag	UNP A0A7M2ZXW7
E	709	VAL	GLU	conflict	UNP A0A7M2ZXW7
E	863	VAL	ILE	conflict	UNP A0A7M2ZXW7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	1049	ALA	VAL	conflict	UNP A0A7M2ZXW7
E	1155	THR	ALA	conflict	UNP A0A7M2ZXW7
E	1163	ASP	-	expression tag	UNP A0A7M2ZXW7
E	1164	GLU	-	expression tag	UNP A0A7M2ZXW7
E	1165	ILE	-	expression tag	UNP A0A7M2ZXW7
E	1166	ALA	-	expression tag	UNP A0A7M2ZXW7
E	1167	SER	-	expression tag	UNP A0A7M2ZXW7
E	1168	ASN	-	expression tag	UNP A0A7M2ZXW7
E	1169	ASN	-	expression tag	UNP A0A7M2ZXW7
E	1170	TYR	-	expression tag	UNP A0A7M2ZXW7
E	1171	LYS	-	expression tag	UNP A0A7M2ZXW7
E	1172	PRO	-	expression tag	UNP A0A7M2ZXW7
E	1173	ASP	-	expression tag	UNP A0A7M2ZXW7
E	1174	ARG	-	expression tag	UNP A0A7M2ZXW7
E	1175	THR	-	expression tag	UNP A0A7M2ZXW7
E	1176	GLY	-	expression tag	UNP A0A7M2ZXW7
E	1177	TRP	-	expression tag	UNP A0A7M2ZXW7
E	1178	ARG	-	expression tag	UNP A0A7M2ZXW7
E	1179	LEU	-	expression tag	UNP A0A7M2ZXW7
E	1180	ASP	-	expression tag	UNP A0A7M2ZXW7
E	1181	LYS	-	expression tag	UNP A0A7M2ZXW7
E	1182	ASN	-	expression tag	UNP A0A7M2ZXW7
E	1183	GLY	-	expression tag	UNP A0A7M2ZXW7
E	1184	ARG	-	expression tag	UNP A0A7M2ZXW7
E	1185	PHE	-	expression tag	UNP A0A7M2ZXW7
E	1186	GLU	-	expression tag	UNP A0A7M2ZXW7
E	1187	ILE	-	expression tag	UNP A0A7M2ZXW7
E	1188	ASN	-	expression tag	UNP A0A7M2ZXW7
E	1189	GLY	-	expression tag	UNP A0A7M2ZXW7
E	1190	VAL	-	expression tag	UNP A0A7M2ZXW7
E	1191	VAL	-	expression tag	UNP A0A7M2ZXW7
E	1192	GLU	-	expression tag	UNP A0A7M2ZXW7
E	1193	GLY	-	expression tag	UNP A0A7M2ZXW7
E	1194	GLN	-	expression tag	UNP A0A7M2ZXW7
E	1195	GLY	-	expression tag	UNP A0A7M2ZXW7
E	1196	ARG	-	expression tag	UNP A0A7M2ZXW7
E	1197	VAL	-	expression tag	UNP A0A7M2ZXW7
E	1198	THR	-	expression tag	UNP A0A7M2ZXW7
E	1199	ILE	-	expression tag	UNP A0A7M2ZXW7
E	1200	THR	-	expression tag	UNP A0A7M2ZXW7
E	1201	ASN	-	expression tag	UNP A0A7M2ZXW7
E	1202	ASN	-	expression tag	UNP A0A7M2ZXW7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	1203	ALA	-	expression tag	UNP A0A7M2ZXW7
E	1204	ILE	-	expression tag	UNP A0A7M2ZXW7
E	1205	LYS	-	expression tag	UNP A0A7M2ZXW7
E	1206	VAL	-	expression tag	UNP A0A7M2ZXW7
E	1207	PHE	-	expression tag	UNP A0A7M2ZXW7
E	1208	ASP	-	expression tag	UNP A0A7M2ZXW7
E	1209	ASP	-	expression tag	UNP A0A7M2ZXW7
E	1210	LYS	-	expression tag	UNP A0A7M2ZXW7
E	1211	GLY	-	expression tag	UNP A0A7M2ZXW7
E	1212	ARG	-	expression tag	UNP A0A7M2ZXW7
E	1213	LEU	-	expression tag	UNP A0A7M2ZXW7
E	1214	ARG	-	expression tag	UNP A0A7M2ZXW7
E	1215	VAL	-	expression tag	UNP A0A7M2ZXW7
E	1216	HIS	-	expression tag	UNP A0A7M2ZXW7
E	1217	ILE	-	expression tag	UNP A0A7M2ZXW7
E	1218	GLY	-	expression tag	UNP A0A7M2ZXW7
E	1219	ASP	-	expression tag	UNP A0A7M2ZXW7
E	1220	LEU	-	expression tag	UNP A0A7M2ZXW7
E	1221	SER	-	expression tag	UNP A0A7M2ZXW7
E	1222	ALA	-	expression tag	UNP A0A7M2ZXW7
I	709	VAL	GLU	conflict	UNP A0A7M2ZXW7
I	863	VAL	ILE	conflict	UNP A0A7M2ZXW7
I	1049	ALA	VAL	conflict	UNP A0A7M2ZXW7
I	1155	THR	ALA	conflict	UNP A0A7M2ZXW7
I	1163	ASP	-	expression tag	UNP A0A7M2ZXW7
I	1164	GLU	-	expression tag	UNP A0A7M2ZXW7
I	1165	ILE	-	expression tag	UNP A0A7M2ZXW7
I	1166	ALA	-	expression tag	UNP A0A7M2ZXW7
I	1167	SER	-	expression tag	UNP A0A7M2ZXW7
I	1168	ASN	-	expression tag	UNP A0A7M2ZXW7
I	1169	ASN	-	expression tag	UNP A0A7M2ZXW7
I	1170	TYR	-	expression tag	UNP A0A7M2ZXW7
I	1171	LYS	-	expression tag	UNP A0A7M2ZXW7
I	1172	PRO	-	expression tag	UNP A0A7M2ZXW7
I	1173	ASP	-	expression tag	UNP A0A7M2ZXW7
I	1174	ARG	-	expression tag	UNP A0A7M2ZXW7
I	1175	THR	-	expression tag	UNP A0A7M2ZXW7
I	1176	GLY	-	expression tag	UNP A0A7M2ZXW7
I	1177	TRP	-	expression tag	UNP A0A7M2ZXW7
I	1178	ARG	-	expression tag	UNP A0A7M2ZXW7
I	1179	LEU	-	expression tag	UNP A0A7M2ZXW7
I	1180	ASP	-	expression tag	UNP A0A7M2ZXW7

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Chain	Residue	Modelled	Actual	Comment	Reference
I	1181	LYS	-	expression tag	UNP A0A7M2ZXW7
I	1182	ASN	-	expression tag	UNP A0A7M2ZXW7
I	1183	GLY	-	expression tag	UNP A0A7M2ZXW7
I	1184	ARG	-	expression tag	UNP A0A7M2ZXW7
I	1185	PHE	-	expression tag	UNP A0A7M2ZXW7
I	1186	GLU	-	expression tag	UNP A0A7M2ZXW7
I	1187	ILE	-	expression tag	UNP A0A7M2ZXW7
I	1188	ASN	-	expression tag	UNP A0A7M2ZXW7
I	1189	GLY	-	expression tag	UNP A0A7M2ZXW7
I	1190	VAL	-	expression tag	UNP A0A7M2ZXW7
I	1191	VAL	-	expression tag	UNP A0A7M2ZXW7
I	1192	GLU	-	expression tag	UNP A0A7M2ZXW7
I	1193	GLY	-	expression tag	UNP A0A7M2ZXW7
I	1194	GLN	-	expression tag	UNP A0A7M2ZXW7
I	1195	GLY	-	expression tag	UNP A0A7M2ZXW7
I	1196	ARG	-	expression tag	UNP A0A7M2ZXW7
I	1197	VAL	-	expression tag	UNP A0A7M2ZXW7
I	1198	THR	-	expression tag	UNP A0A7M2ZXW7
I	1199	ILE	-	expression tag	UNP A0A7M2ZXW7
I	1200	THR	-	expression tag	UNP A0A7M2ZXW7
I	1201	ASN	-	expression tag	UNP A0A7M2ZXW7
I	1202	ASN	-	expression tag	UNP A0A7M2ZXW7
I	1203	ALA	-	expression tag	UNP A0A7M2ZXW7
I	1204	ILE	-	expression tag	UNP A0A7M2ZXW7
I	1205	LYS	-	expression tag	UNP A0A7M2ZXW7
I	1206	VAL	-	expression tag	UNP A0A7M2ZXW7
I	1207	PHE	-	expression tag	UNP A0A7M2ZXW7
I	1208	ASP	-	expression tag	UNP A0A7M2ZXW7
I	1209	ASP	-	expression tag	UNP A0A7M2ZXW7
I	1210	LYS	-	expression tag	UNP A0A7M2ZXW7
I	1211	GLY	-	expression tag	UNP A0A7M2ZXW7
I	1212	ARG	-	expression tag	UNP A0A7M2ZXW7
I	1213	LEU	-	expression tag	UNP A0A7M2ZXW7
I	1214	ARG	-	expression tag	UNP A0A7M2ZXW7
I	1215	VAL	-	expression tag	UNP A0A7M2ZXW7
I	1216	HIS	-	expression tag	UNP A0A7M2ZXW7
I	1217	ILE	-	expression tag	UNP A0A7M2ZXW7
I	1218	GLY	-	expression tag	UNP A0A7M2ZXW7
I	1219	ASP	-	expression tag	UNP A0A7M2ZXW7
I	1220	LEU	-	expression tag	UNP A0A7M2ZXW7
I	1221	SER	-	expression tag	UNP A0A7M2ZXW7
I	1222	ALA	-	expression tag	UNP A0A7M2ZXW7

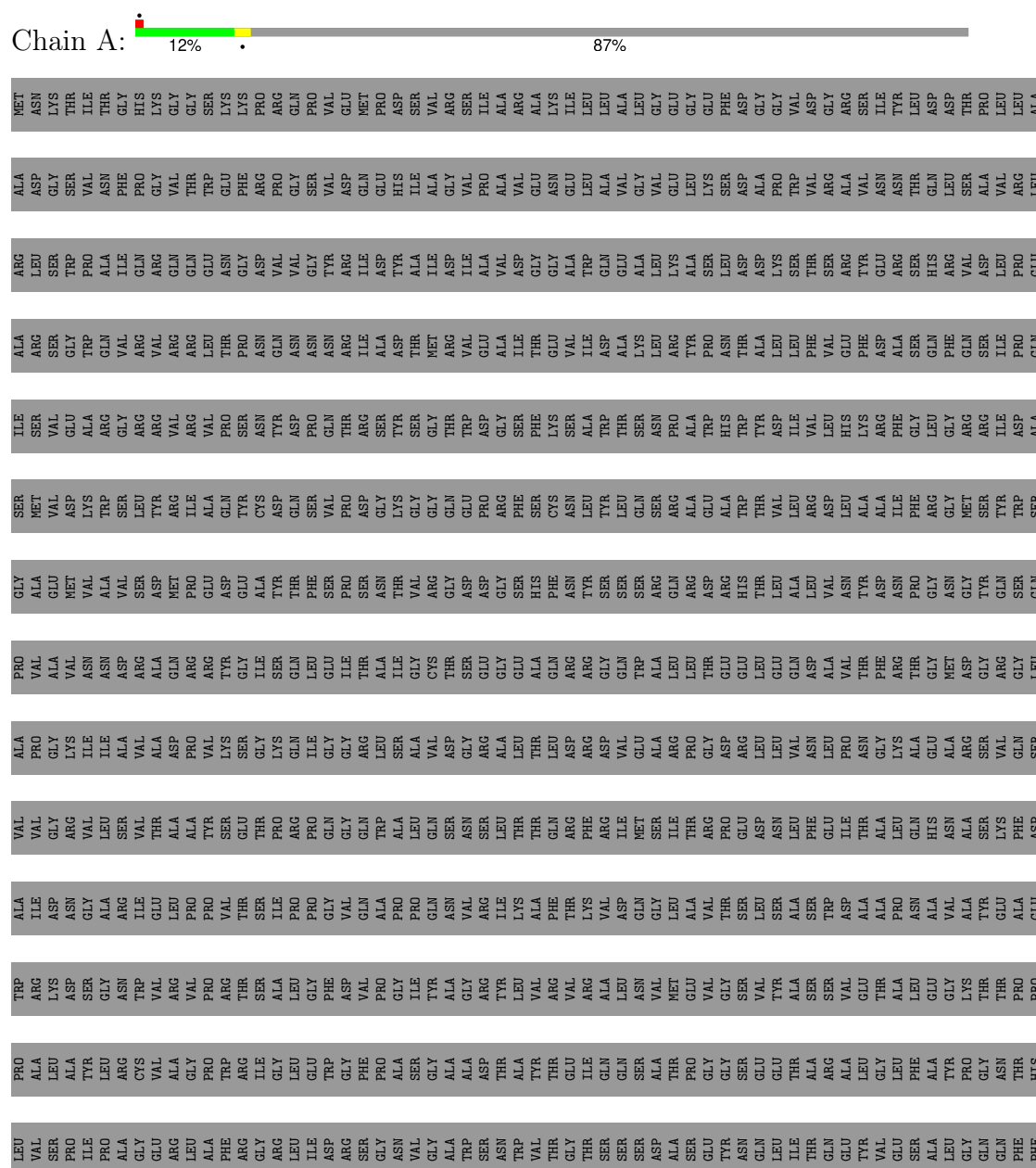
- Molecule 2 is a protein called Side fiber protein of F7 pyocin.

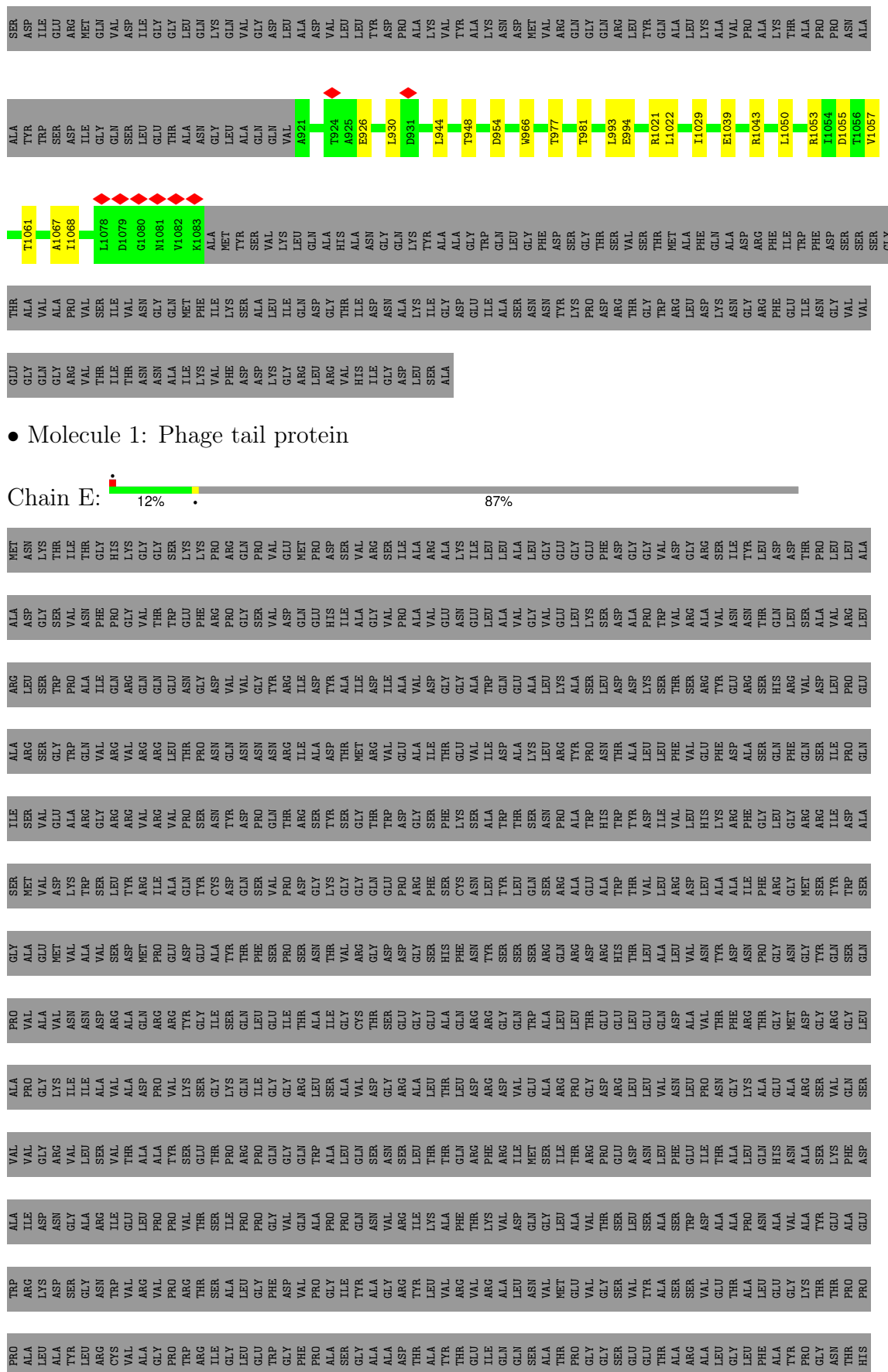
Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	463	Total	C	N	O	S	0	0
			3348	2095	596	640	17		
2	C	463	Total	C	N	O	S	0	0
			3348	2095	596	640	17		
2	D	463	Total	C	N	O	S	0	0
			3348	2095	596	640	17		
2	F	463	Total	C	N	O	S	0	0
			3348	2095	596	640	17		
2	G	463	Total	C	N	O	S	0	0
			3348	2095	596	640	17		
2	H	463	Total	C	N	O	S	0	0
			3348	2095	596	640	17		
2	J	463	Total	C	N	O	S	0	0
			3348	2095	596	640	17		
2	K	463	Total	C	N	O	S	0	0
			3348	2095	596	640	17		
2	L	463	Total	C	N	O	S	0	0
			3348	2095	596	640	17		

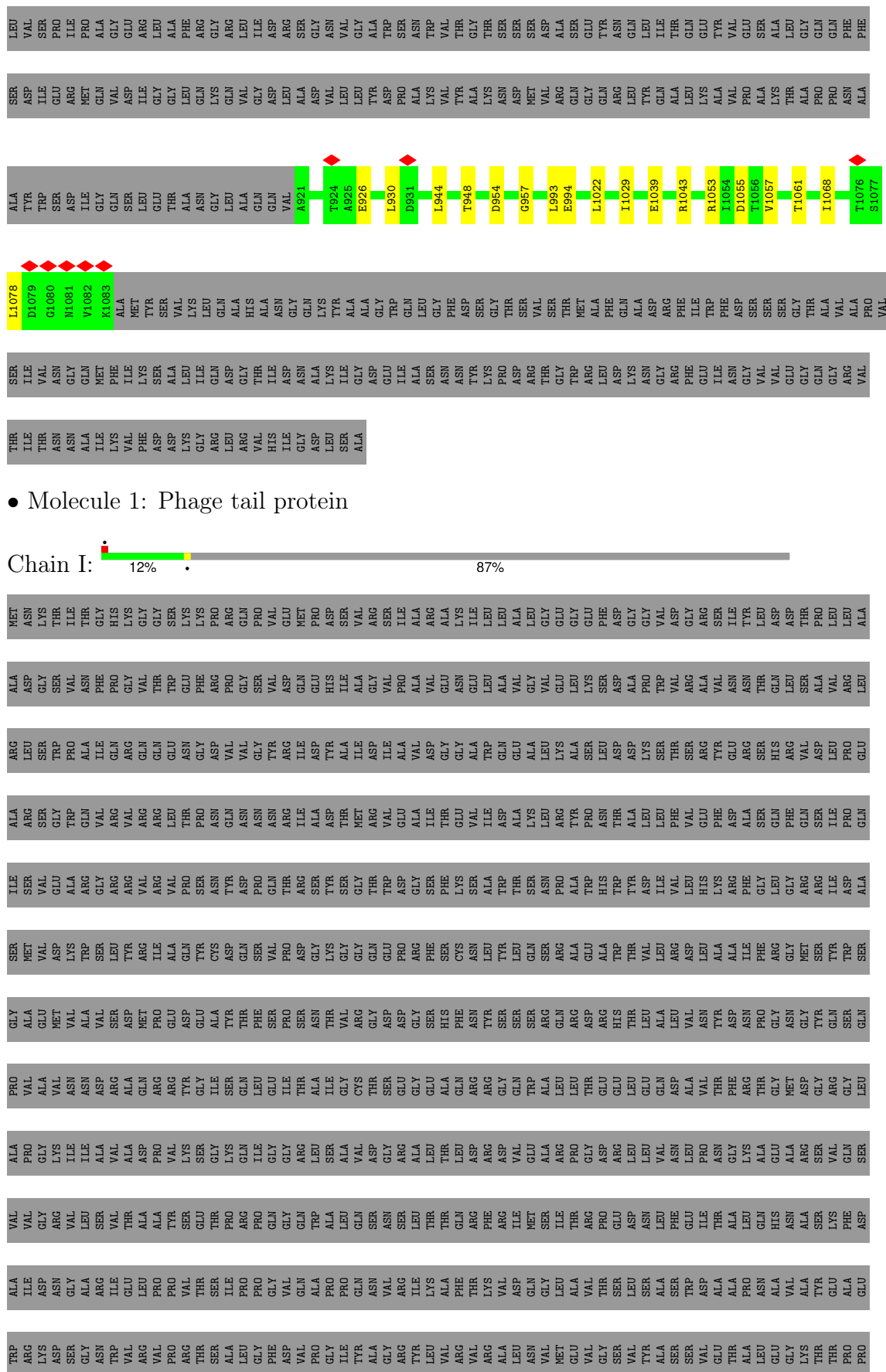
3 Residue-property plots

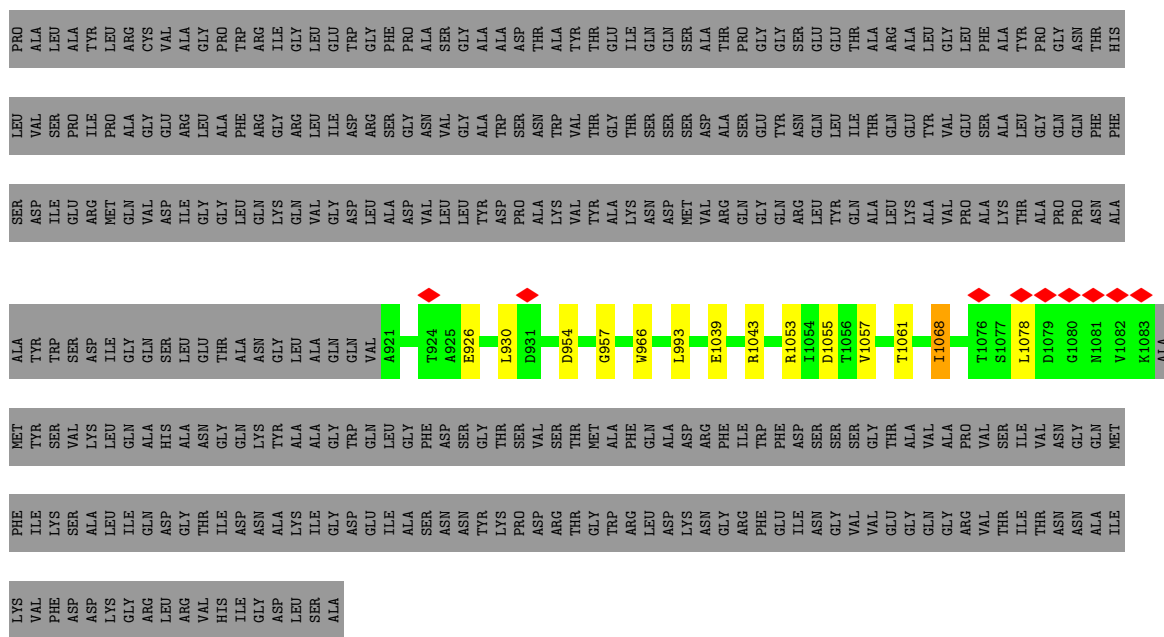
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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: Phage tail protein

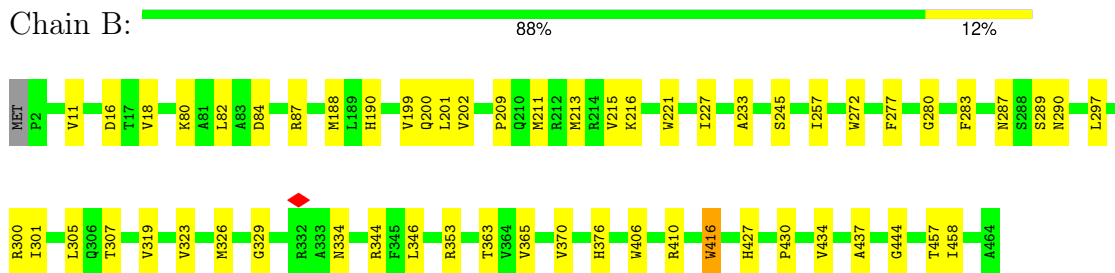




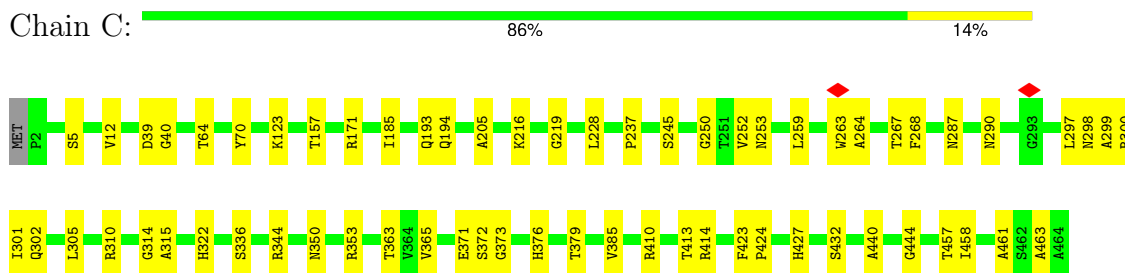




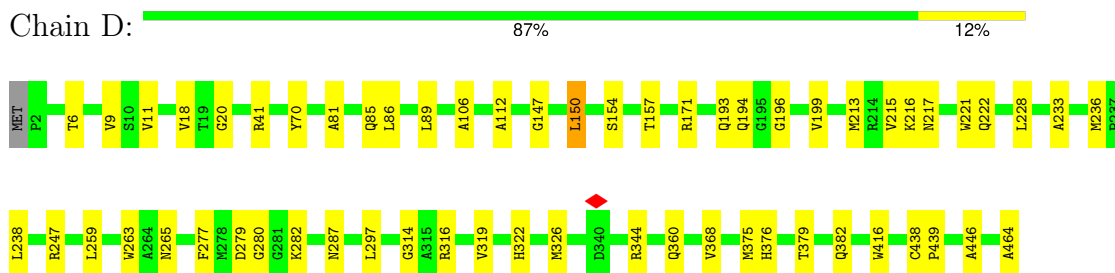
- Molecule 2: Side fiber protein of F7 pyocin



- Molecule 2: Side fiber protein of F7 pyocin

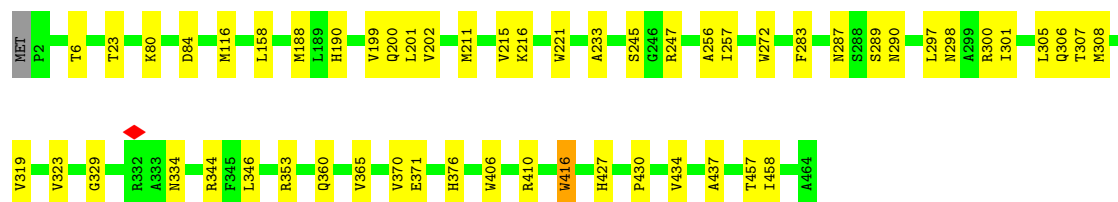


- Molecule 2: Side fiber protein of F7 pyocin



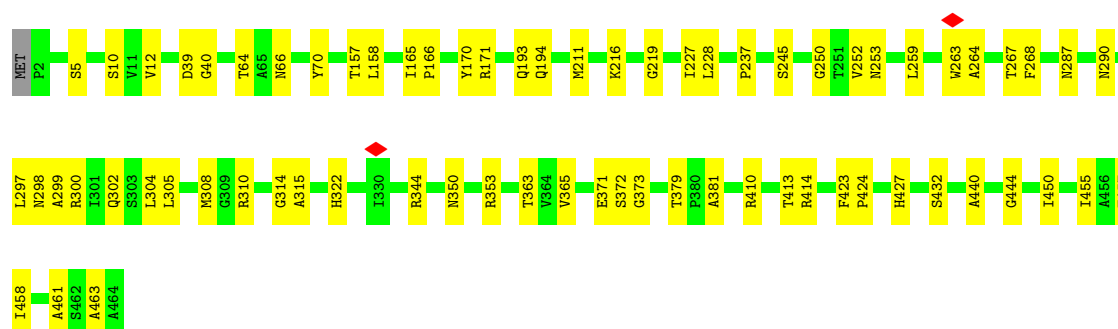
- Molecule 2: Side fiber protein of F7 pyocin

Chain F:  88% 12%



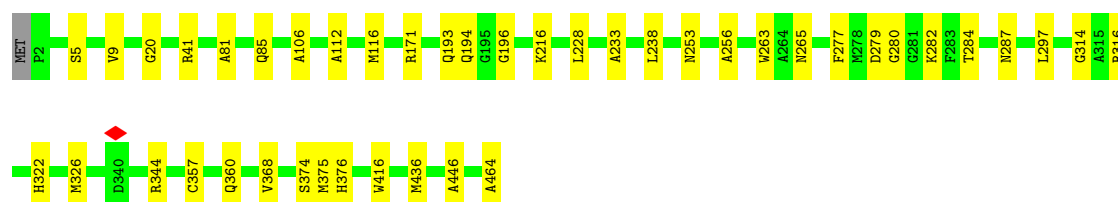
- Molecule 2: Side fiber protein of F7 pyocin

Chain G:  85% 15%



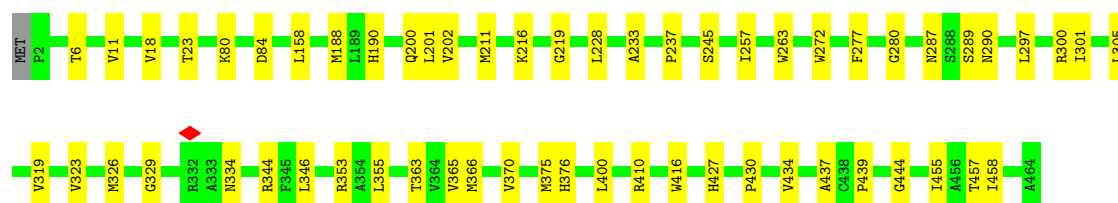
- Molecule 2: Side fiber protein of F7 pyocin

Chain H:  91% 9%



- Molecule 2: Side fiber protein of F7 pyocin

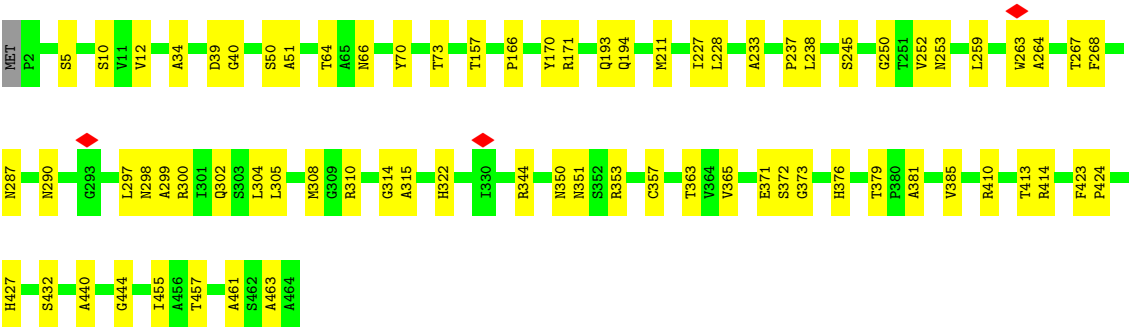
Chain J:  87% 12%



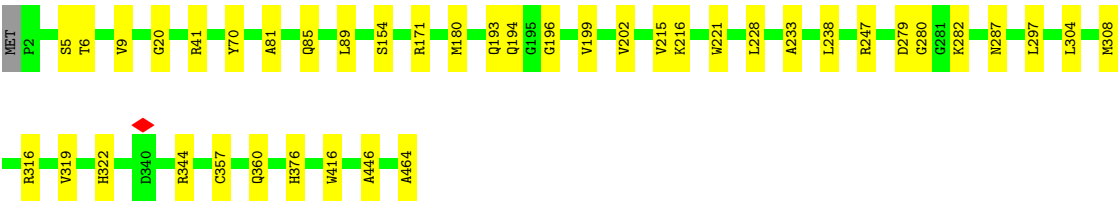
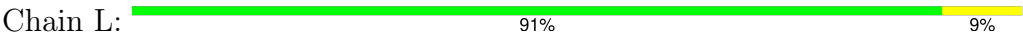
- Molecule 2: Side fiber protein of F7 pyocin

Chain K:  84% 16%





• Molecule 2: Side fiber protein of F7 pyocin



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	359107	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 ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	Not provided	
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	0.699	Depositor
Minimum map value	-0.275	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.016	Depositor
Recommended contour level	0.12	Depositor
Map size ($\text{\AA}$)	440.0, 440.0, 440.0	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	1.1, 1.1, 1.1	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.23	0/1180	0.29	0/1602
1	E	0.22	0/1180	0.29	0/1602
1	I	0.23	0/1180	0.30	0/1602
2	B	0.17	0/3429	0.34	0/4688
2	C	0.14	0/3429	0.29	0/4688
2	D	0.13	0/3429	0.27	0/4688
2	F	0.21	0/3429	0.38	0/4688
2	G	0.12	0/3429	0.27	0/4688
2	H	0.13	0/3429	0.28	0/4688
2	J	0.19	0/3429	0.34	0/4688
2	K	0.12	0/3429	0.27	0/4688
2	L	0.12	0/3429	0.27	0/4688
All	All	0.16	0/34401	0.30	0/46998

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1178	0	1144	19	0
1	E	1178	0	1144	17	0
1	I	1178	0	1144	13	0
2	B	3348	0	3241	38	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	C	3348	0	3241	43	0
2	D	3348	0	3241	40	0
2	F	3348	0	3241	35	0
2	G	3348	0	3241	46	0
2	H	3348	0	3241	29	0
2	J	3348	0	3241	35	0
2	K	3348	0	3241	49	0
2	L	3348	0	3241	29	0
All	All	33666	0	32601	341	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 5.

All (341) 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:E:1068:ILE:HG13	1:I:1068:ILE:HD11	1.74	0.68
2:C:365:VAL:HG12	2:C:410:ARG:HG3	1.78	0.65
2:G:228:LEU:HD11	2:H:233:ALA:HB2	1.79	0.64
2:G:365:VAL:HG12	2:G:410:ARG:HG3	1.80	0.64
2:B:370:VAL:HG22	2:B:434:VAL:HG22	1.80	0.62
2:C:228:LEU:HD11	2:D:233:ALA:HB2	1.81	0.62
2:G:379:THR:HG22	2:G:381:ALA:H	1.65	0.62
2:K:379:THR:HG22	2:K:381:ALA:H	1.64	0.61
2:G:12:VAL:HA	2:G:64:THR:HG23	1.81	0.61
2:J:216:LYS:HE3	2:J:219:GLY:HA2	1.82	0.61
2:B:188:MET:HB2	2:B:202:VAL:HG23	1.83	0.61
2:K:12:VAL:HA	2:K:64:THR:HG23	1.83	0.61
2:K:50:SER:OG	2:K:51:ALA:N	2.33	0.61
2:C:12:VAL:HA	2:C:64:THR:HG23	1.82	0.60
2:B:233:ALA:HB2	2:D:228:LEU:HD11	1.83	0.60
2:C:253:ASN:HD21	2:C:463:ALA:H	1.49	0.60
2:F:300:ARG:NH2	2:F:371:GLU:OE2	2.35	0.59
2:K:253:ASN:HD21	2:K:463:ALA:H	1.50	0.59
1:A:1068:ILE:HD11	1:I:1068:ILE:HG13	1.84	0.59
2:K:413:THR:OG1	2:K:414:ARG:N	2.36	0.59
1:A:1053:ARG:NH2	1:E:1055:ASP:OD1	2.36	0.59
2:J:233:ALA:HB2	2:L:228:LEU:HD11	1.85	0.59
2:F:233:ALA:HB2	2:H:228:LEU:HD11	1.85	0.58
2:K:252:VAL:HG21	2:K:259:LEU:HD23	1.85	0.58
2:B:201:LEU:HD11	2:B:211:MET:HE3	1.86	0.58

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:1055:ASP:OD1	1:I:1053:ARG:NH2	2.37	0.58
2:K:263:TRP:HB3	2:K:322:HIS:CD2	2.38	0.58
2:K:297:LEU:HD23	2:K:302:GLN:HG3	1.86	0.58
2:J:370:VAL:HG22	2:J:434:VAL:HG22	1.85	0.58
2:G:253:ASN:HD21	2:G:463:ALA:H	1.50	0.58
2:J:257:ILE:HG22	2:J:319:VAL:HG22	1.85	0.58
1:E:1053:ARG:NH2	1:I:1055:ASP:OD1	2.37	0.58
2:F:272:TRP:HH2	2:F:334:ASN:HD21	1.51	0.57
2:F:257:ILE:HG22	2:F:319:VAL:HG22	1.86	0.57
2:K:10:SER:OG	2:K:66:ASN:OD1	2.18	0.57
2:K:365:VAL:HG22	2:K:440:ALA:HB3	1.84	0.57
2:G:263:TRP:HB3	2:G:322:HIS:CD2	2.40	0.57
2:G:363:THR:OG1	2:G:444:GLY:O	2.21	0.57
2:B:272:TRP:HH2	2:B:334:ASN:HD21	1.51	0.57
2:C:263:TRP:HB3	2:C:322:HIS:CD2	2.39	0.57
2:G:252:VAL:HG21	2:G:259:LEU:HD23	1.87	0.57
2:K:365:VAL:HG12	2:K:410:ARG:HG3	1.85	0.57
2:F:188:MET:HB2	2:F:202:VAL:HG23	1.87	0.57
2:G:300:ARG:NH1	2:G:371:GLU:OE2	2.39	0.56
2:K:263:TRP:HB3	2:K:322:HIS:HD2	1.70	0.56
2:C:252:VAL:HG21	2:C:259:LEU:HD23	1.87	0.56
2:K:300:ARG:NH1	2:K:371:GLU:OE2	2.39	0.56
2:J:301:ILE:O	2:J:305:LEU:HG	2.06	0.55
2:K:228:LEU:HD11	2:L:233:ALA:HB2	1.87	0.55
2:J:272:TRP:HH2	2:J:334:ASN:HD21	1.54	0.55
2:C:157:THR:OG1	2:C:171:ARG:NE	2.39	0.55
2:G:157:THR:OG1	2:G:171:ARG:NE	2.37	0.55
2:J:201:LEU:HD11	2:J:211:MET:HE3	1.89	0.55
2:C:297:LEU:HD23	2:C:302:GLN:HG3	1.89	0.55
2:L:154:SER:O	2:L:154:SER:OG	2.21	0.55
2:C:365:VAL:HG22	2:C:440:ALA:HB3	1.89	0.54
2:C:300:ARG:NH1	2:C:371:GLU:OE2	2.39	0.54
2:J:363:THR:OG1	2:J:444:GLY:O	2.21	0.54
2:F:370:VAL:HG22	2:F:434:VAL:HG22	1.89	0.54
2:G:237:PRO:HG2	2:H:238:LEU:HD22	1.89	0.54
2:K:157:THR:OG1	2:K:171:ARG:NE	2.39	0.54
2:B:287:ASN:HD21	2:B:289:SER:HB2	1.73	0.54
2:C:263:TRP:HB3	2:C:322:HIS:HD2	1.71	0.54
2:G:263:TRP:HB3	2:G:322:HIS:HD2	1.71	0.54
2:C:185:ILE:HD12	2:C:205:ALA:HB3	1.90	0.54
2:D:147:GLY:HA2	2:D:150:LEU:HD12	1.89	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:K:363:THR:OG1	2:K:444:GLY:O	2.25	0.54
2:F:287:ASN:HD21	2:F:289:SER:HB2	1.72	0.53
2:G:365:VAL:HG22	2:G:440:ALA:HB3	1.89	0.53
2:F:201:LEU:HD11	2:F:211:MET:HE3	1.89	0.53
2:C:237:PRO:HG2	2:D:238:LEU:HD22	1.91	0.53
2:B:209:PRO:HB2	2:D:213:MET:HE3	1.90	0.53
2:B:257:ILE:HG22	2:B:319:VAL:HG22	1.89	0.53
2:K:237:PRO:HG2	2:L:238:LEU:HD22	1.91	0.53
2:B:190:HIS:CE1	2:B:200:GLN:HE21	2.27	0.53
2:B:363:THR:OG1	2:B:444:GLY:O	2.21	0.53
2:G:158:LEU:HD11	2:G:170:TYR:HB3	1.92	0.52
2:J:190:HIS:CE1	2:J:200:GLN:HE21	2.27	0.52
2:B:307:THR:HG21	2:B:406:TRP:HB2	1.92	0.51
2:K:455:ILE:HG23	2:L:357:CYS:HB3	1.93	0.51
2:J:287:ASN:HD21	2:J:289:SER:HB2	1.73	0.51
2:G:211:MET:HG2	2:G:227:ILE:HD13	1.92	0.51
2:B:283:PHE:HE2	2:B:305:LEU:HD21	1.76	0.51
2:K:34:ALA:HB3	2:K:73:THR:HB	1.93	0.50
2:G:304:LEU:HG	2:G:308:MET:HE3	1.94	0.50
2:J:228:LEU:HD11	2:K:233:ALA:HB2	1.93	0.50
2:C:413:THR:OG1	2:C:414:ARG:N	2.38	0.50
2:G:193:GLN:HG2	2:G:194:GLN:H	1.76	0.50
2:B:84:ASP:HB2	1:I:957:GLY:HA3	1.93	0.50
2:B:353:ARG:NH2	2:D:464:ALA:O	2.40	0.50
2:K:193:GLN:HG2	2:K:194:GLN:H	1.76	0.50
2:L:304:LEU:O	2:L:308:MET:HG3	2.12	0.50
1:I:1057:VAL:O	1:I:1061:THR:HG23	2.12	0.50
2:F:190:HIS:CE1	2:F:200:GLN:HE21	2.30	0.49
2:F:287:ASN:ND2	2:F:290:ASN:OD1	2.44	0.49
2:J:297:LEU:HD13	2:J:301:ILE:HG21	1.95	0.49
2:C:372:SER:O	2:C:432:SER:OG	2.27	0.49
2:B:287:ASN:ND2	2:B:290:ASN:OD1	2.45	0.49
1:E:1057:VAL:O	1:E:1061:THR:HG23	2.12	0.49
1:I:954:ASP:HB3	2:J:80:LYS:HD2	1.94	0.49
1:A:1039:GLU:OE1	1:E:1043:ARG:NH1	2.46	0.49
1:A:1057:VAL:O	1:A:1061:THR:HG23	2.12	0.49
2:J:277:PHE:CZ	2:J:326:MET:HG3	2.47	0.49
2:C:305:LEU:HD13	2:C:315:ALA:HA	1.94	0.49
2:G:287:ASN:HD21	2:G:290:ASN:HD22	1.59	0.49
2:C:39:ASP:OD1	2:C:39:ASP:N	2.40	0.49
2:K:267:THR:HG21	2:K:353:ARG:HH21	1.78	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:B:84:ASP:HA	2:B:87:ARG:HB2	1.95	0.48
2:F:297:LEU:HD13	2:F:301:ILE:HG21	1.95	0.48
2:F:344:ARG:HB2	2:F:376:HIS:CE1	2.48	0.48
2:C:363:THR:OG1	2:C:444:GLY:O	2.28	0.48
2:G:305:LEU:HD13	2:G:315:ALA:HA	1.95	0.48
2:J:188:MET:HB2	2:J:202:VAL:HG23	1.94	0.48
2:B:277:PHE:CZ	2:B:326:MET:HG3	2.49	0.48
2:F:283:PHE:HE2	2:F:305:LEU:HD21	1.79	0.48
2:G:413:THR:OG1	2:G:414:ARG:N	2.47	0.48
2:G:237:PRO:HG3	2:H:446:ALA:HB2	1.96	0.48
2:G:253:ASN:ND2	2:G:461:ALA:O	2.44	0.48
2:J:280:GLY:O	2:J:300:ARG:NH1	2.47	0.48
2:K:310:ARG:HB3	2:K:314:GLY:HA3	1.95	0.48
1:E:954:ASP:HB3	2:F:80:LYS:HD2	1.96	0.48
2:F:346:LEU:HA	2:F:427:HIS:HD2	1.78	0.47
2:J:344:ARG:HB2	2:J:376:HIS:CE1	2.49	0.47
2:B:344:ARG:HD3	2:B:376:HIS:ND1	2.29	0.47
2:G:267:THR:HG21	2:G:353:ARG:HH21	1.79	0.47
2:L:81:ALA:O	2:L:85:GLN:HG3	2.14	0.47
2:B:457:THR:OG1	2:B:458:ILE:N	2.47	0.47
2:D:199:VAL:HG23	2:D:215:VAL:HG12	1.96	0.47
2:F:344:ARG:HD3	2:F:376:HIS:ND1	2.29	0.47
2:H:81:ALA:O	2:H:85:GLN:HG3	2.15	0.47
2:H:287:ASN:OD1	2:H:287:ASN:N	2.41	0.47
2:K:237:PRO:HG3	2:L:446:ALA:HB2	1.97	0.47
2:B:297:LEU:HD13	2:B:301:ILE:HG21	1.96	0.47
1:E:1039:GLU:OE1	1:I:1043:ARG:NH1	2.48	0.47
2:G:297:LEU:HD23	2:G:302:GLN:HG3	1.96	0.47
2:K:305:LEU:HD13	2:K:315:ALA:HA	1.97	0.47
2:F:245:SER:HB2	2:F:457:THR:HG22	1.95	0.47
2:K:287:ASN:HD21	2:K:290:ASN:HD22	1.60	0.47
2:C:5:SER:HB2	2:C:70:TYR:O	2.14	0.47
2:B:199:VAL:HG23	2:B:215:VAL:HG12	1.97	0.47
2:H:9:VAL:HA	2:H:20:GLY:HA2	1.97	0.47
2:J:344:ARG:HD3	2:J:376:HIS:ND1	2.30	0.47
2:J:366:MET:HB3	2:J:439:PRO:HA	1.97	0.47
2:K:253:ASN:ND2	2:K:461:ALA:O	2.42	0.47
2:F:457:THR:OG1	2:F:458:ILE:N	2.48	0.47
2:K:250:GLY:H	2:K:264:ALA:HB3	1.80	0.47
2:B:280:GLY:O	2:B:300:ARG:NH1	2.48	0.47
2:B:416:TRP:CZ3	2:D:314:GLY:HA3	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:D:81:ALA:O	2:D:85:GLN:HG3	2.15	0.47
2:J:457:THR:OG1	2:J:458:ILE:N	2.47	0.47
2:F:306:GLN:O	2:F:307:THR:C	2.57	0.46
2:C:253:ASN:ND2	2:C:461:ALA:O	2.43	0.46
2:G:310:ARG:HB3	2:G:314:GLY:HA3	1.97	0.46
2:D:287:ASN:HD22	2:D:297:LEU:HG	1.80	0.46
2:C:336:SER:OG	2:C:379:THR:OG1	2.22	0.46
2:L:344:ARG:HD2	2:L:376:HIS:ND1	2.30	0.46
2:B:11:VAL:HG12	2:B:18:VAL:HG22	1.98	0.46
2:C:385:VAL:HG22	2:C:413:THR:HB	1.98	0.46
2:C:310:ARG:HB3	2:C:314:GLY:HA3	1.98	0.46
2:F:199:VAL:HG23	2:F:215:VAL:HG12	1.98	0.46
2:B:216:LYS:HB3	2:B:221:TRP:CE3	2.51	0.46
2:G:40:GLY:HA3	2:H:41:ARG:HG2	1.98	0.46
2:D:368:VAL:HG11	2:D:375:MET:HE1	1.98	0.46
2:H:360:GLN:HG2	2:H:416:TRP:HA	1.98	0.46
2:J:365:VAL:HG12	2:J:410:ARG:HG3	1.98	0.46
2:G:245:SER:HB2	2:G:457:THR:HG22	1.98	0.46
1:A:1067:ALA:HB3	1:E:1068:ILE:HD13	1.98	0.45
2:C:193:GLN:HG2	2:C:194:GLN:H	1.80	0.45
2:D:9:VAL:HA	2:D:20:GLY:HA2	1.97	0.45
2:G:344:ARG:HB3	2:G:427:HIS:HB3	1.98	0.45
2:C:344:ARG:HB3	2:C:427:HIS:HB3	1.99	0.45
2:D:344:ARG:HD2	2:D:376:HIS:ND1	2.31	0.45
2:H:287:ASN:HD22	2:H:297:LEU:HG	1.81	0.45
2:L:287:ASN:HD22	2:L:297:LEU:HG	1.81	0.45
2:D:11:VAL:HG12	2:D:18:VAL:HG22	1.98	0.45
2:G:193:GLN:CD	2:H:171:ARG:HE	2.25	0.45
2:F:323:VAL:HG13	2:F:437:ALA:HB2	1.97	0.45
2:H:106:ALA:HB3	2:H:112:ALA:HB2	1.97	0.45
2:B:344:ARG:HB2	2:B:376:HIS:CE1	2.51	0.45
2:F:329:GLY:HA2	2:F:430:PRO:HB3	1.98	0.45
2:J:455:ILE:HA	2:K:357:CYS:HB3	1.99	0.45
2:C:245:SER:HB2	2:C:457:THR:HG22	1.97	0.45
2:C:267:THR:HG21	2:C:353:ARG:HH21	1.82	0.45
2:F:84:ASP:OD2	2:F:84:ASP:C	2.60	0.45
2:G:250:GLY:H	2:G:264:ALA:HB3	1.81	0.45
2:B:346:LEU:HA	2:B:427:HIS:HD2	1.82	0.45
2:K:245:SER:HB2	2:K:457:THR:HG22	1.98	0.45
2:H:193:GLN:HG2	2:H:194:GLN:N	2.31	0.45
2:L:6:THR:O	2:L:70:TYR:OH	2.31	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:957:GLY:HA3	2:J:84:ASP:HB2	1.99	0.45
2:F:365:VAL:HG12	2:F:410:ARG:HG3	1.98	0.45
2:F:416:TRP:CZ3	2:H:314:GLY:HA3	2.52	0.45
2:J:263:TRP:HH2	2:J:277:PHE:HB3	1.82	0.45
2:F:307:THR:HG21	2:F:406:TRP:HB2	1.99	0.44
2:K:344:ARG:HB3	2:K:427:HIS:HB3	1.99	0.44
2:G:5:SER:HB2	2:G:70:TYR:O	2.16	0.44
1:I:1078:LEU:HD12	1:I:1078:LEU:HA	1.85	0.44
2:L:287:ASN:OD1	2:L:287:ASN:N	2.42	0.44
2:H:5:SER:O	2:H:5:SER:OG	2.33	0.44
2:H:253:ASN:HD22	2:H:256:ALA:H	1.66	0.44
2:H:368:VAL:HG11	2:H:375:MET:HE1	1.98	0.44
2:J:287:ASN:ND2	2:J:290:ASN:OD1	2.50	0.44
2:F:360:GLN:HG2	2:F:416:TRP:HA	1.98	0.44
2:J:11:VAL:HG12	2:J:18:VAL:HG22	1.99	0.44
2:K:211:MET:HG2	2:K:227:ILE:HD13	1.99	0.44
2:J:323:VAL:HG13	2:J:437:ALA:HB2	1.99	0.44
2:D:379:THR:HG22	2:D:382:GLN:HG3	1.99	0.44
2:K:5:SER:HB2	2:K:70:TYR:O	2.17	0.44
2:K:372:SER:O	2:K:432:SER:OG	2.26	0.44
2:D:193:GLN:HG2	2:D:194:GLN:N	2.33	0.44
2:D:287:ASN:OD1	2:D:287:ASN:N	2.40	0.44
2:G:157:THR:HG1	2:G:171:ARG:HE	1.63	0.44
2:H:287:ASN:OD1	2:H:316:ARG:HA	2.18	0.44
2:L:193:GLN:HG2	2:L:194:GLN:N	2.33	0.43
2:L:304:LEU:HD11	2:L:308:MET:HE3	2.00	0.43
2:D:259:LEU:HD23	2:D:259:LEU:HA	1.87	0.43
2:G:268:PHE:O	2:G:350:ASN:HA	2.18	0.43
2:H:277:PHE:CE2	2:H:326:MET:HB2	2.54	0.43
1:I:966:TRP:CD1	2:L:89:LEU:HD21	2.54	0.43
1:A:1050:LEU:HA	1:A:1050:LEU:HD23	1.83	0.43
2:D:247:ARG:HH21	2:D:319:VAL:HG13	1.83	0.43
2:K:304:LEU:O	2:K:308:MET:HG3	2.18	0.43
2:D:154:SER:O	2:D:154:SER:OG	2.28	0.43
2:D:263:TRP:CH2	2:D:265:ASN:HB2	2.53	0.43
2:G:450:ILE:HD13	2:G:450:ILE:HA	1.87	0.43
2:H:284:THR:H	2:H:297:LEU:HD21	1.84	0.43
2:K:298:ASN:OD1	2:K:299:ALA:N	2.51	0.43
1:A:1068:ILE:HG13	1:E:1068:ILE:HD11	1.99	0.43
2:G:10:SER:OG	2:G:66:ASN:OD1	2.21	0.43
2:G:372:SER:O	2:G:432:SER:OG	2.26	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:H:344:ARG:HD2	2:H:376:HIS:ND1	2.33	0.43
2:K:423:PHE:CD1	2:K:424:PRO:HA	2.53	0.43
2:L:180:MET:HE3	2:L:202:VAL:HG12	2.01	0.43
1:A:994:GLU:HG2	1:I:993:LEU:HD21	2.00	0.43
1:A:1043:ARG:NH1	1:I:1039:GLU:OE1	2.52	0.43
2:F:158:LEU:HD23	2:F:158:LEU:HA	1.91	0.43
2:G:298:ASN:OD1	2:G:299:ALA:N	2.51	0.43
1:I:926:GLU:O	1:I:930:LEU:HG	2.19	0.43
2:K:268:PHE:O	2:K:350:ASN:HA	2.19	0.43
2:C:250:GLY:H	2:C:264:ALA:HB3	1.82	0.43
2:G:457:THR:OG1	2:G:458:ILE:N	2.52	0.43
2:J:375:MET:HE2	2:J:400:LEU:HD12	2.01	0.43
2:D:154:SER:OG	2:D:157:THR:OG1	2.36	0.43
2:D:360:GLN:HG2	2:D:416:TRP:HA	2.01	0.43
2:L:9:VAL:HA	2:L:20:GLY:HA2	2.00	0.43
2:G:157:THR:HG21	2:G:171:ARG:HH21	1.84	0.42
2:J:329:GLY:HA2	2:J:430:PRO:HB3	2.01	0.42
2:K:40:GLY:HA3	2:L:41:ARG:HG2	2.01	0.42
2:C:40:GLY:HA3	2:D:41:ARG:HG2	2.01	0.42
2:D:279:ASP:OD1	2:D:280:GLY:N	2.52	0.42
2:F:353:ARG:NH2	2:H:464:ALA:O	2.43	0.42
2:K:39:ASP:OD1	2:K:39:ASP:N	2.43	0.42
2:K:193:GLN:CD	2:L:171:ARG:HE	2.26	0.42
1:A:1021:ARG:HA	1:A:1021:ARG:HD2	1.84	0.42
2:F:116:MET:HE3	2:H:116:MET:HE1	2.01	0.42
2:G:39:ASP:OD1	2:G:39:ASP:N	2.41	0.42
2:C:193:GLN:CD	2:D:171:ARG:HE	2.27	0.42
2:D:196:GLY:O	2:D:216:LYS:HG2	2.19	0.42
2:D:216:LYS:HD3	2:D:221:TRP:CE2	2.55	0.42
2:D:277:PHE:CE2	2:D:326:MET:HB2	2.55	0.42
2:J:245:SER:HB2	2:J:457:THR:HG22	2.01	0.42
2:C:268:PHE:O	2:C:350:ASN:HA	2.19	0.42
2:K:373:GLY:HA3	2:K:432:SER:HB2	2.02	0.42
2:D:6:THR:O	2:D:70:TYR:OH	2.29	0.42
1:E:926:GLU:O	1:E:930:LEU:HG	2.19	0.42
2:H:436:MET:HE3	2:H:436:MET:HB2	1.93	0.42
2:K:166:PRO:O	2:K:170:TYR:OH	2.31	0.42
2:B:416:TRP:HZ3	2:D:314:GLY:HA3	1.85	0.42
2:C:457:THR:OG1	2:C:458:ILE:N	2.52	0.42
2:D:106:ALA:HB3	2:D:112:ALA:HB2	2.02	0.42
2:G:216:LYS:HE2	2:G:219:GLY:HA2	2.02	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:L:5:SER:O	2:L:5:SER:OG	2.34	0.42
2:L:196:GLY:O	2:L:216:LYS:HG2	2.20	0.42
1:A:926:GLU:O	1:A:930:LEU:HG	2.19	0.41
2:B:16:ASP:OD1	2:B:16:ASP:N	2.51	0.41
2:D:217:ASN:HB3	2:D:222:GLN:HE21	1.85	0.41
2:H:196:GLY:O	2:H:216:LYS:HG2	2.20	0.41
2:H:282:LYS:HB2	2:H:322:HIS:CE1	2.55	0.41
1:A:954:ASP:HB3	2:B:80:LYS:HD2	2.01	0.41
2:H:263:TRP:CH2	2:H:265:ASN:HB2	2.55	0.41
2:L:199:VAL:HG23	2:L:215:VAL:HG12	2.00	0.41
1:A:966:TRP:CD1	2:D:89:LEU:HD21	2.55	0.41
2:B:283:PHE:CE2	2:B:305:LEU:HD21	2.55	0.41
2:C:423:PHE:CD1	2:C:424:PRO:HA	2.55	0.41
1:E:1078:LEU:HD12	1:E:1078:LEU:HA	1.85	0.41
2:G:165:ILE:HG22	2:G:166:PRO:HD2	2.01	0.41
2:C:237:PRO:HG3	2:D:446:ALA:HB2	2.02	0.41
1:E:1022:LEU:HD23	1:E:1022:LEU:HA	1.93	0.41
2:L:360:GLN:HG2	2:L:416:TRP:HA	2.02	0.41
2:B:323:VAL:HG13	2:B:437:ALA:HB2	2.01	0.41
2:B:365:VAL:HG12	2:B:410:ARG:HG3	2.02	0.41
2:C:301:ILE:O	2:C:305:LEU:HG	2.21	0.41
2:F:247:ARG:NH2	2:F:256:ALA:O	2.53	0.41
2:G:455:ILE:HG23	2:H:357:CYS:HB3	2.02	0.41
2:C:298:ASN:OD1	2:C:299:ALA:N	2.53	0.41
1:E:993:LEU:HD23	1:E:993:LEU:HA	1.83	0.41
2:J:158:LEU:HD23	2:J:158:LEU:HA	1.91	0.41
2:K:385:VAL:HG22	2:K:413:THR:HB	2.02	0.41
2:B:245:SER:HB2	2:B:457:THR:HG22	2.02	0.41
2:B:329:GLY:HA2	2:B:430:PRO:HB3	2.01	0.41
2:D:282:LYS:HB2	2:D:322:HIS:CE1	2.55	0.41
2:J:353:ARG:NH2	2:L:464:ALA:O	2.45	0.41
2:B:82:LEU:HD21	2:D:86:LEU:HD23	2.03	0.41
2:C:287:ASN:HD21	2:C:290:ASN:HD22	1.68	0.41
1:E:944:LEU:O	1:E:948:THR:HG23	2.21	0.41
2:F:216:LYS:HB3	2:F:221:TRP:CE3	2.56	0.41
2:G:373:GLY:HA3	2:G:432:SER:HB2	2.02	0.41
2:J:237:PRO:HG2	2:K:238:LEU:HD21	2.03	0.41
2:L:287:ASN:OD1	2:L:316:ARG:HA	2.21	0.41
1:A:977:THR:O	1:A:981:THR:HG23	2.21	0.41
2:C:123:LYS:HB2	2:C:123:LYS:HE2	1.91	0.41
2:C:216:LYS:HE2	2:C:219:GLY:HA2	2.03	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:344:ARG:HB2	2:C:376:HIS:CE1	2.55	0.41
2:C:373:GLY:HA3	2:C:432:SER:HB2	2.03	0.41
2:D:438:CYS:N	2:D:439:PRO:HD3	2.36	0.41
2:J:346:LEU:HA	2:J:427:HIS:HD2	1.86	0.41
2:L:279:ASP:OD1	2:L:280:GLY:N	2.53	0.41
1:A:944:LEU:O	1:A:948:THR:HG23	2.21	0.41
2:C:287:ASN:OD1	2:C:287:ASN:N	2.54	0.41
2:F:6:THR:O	2:F:23:THR:HG22	2.21	0.41
2:H:279:ASP:OD1	2:H:280:GLY:N	2.54	0.41
2:K:287:ASN:N	2:K:287:ASN:OD1	2.53	0.41
1:A:1029:ILE:HD13	1:E:1029:ILE:HD13	2.03	0.40
2:B:365:VAL:HG12	2:B:410:ARG:CG	2.51	0.40
2:D:287:ASN:OD1	2:D:316:ARG:HA	2.21	0.40
2:F:308:MET:HE2	2:F:308:MET:HB2	1.96	0.40
2:G:158:LEU:HD11	2:G:170:TYR:CB	2.51	0.40
2:L:216:LYS:HB2	2:L:221:TRP:CZ3	2.56	0.40
1:A:993:LEU:HD21	1:E:994:GLU:HG2	2.03	0.40
2:D:233:ALA:HA	2:D:236:MET:HG3	2.02	0.40
2:F:298:ASN:OD1	2:F:298:ASN:N	2.54	0.40
1:A:1022:LEU:HD23	1:A:1022:LEU:HA	1.93	0.40
2:B:213:MET:HE2	2:B:227:ILE:HD11	2.03	0.40
2:C:287:ASN:ND2	2:C:297:LEU:HD22	2.36	0.40
2:K:287:ASN:HD22	2:K:297:LEU:HD22	1.87	0.40
2:K:344:ARG:HB2	2:K:376:HIS:CE1	2.56	0.40
2:J:6:THR:O	2:J:23:THR:HG22	2.22	0.40
2:K:267:THR:HG23	2:K:351:ASN:HD22	1.86	0.40
2:L:247:ARG:HH21	2:L:319:VAL:HG13	1.86	0.40
2:G:423:PHE:CD1	2:G:424:PRO:HA	2.57	0.40
2:L:282:LYS:HB2	2:L:322:HIS:CE1	2.56	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	161/1222 (13%)	161 (100%)	0	0	100	100
1	E	161/1222 (13%)	161 (100%)	0	0	100	100
1	I	161/1222 (13%)	161 (100%)	0	0	100	100
2	B	461/464 (99%)	449 (97%)	12 (3%)	0	100	100
2	C	461/464 (99%)	449 (97%)	12 (3%)	0	100	100
2	D	461/464 (99%)	448 (97%)	13 (3%)	0	100	100
2	F	461/464 (99%)	450 (98%)	11 (2%)	0	100	100
2	G	461/464 (99%)	454 (98%)	7 (2%)	0	100	100
2	H	461/464 (99%)	448 (97%)	13 (3%)	0	100	100
2	J	461/464 (99%)	449 (97%)	12 (3%)	0	100	100
2	K	461/464 (99%)	453 (98%)	8 (2%)	0	100	100
2	L	461/464 (99%)	448 (97%)	13 (3%)	0	100	100
All	All	4632/7842 (59%)	4531 (98%)	101 (2%)	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	119/974 (12%)	119 (100%)	0	100	100
1	E	119/974 (12%)	119 (100%)	0	100	100
1	I	119/974 (12%)	118 (99%)	1 (1%)	73	81
2	B	333/334 (100%)	332 (100%)	1 (0%)	86	87
2	C	333/334 (100%)	333 (100%)	0	100	100
2	D	333/334 (100%)	332 (100%)	1 (0%)	86	87
2	F	333/334 (100%)	332 (100%)	1 (0%)	86	87
2	G	333/334 (100%)	333 (100%)	0	100	100
2	H	333/334 (100%)	332 (100%)	1 (0%)	86	87

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
2	J	333/334 (100%)	331 (99%)	2 (1%)	78	83
2	K	333/334 (100%)	333 (100%)	0	100	100
2	L	333/334 (100%)	333 (100%)	0	100	100
All	All	3354/5928 (57%)	3347 (100%)	7 (0%)	85	89

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	416	TRP
2	D	150	LEU
2	F	416	TRP
2	H	374	SER
1	I	1068	ILE
2	J	355	LEU
2	J	416	TRP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (36) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	1040	GLN
2	B	61	GLN
2	B	187	GLN
2	B	334	ASN
2	B	360	GLN
2	C	162	ASN
2	C	193	GLN
2	C	217	ASN
2	C	290	ASN
2	C	322	HIS
2	D	66	ASN
2	D	217	ASN
2	D	222	GLN
1	E	1040	GLN
2	F	61	GLN
2	F	68	GLN
2	F	187	GLN
2	F	334	ASN
2	F	360	GLN
2	G	47	ASN
2	G	162	ASN

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Mol	Chain	Res	Type
2	G	217	ASN
2	G	290	ASN
2	G	322	HIS
2	H	217	ASN
2	H	306	GLN
2	J	61	GLN
2	J	187	GLN
2	J	261	GLN
2	J	360	GLN
2	K	47	ASN
2	K	162	ASN
2	K	290	ASN
2	K	322	HIS
2	L	66	ASN
2	L	217	ASN

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-72915. These allow visual inspection of the internal detail of the map and identification of artifacts.

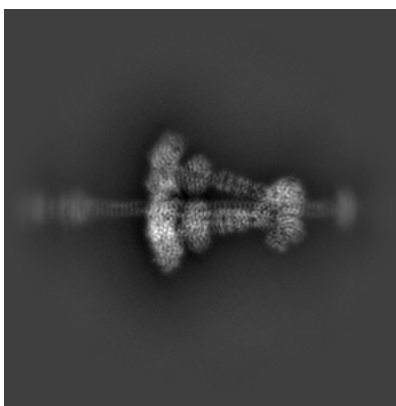
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

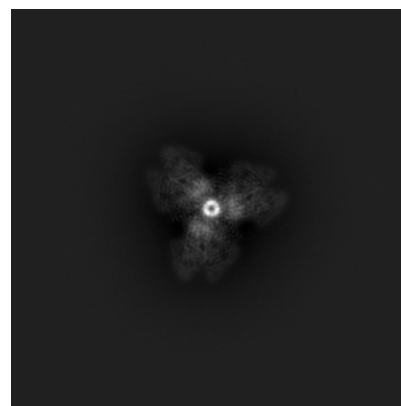
6.1.1 Primary map



X

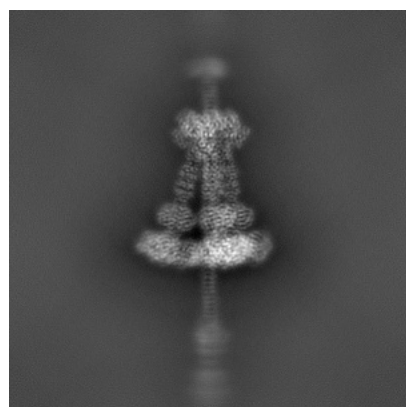


Y

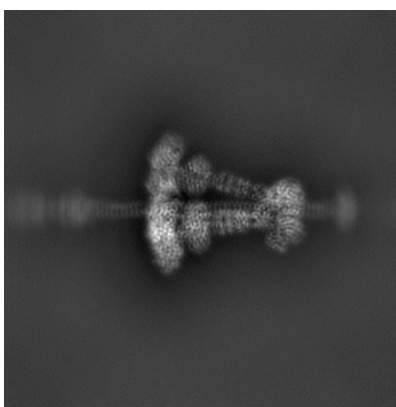


Z

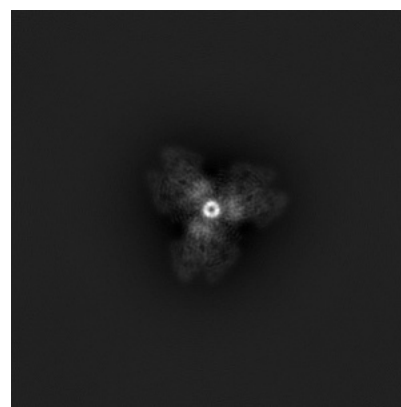
6.1.2 Raw map



X



Y



Z

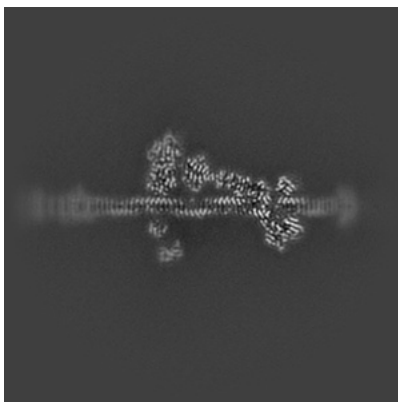
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

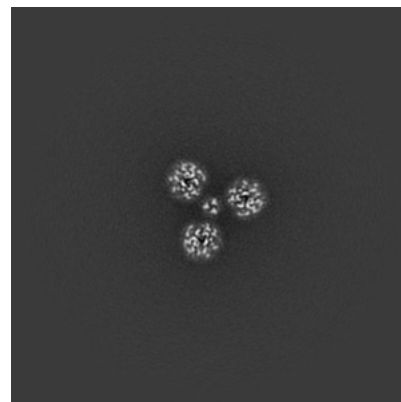
6.2.1 Primary map



X Index: 200

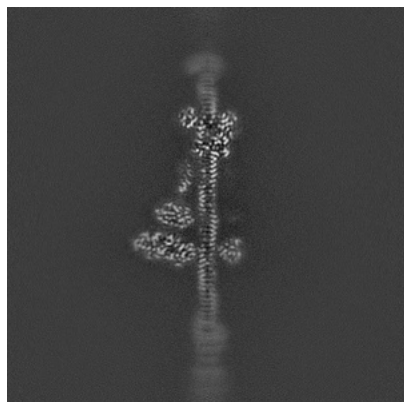


Y Index: 200

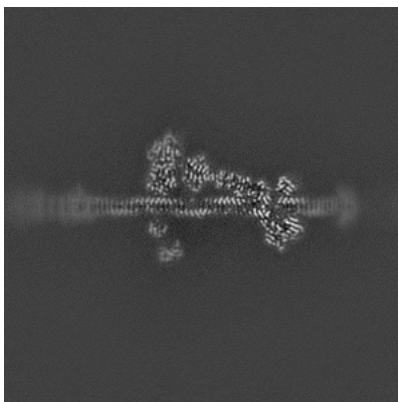


Z Index: 200

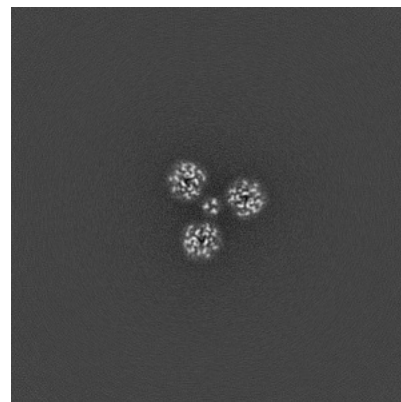
6.2.2 Raw map



X Index: 200



Y Index: 200

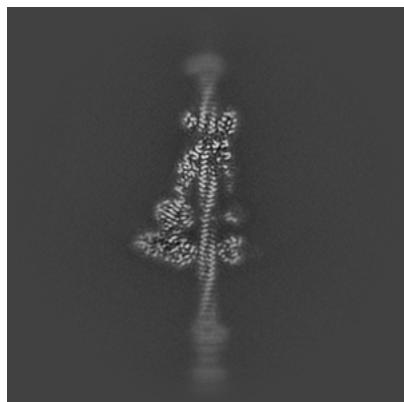


Z Index: 200

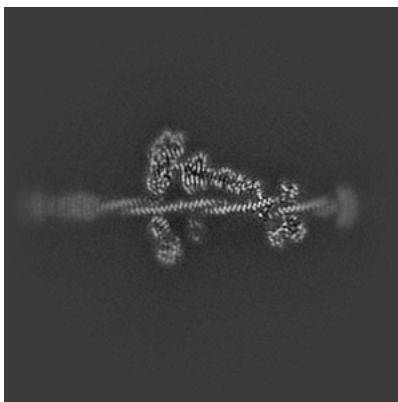
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

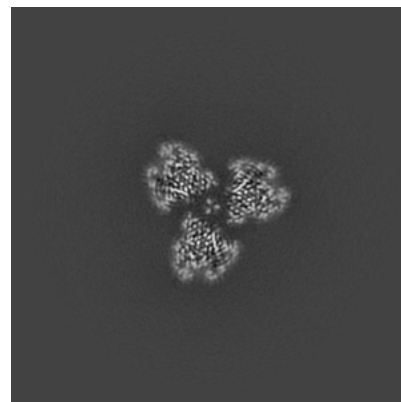
6.3.1 Primary map



X Index: 196

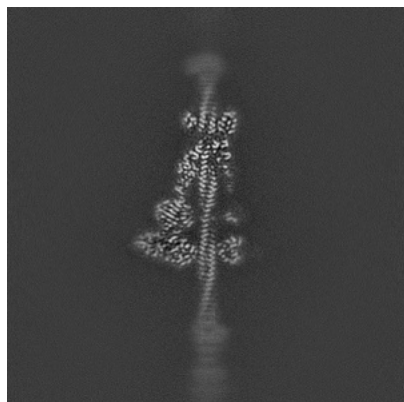


Y Index: 205

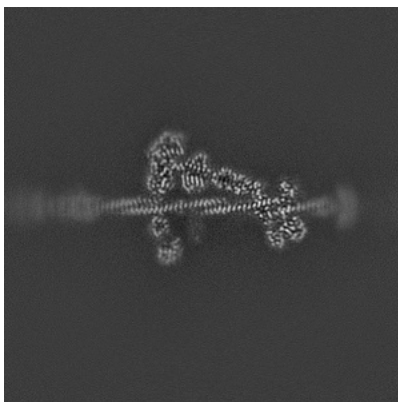


Z Index: 165

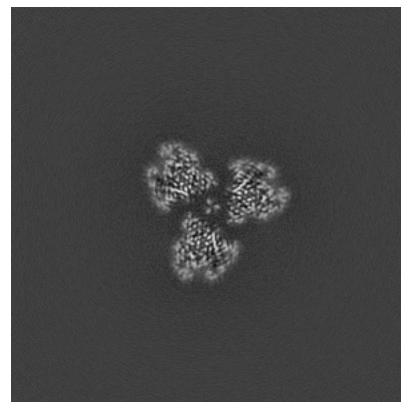
6.3.2 Raw map



X Index: 196



Y Index: 204



Z Index: 165

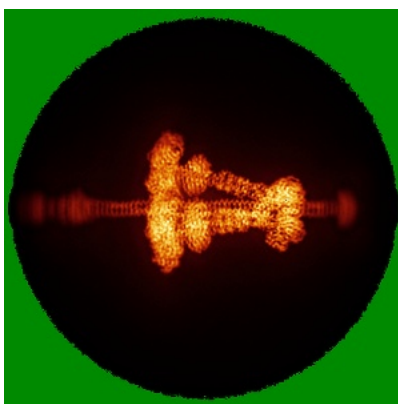
The images above show the largest variance slices of the map in three orthogonal directions.

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

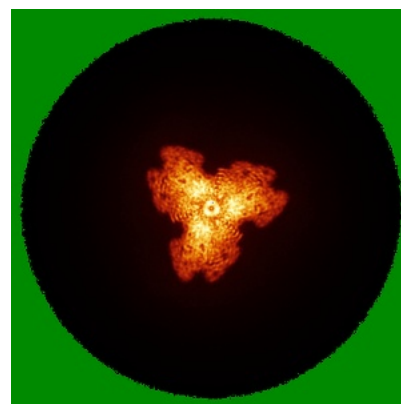
6.4.1 Primary map



X



Y

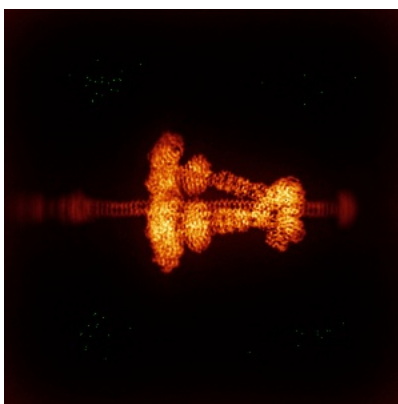


Z

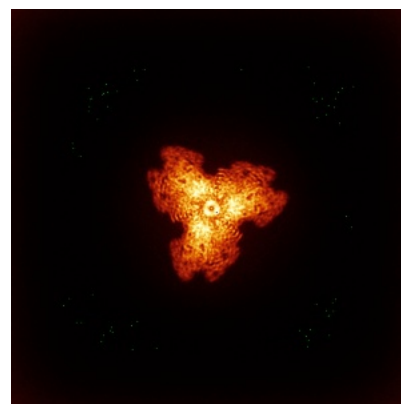
6.4.2 Raw map



X



Y

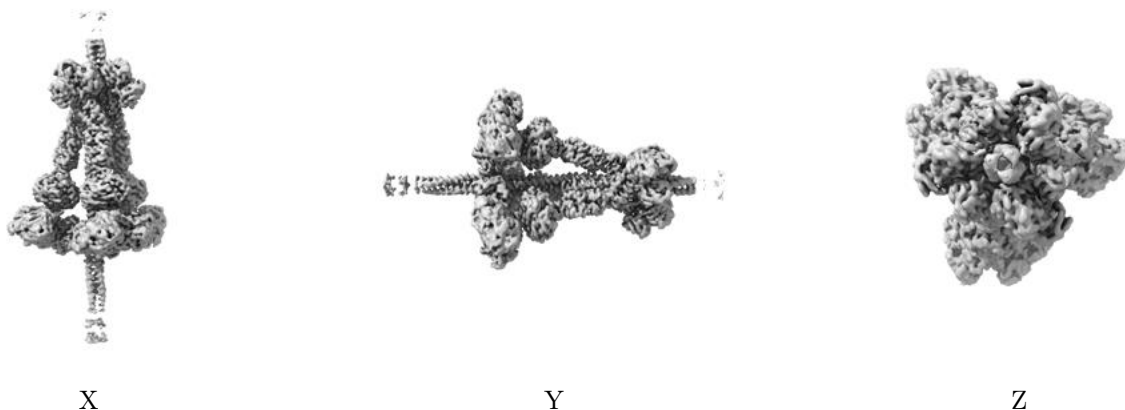


Z

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.

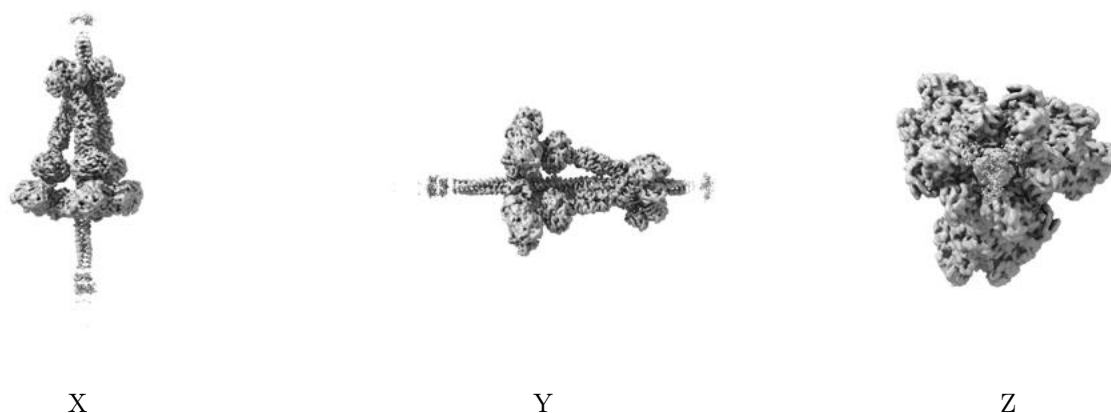
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.12. 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.

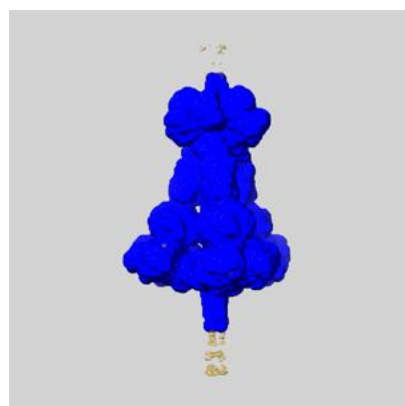
6.6 Mask visualisation [i](#)

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

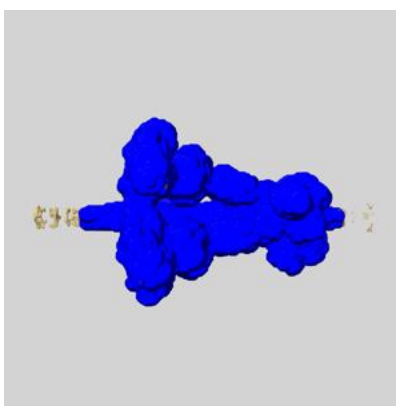
A mask typically either:

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

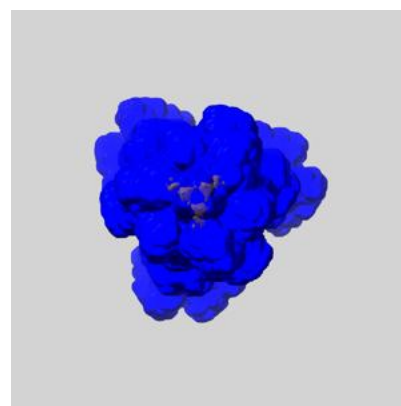
6.6.1 emd_72915_msk_1.map [i](#)



X



Y

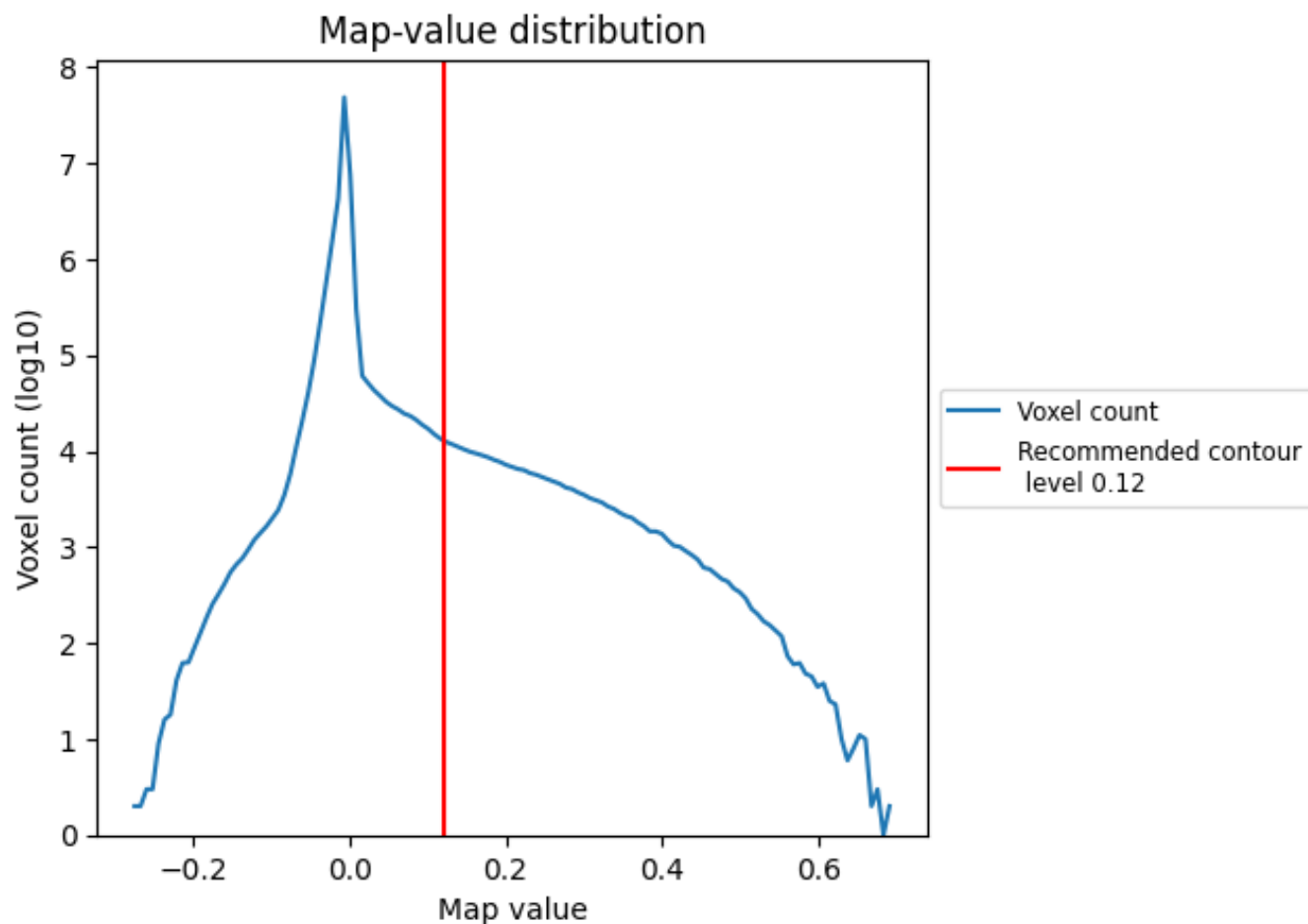


Z

7 Map analysis [i](#)

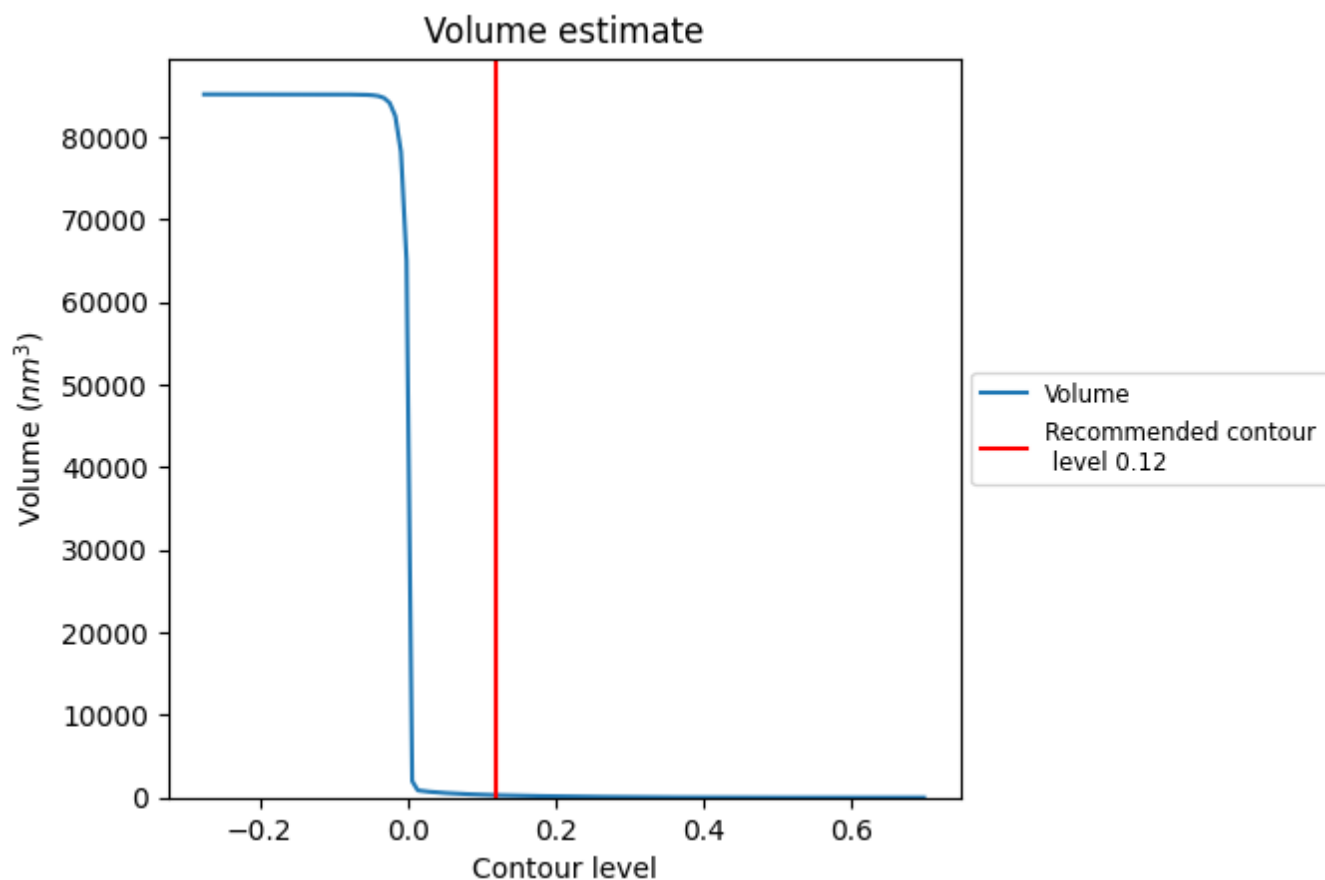
This section contains the results of statistical analysis of the map.

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.

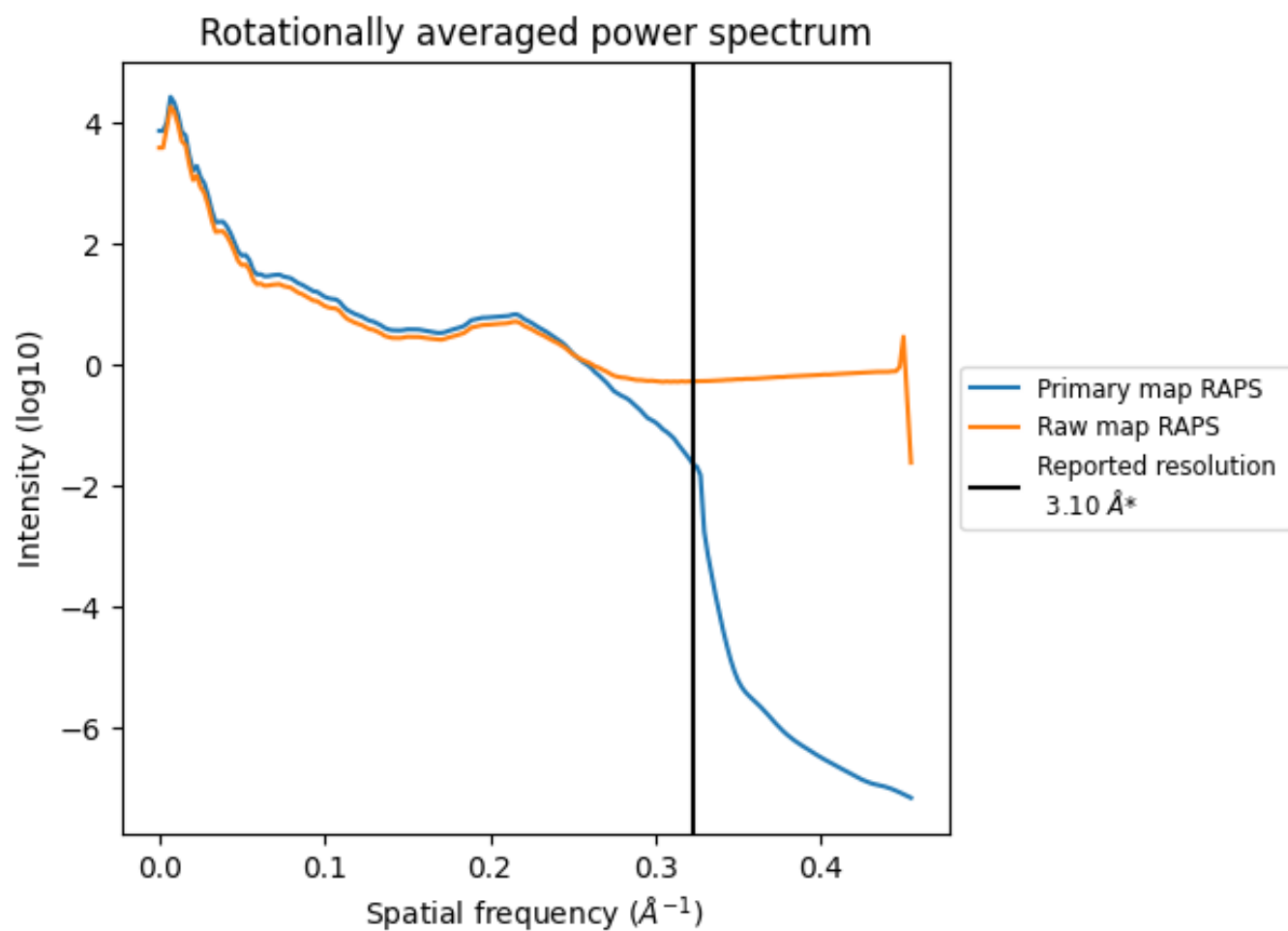
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 291 nm^3 ; this corresponds to an approximate mass of 263 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 ⓘ

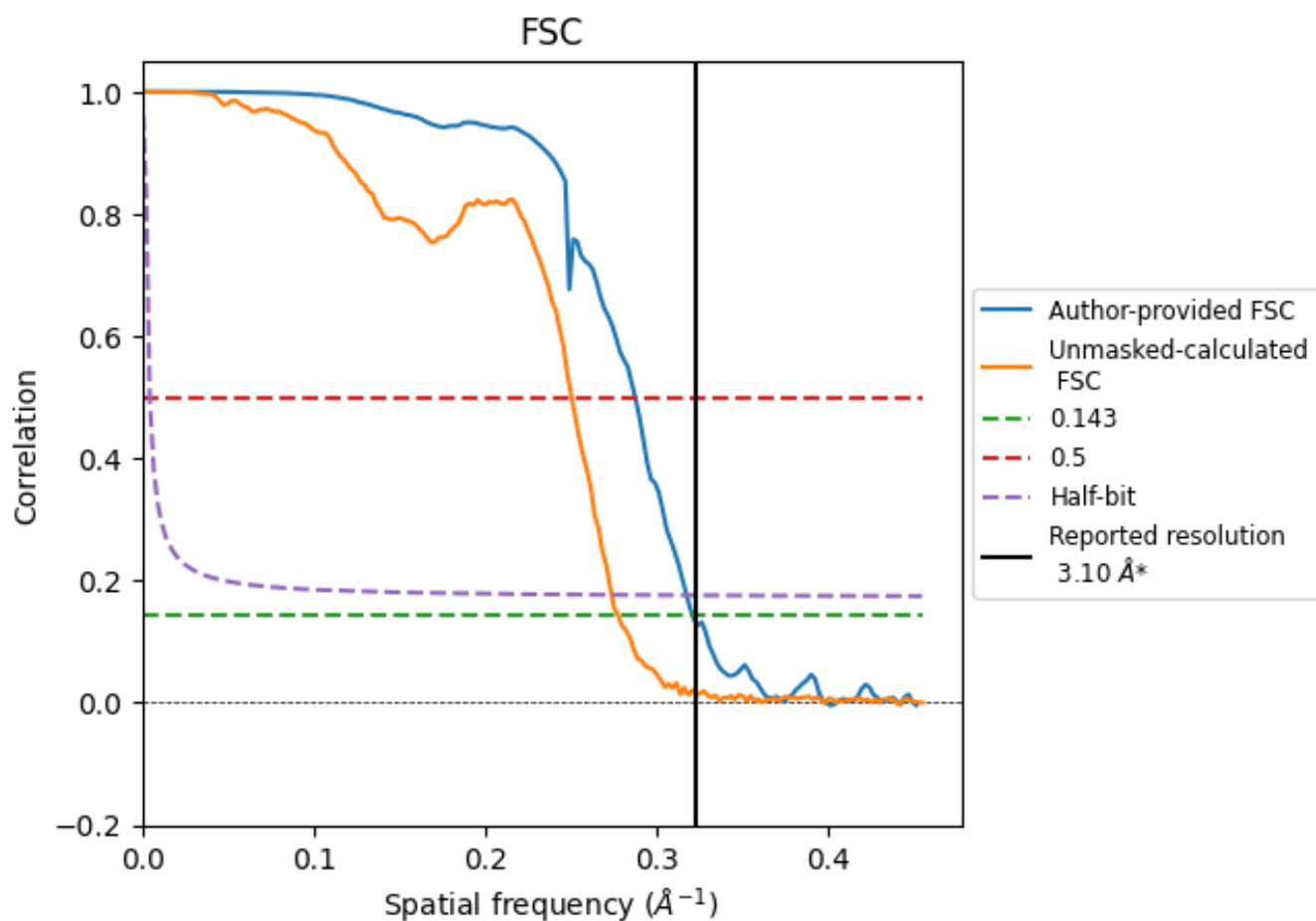


*Reported resolution corresponds to spatial frequency of 0.323 Å⁻¹

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 0.323 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

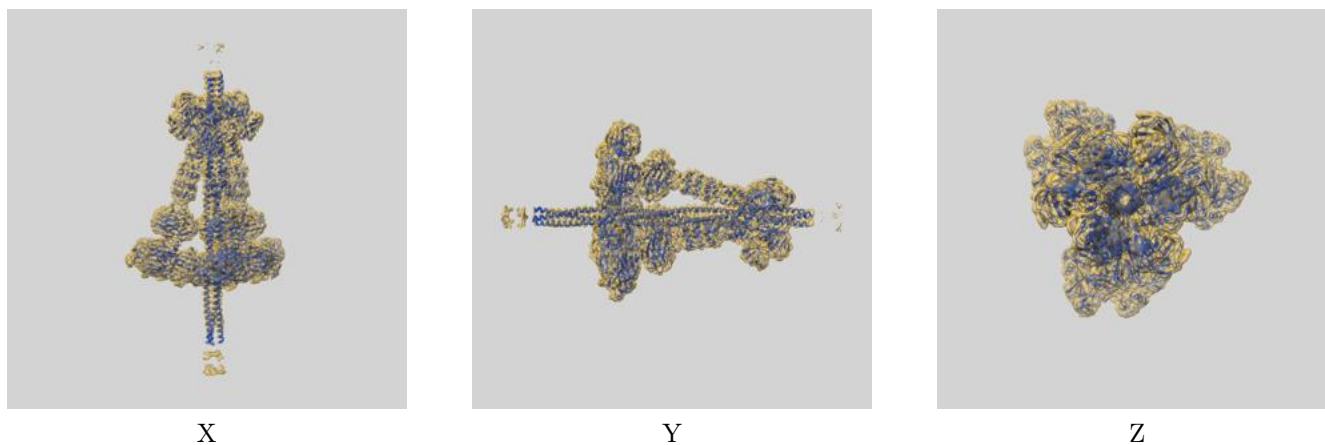
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	3.12	3.48	3.15
Unmasked-calculated*	3.61	4.00	3.65

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

9 Map-model fit [i](#)

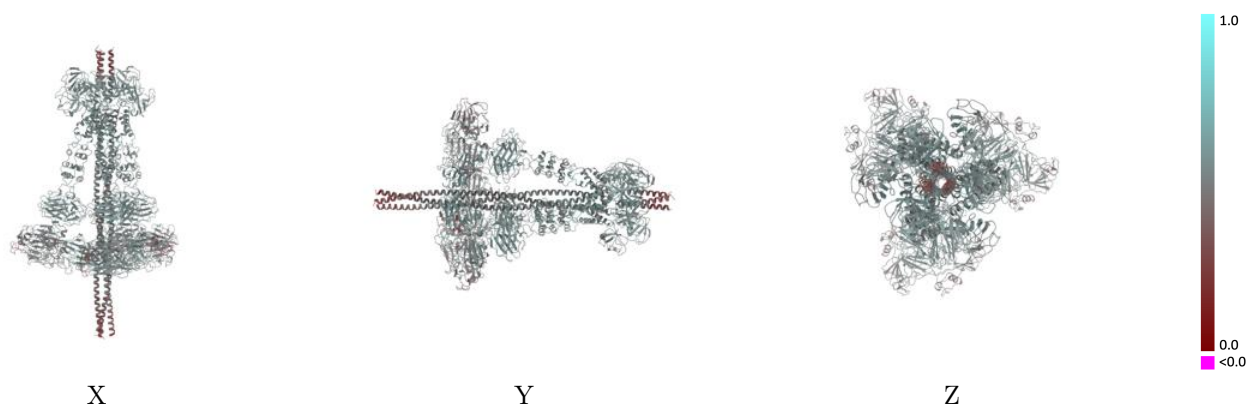
This section contains information regarding the fit between EMDB map EMD-72915 and PDB model 9YG8. Per-residue inclusion information can be found in section [3](#) on page [10](#).

9.1 Map-model overlay [i](#)



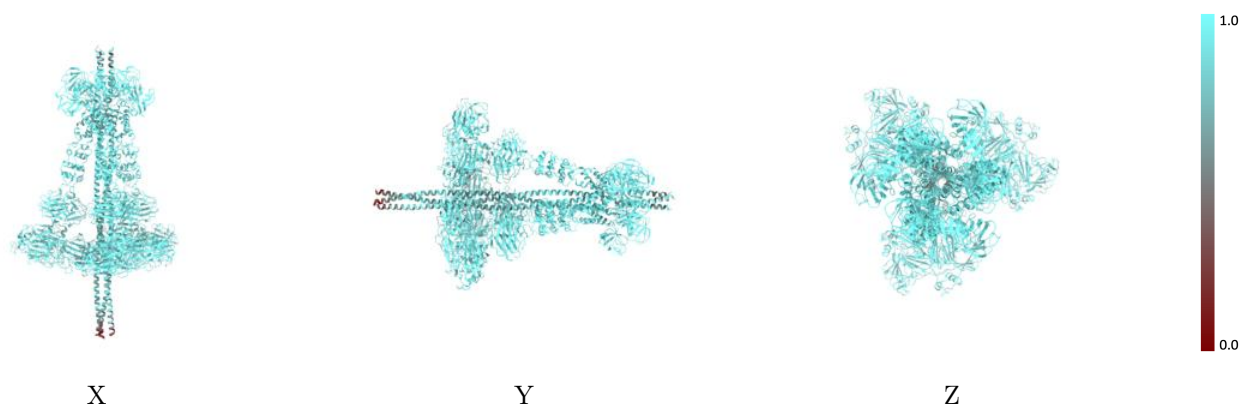
The images above show the 3D surface view of the map at the recommended contour level 0.12 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



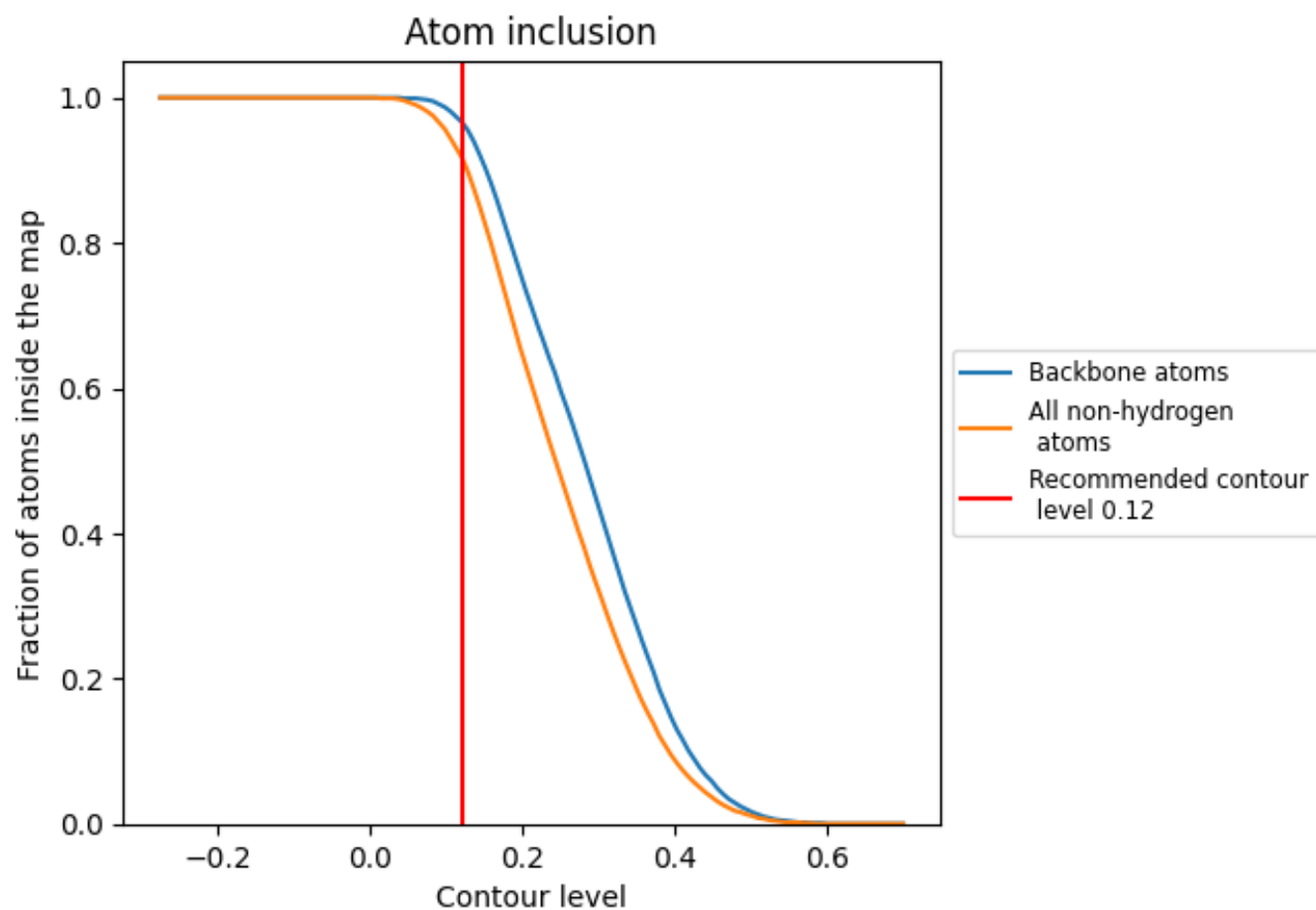
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions at which atoms can be resolved. Low Q-score values may therefore be expected for many entries.

9.3 Atom inclusion mapped to coordinate model [i](#)



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.12).

9.4 Atom inclusion [i](#)



At the recommended contour level, 97% of all backbone atoms, 92% 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 (0.12) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9180	0.5180
A	0.8130	0.4620
B	0.9280	0.5180
C	0.9280	0.5130
D	0.9370	0.5430
E	0.8080	0.4600
F	0.9310	0.5180
G	0.9240	0.5130
H	0.9370	0.5410
I	0.8110	0.4620
J	0.9310	0.5180
K	0.9240	0.5140
L	0.9360	0.5430

