



Full wwPDB EM Validation Report ⓘ

Aug 21, 2025 – 11:15 AM EDT

PDB ID : 9E01 / pdb_00009e01
EMDB ID : EMD-47345
Title : Cryo-EM structure of a TatBC-MdoD complex from Escherichia coli
Authors : Deme, J.C.; Bryant, O.J.; Berks, B.C.; Lea, S.M.
Deposited on : 2024-10-17
Resolution : 2.40 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.dev126
MolProbity : 4-5-2 with Phenix2.0rc1
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) : 2.45.1

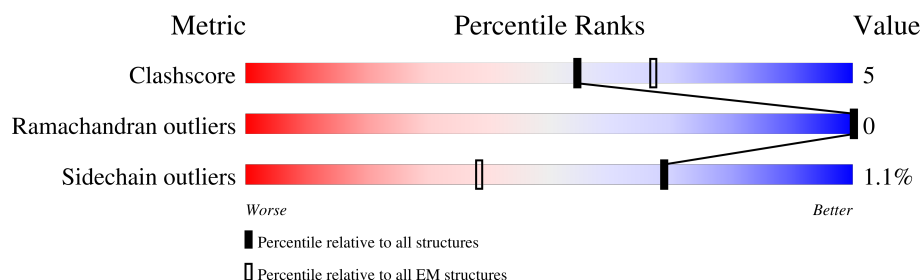
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.40 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.



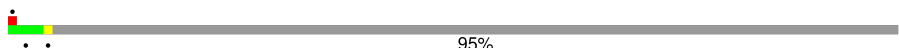
Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	210492	15764
Ramachandran outliers	207382	16835
Sidechain outliers	206894	16415

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 ≥ 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	B	171	<div> <div>5%</div> <div>34%</div> <div>63%</div> </div>
1	D	171	<div> <div>34%</div> <div>63%</div> </div>
1	F	171	<div> <div>6%</div> <div>33%</div> <div>5%</div> <div>63%</div> </div>
2	A	266	<div> <div>78%</div> <div>9%</div> <div>13%</div> </div>
2	C	266	<div> <div>74%</div> <div>13%</div> <div>13%</div> </div>
2	E	266	<div> <div>76%</div> <div>12%</div> <div>13%</div> </div>
3	G	552	<div> <div>95%</div> </div>
3	H	552	<div> <div>95%</div> </div>

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Mol	Chain	Length	Quality of chain
3	I	552	 95%

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 7638 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 Sec-independent protein translocase protein TatB.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	B	64	Total	C	N	O	S	0	0
			504	331	83	89	1		
1	D	64	Total	C	N	O	S	0	0
			504	331	83	89	1		
1	F	64	Total	C	N	O	S	0	0
			504	331	83	89	1		

- Molecule 2 is a protein called Sec-independent protein translocase protein TatC.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	A	232	Total	C	N	O	S	2	0
			1845	1252	279	302	12		
2	C	232	Total	C	N	O	S	2	0
			1845	1252	279	302	12		
2	E	232	Total	C	N	O	S	2	0
			1845	1252	279	302	12		

There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	259	ARG	-	expression tag	UNP C3SK12
A	260	SER	-	expression tag	UNP C3SK12
A	261	HIS	-	expression tag	UNP C3SK12
A	262	HIS	-	expression tag	UNP C3SK12
A	263	HIS	-	expression tag	UNP C3SK12
A	264	HIS	-	expression tag	UNP C3SK12
A	265	HIS	-	expression tag	UNP C3SK12
A	266	HIS	-	expression tag	UNP C3SK12
C	259	ARG	-	expression tag	UNP C3SK12
C	260	SER	-	expression tag	UNP C3SK12
C	261	HIS	-	expression tag	UNP C3SK12
C	262	HIS	-	expression tag	UNP C3SK12
C	263	HIS	-	expression tag	UNP C3SK12

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Chain	Residue	Modelled	Actual	Comment	Reference
C	264	HIS	-	expression tag	UNP C3SK12
C	265	HIS	-	expression tag	UNP C3SK12
C	266	HIS	-	expression tag	UNP C3SK12
E	259	ARG	-	expression tag	UNP C3SK12
E	260	SER	-	expression tag	UNP C3SK12
E	261	HIS	-	expression tag	UNP C3SK12
E	262	HIS	-	expression tag	UNP C3SK12
E	263	HIS	-	expression tag	UNP C3SK12
E	264	HIS	-	expression tag	UNP C3SK12
E	265	HIS	-	expression tag	UNP C3SK12
E	266	HIS	-	expression tag	UNP C3SK12

- Molecule 3 is a protein called Glucans biosynthesis protein D.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	G	27	Total	C	N	O	S	0	0
			196	121	37	34	4		
3	H	27	Total	C	N	O	S	0	0
			196	121	37	34	4		
3	I	27	Total	C	N	O	S	0	0
			196	121	37	34	4		

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
G	0	SER	-	expression tag	UNP A0A138L4Y6
H	0	SER	-	expression tag	UNP A0A138L4Y6
I	0	SER	-	expression tag	UNP A0A138L4Y6

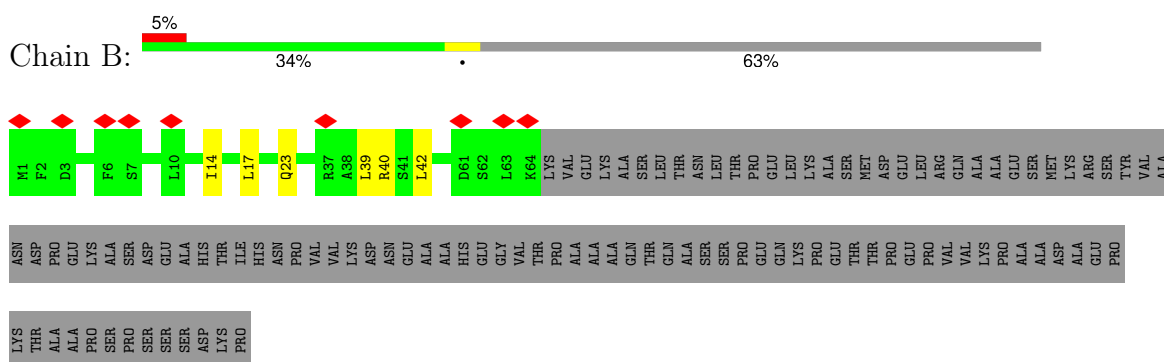
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		AltConf
4	A	1	Total	O	0
			1	1	
4	C	1	Total	O	0
			1	1	
4	E	1	Total	O	0
			1	1	

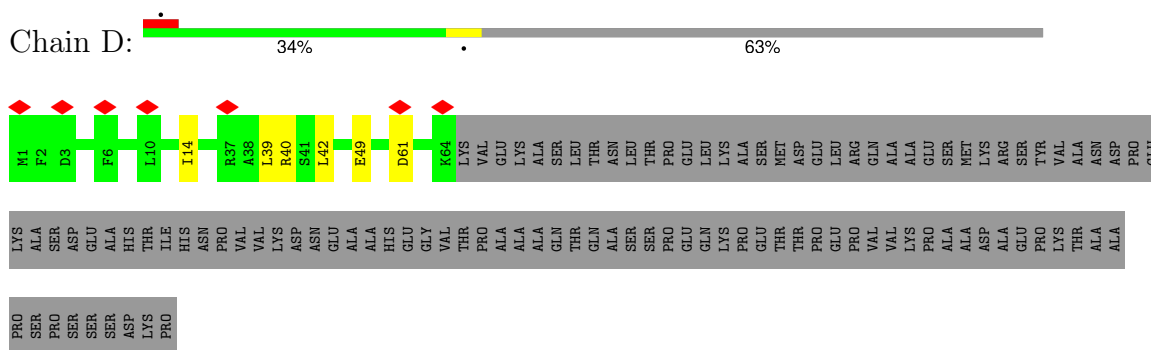
3 Residue-property plots [i](#)

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.

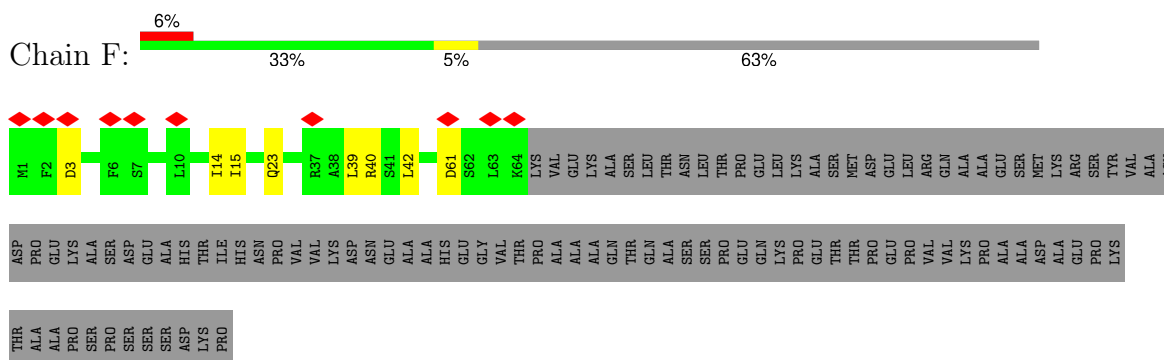
- Molecule 1: Sec-independent protein translocase protein TatB



- Molecule 1: Sec-independent protein translocase protein TatB



- Molecule 1: Sec-independent protein translocase protein TatB



- [illegible]

[illegible]

- Molecule 3: Glucans biosynthesis protein D

Chain H: 95%

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

- Molecule 3: Glucans biosynthesis protein D

Chain I:  95%

ALA	THR	THR	THR	PRO	GLN	ALA	ALA	THR	ASN	ILE	GLN	TYR	ASP	ALA	GLY	LYS	LEU	TRP	HIS	ASN	VAL	GLU	ASN	ARG	GLN	LEU	ASP	GLN	THR	GLN	ARG	PHE	PHE	HIS	MET	GLY	MET	GLY	PHE	ARG	ARG	VAL	ARG	MET	PHE	SER	VAL	PRO	ALA	THR	HIS	LEU	ALA	GLU	ILE	HIS	THR	PHE
SER	M1	R5	F6	I7	K8			M11		I22	A23		S27	GLN	ALA	ALA	PHE	ALA	ASP	SER	ASP	ILE	ALA	ASP	GLY	GLN	THR	GLN	ARG	PHE	PHE	ASP	PHE	SER	ILE	LEU	GLN	SER	MET	ALA	HIS	ASP	LEU	ALA	GLN	THR	ALA	TRP	ARG	GLY	ALA	PRO	PRO	LEU	PRO	ASP	ILE	THR



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	1201445	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 ($e^-/\text{\AA}^2$)	52.1	Depositor
Minimum defocus (nm)	500	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	1.965	Depositor
Minimum map value	-0.043	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.012	Depositor
Recommended contour level	0.1	Depositor
Map size (Å)	354.816, 354.816, 354.816	wwPDB
Map dimensions	512, 512, 512	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.693, 0.693, 0.693	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	B	0.07	0/511	0.15	0/692
1	D	0.07	0/511	0.14	0/692
1	F	0.08	0/511	0.16	0/692
2	A	0.10	0/1906	0.23	0/2606
2	C	0.11	0/1906	0.25	0/2606
2	E	0.10	0/1906	0.24	0/2606
3	G	0.08	0/197	0.21	0/259
3	H	0.09	0/197	0.24	0/259
3	I	0.09	0/197	0.22	0/259
All	All	0.10	0/7842	0.22	0/10671

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	B	504	0	541	5	0
1	D	504	0	541	5	0
1	F	504	0	541	9	0
2	A	1845	0	1927	15	0
2	C	1845	0	1927	22	0
2	E	1845	0	1927	23	0
3	G	196	0	206	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	H	196	0	206	5	0
3	I	196	0	206	6	0
4	A	1	0	0	0	0
4	C	1	0	0	0	0
4	E	1	0	0	0	0
All	All	7638	0	8022	77	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 (77) 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:A:63:ASP:HB2	2:A:66:SER:HB2	1.67	0.77
2:C:63:ASP:HB2	2:C:66:SER:HB2	1.68	0.76
2:E:63:ASP:HB2	2:E:66:SER:HB2	1.70	0.72
2:E:225:LEU:HD13	2:E:228:ILE:HD12	1.76	0.67
2:C:74:LEU:HA	2:C:77:MET:HE2	1.79	0.65
2:C:65:ALA:HB1	2:C:69:PHE:HD1	1.63	0.63
2:E:186:PRO:HG3	2:E:233:SER:HB3	1.79	0.62
2:A:219:ALA:HA	2:A:222:MET:HE2	1.83	0.61
2:E:65:ALA:HB1	2:E:69:PHE:HD1	1.66	0.61
2:C:186:PRO:HG3	2:C:233:SER:HB3	1.82	0.61
2:A:186:PRO:HG3	2:A:233:SER:HB3	1.84	0.60
2:E:6:THR:OG1	3:I:1:MET:N	2.34	0.59
2:A:65:ALA:HB1	2:A:69:PHE:HD2	1.67	0.59
1:F:42:LEU:HD21	3:I:8:LYS:HG2	1.86	0.57
2:E:201:PHE:HE1	1:F:15:ILE:HG13	1.70	0.56
1:B:42:LEU:HD21	3:G:8:LYS:HG2	1.89	0.55
2:C:177:LEU:O	2:C:181:MET:HG2	2.07	0.54
3:G:6:PHE:O	3:G:10:SER:OG	2.22	0.53
1:D:42:LEU:HD21	3:H:8:LYS:HG2	1.91	0.53
2:A:20:LEU:HD21	1:D:14:ILE:HG12	1.91	0.52
2:A:224:CYS:O	2:A:228:ILE:HG13	2.10	0.52
1:D:49:GLU:OE2	3:H:5:ARG:NH1	2.44	0.51
3:G:27:SER:HA	1:D:40:ARG:HH22	1.77	0.50
3:H:6:PHE:O	3:H:10:SER:OG	2.24	0.50
2:E:65:ALA:HB1	2:E:69:PHE:CD1	2.48	0.49
2:A:181:MET:HG3	2:A:183:ILE:HG13	1.95	0.48
1:F:40:ARG:HH22	3:H:27:SER:HA	1.79	0.48
2:E:224:CYS:O	2:E:228:ILE:HG13	2.14	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:E:106:LEU:HD13	2:E:181:MET:HE2	1.96	0.48
3:I:5:ARG:HH11	3:I:5:ARG:HG3	1.78	0.47
3:G:5:ARG:HG3	3:G:5:ARG:HH11	1.78	0.47
2:E:204:GLY:O	2:E:208:THR:OG1	2.23	0.47
1:B:40:ARG:HH22	3:I:27:SER:HA	1.79	0.47
2:C:102:HIS:CE1	2:C:103:GLU:HG3	2.50	0.47
2:A:102:HIS:CE1	2:A:103:GLU:HG3	2.50	0.47
1:B:14:ILE:HG12	2:E:20:LEU:HD21	1.97	0.47
2:C:20:LEU:HD21	1:F:14:ILE:HG12	1.98	0.46
2:C:224:CYS:O	2:C:228:ILE:HG13	2.15	0.46
2:E:40:ASP:N	2:E:40:ASP:OD1	2.48	0.46
2:A:103:GLU:OE1	3:G:4:ARG:NH2	2.49	0.46
2:C:65:ALA:HB1	2:C:69:PHE:CD1	2.47	0.45
2:A:65:ALA:HB1	2:A:69:PHE:CD2	2.49	0.45
2:C:77:MET:HE3	2:C:159:MET:HE3	1.99	0.45
2:C:228:ILE:HG22	2:C:232:PHE:CE2	2.52	0.45
1:B:23:GLN:NE2	1:F:61:ASP:OD2	2.48	0.45
2:C:204:GLY:O	2:C:208:THR:OG1	2.25	0.44
2:E:108:VAL:HB	2:E:109:PRO:HD3	2.00	0.44
2:E:201:PHE:CE1	1:F:15:ILE:HG13	2.51	0.43
2:A:40:ASP:N	2:A:40:ASP:OD1	2.50	0.43
2:E:70:THR:HB	2:E:71:PRO:HD3	2.00	0.43
2:A:117:LEU:HD12	2:A:170:GLU:HA	1.99	0.43
2:E:210:PRO:HD3	1:F:3:ASP:CG	2.44	0.43
2:C:40:ASP:OD1	2:C:40:ASP:N	2.51	0.43
2:C:68:PHE:HB2	2:E:140:THR:HG21	2.00	0.43
2:E:106:LEU:HB2	2:E:181:MET:HE1	2.00	0.43
2:A:204:GLY:O	2:A:208:THR:OG1	2.23	0.43
2:C:103:GLU:OE1	3:H:4:ARG:NH2	2.52	0.43
2:E:102:HIS:CE1	2:E:103:GLU:HG3	2.54	0.43
2:A:181:MET:H	2:A:181:MET:HG2	1.58	0.42
2:C:117:LEU:HD11	2:C:173:VAL:HG21	2.01	0.42
1:F:39:LEU:HD23	1:F:39:LEU:HA	1.83	0.42
2:E:50:ILE:HD11	2:E:59:MET:HG3	2.02	0.42
2:C:129:VAL:HG21	2:C:165:PHE:CE1	2.54	0.42
1:B:39:LEU:HD12	1:B:39:LEU:HA	1.86	0.42
2:C:52:GLN:HB3	2:C:135:GLY:HA2	2.00	0.42
2:C:218:LEU:C	2:C:221:PRO:HD2	2.45	0.42
3:I:5:ARG:HG3	3:I:5:ARG:NH1	2.34	0.42
3:I:7:ILE:O	3:I:11:MET:HG3	2.20	0.42
2:C:225:LEU:HD23	2:C:228:ILE:HD12	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
3:G:5:ARG:HG3	3:G:5:ARG:NH1	2.34	0.41
2:A:118:PHE:O	2:A:122:MET:HG3	2.20	0.41
1:D:61:ASP:OD2	1:F:23:GLN:NE2	2.53	0.41
2:E:228:ILE:HG13	2:E:228:ILE:H	1.65	0.41
2:C:86:VAL:O	2:C:90:GLN:HG2	2.20	0.41
2:E:221:PRO:O	2:E:225:LEU:HB2	2.21	0.41
2:C:162:PHE:HD1	2:C:162:PHE:HA	1.81	0.40
2:E:88:LEU:HD23	2:E:88:LEU:HA	1.94	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	B	62/171 (36%)	62 (100%)	0	0	100	100
1	D	62/171 (36%)	62 (100%)	0	0	100	100
1	F	62/171 (36%)	62 (100%)	0	0	100	100
2	A	232/266 (87%)	230 (99%)	2 (1%)	0	100	100
2	C	232/266 (87%)	229 (99%)	3 (1%)	0	100	100
2	E	232/266 (87%)	229 (99%)	3 (1%)	0	100	100
3	G	25/552 (4%)	25 (100%)	0	0	100	100
3	H	25/552 (4%)	25 (100%)	0	0	100	100
3	I	25/552 (4%)	25 (100%)	0	0	100	100
All	All	957/2967 (32%)	949 (99%)	8 (1%)	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	B	56/144 (39%)	55 (98%)	1 (2%)	54	73
1	D	56/144 (39%)	55 (98%)	1 (2%)	54	73
1	F	56/144 (39%)	56 (100%)	0	100	100
2	A	205/233 (88%)	203 (99%)	2 (1%)	73	86
2	C	205/233 (88%)	202 (98%)	3 (2%)	60	77
2	E	205/233 (88%)	203 (99%)	2 (1%)	73	86
3	G	20/463 (4%)	20 (100%)	0	100	100
3	H	20/463 (4%)	20 (100%)	0	100	100
3	I	20/463 (4%)	20 (100%)	0	100	100
All	All	843/2520 (34%)	834 (99%)	9 (1%)	69	84

All (9) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	17	LEU
2	A	158	VAL
2	A	169	PHE
2	C	20	LEU
2	C	64	VAL
2	C	158	VAL
2	E	64	VAL
2	E	158	VAL
1	D	39	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
2	A	52	GLN
2	A	102	HIS
2	C	102	HIS
2	E	102	HIS

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.

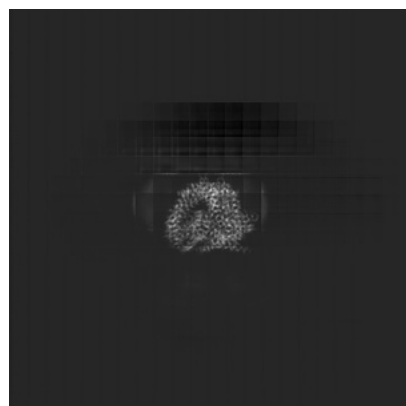
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-47345. These allow visual inspection of the internal detail of the map and identification of artifacts.

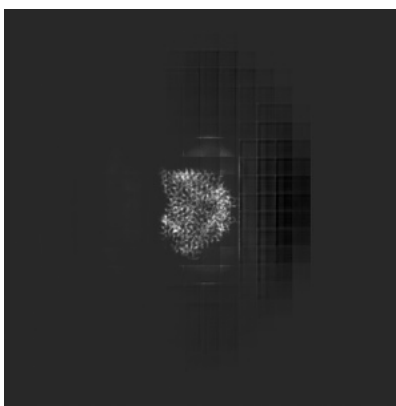
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

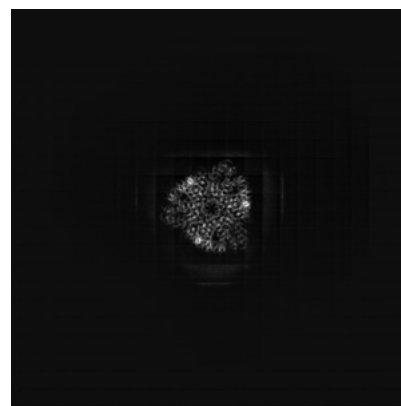
6.1.1 Primary map



X

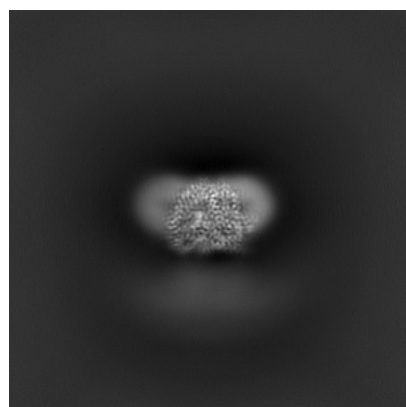


Y

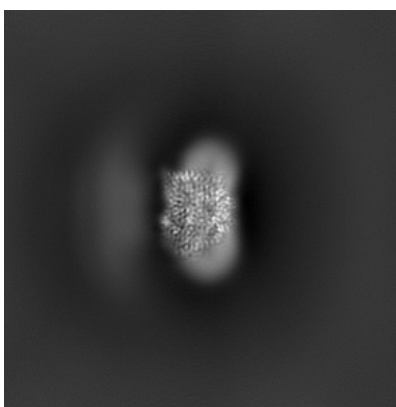


Z

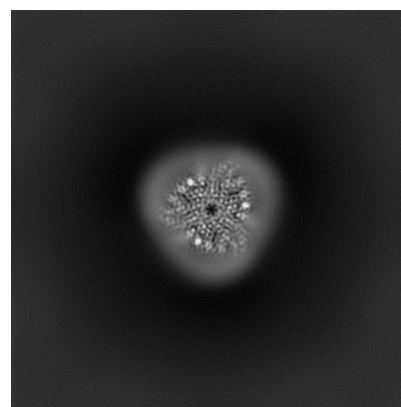
6.1.2 Raw map



X



Y

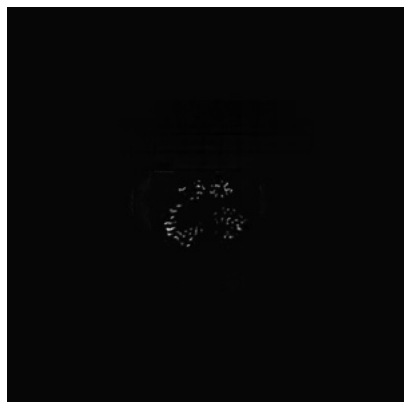


Z

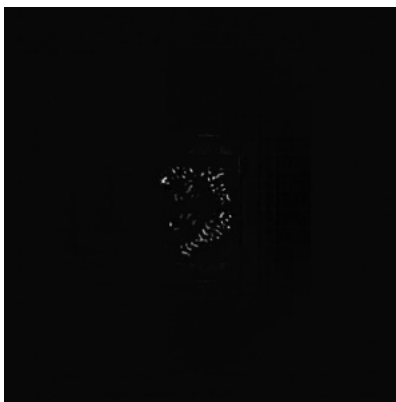
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

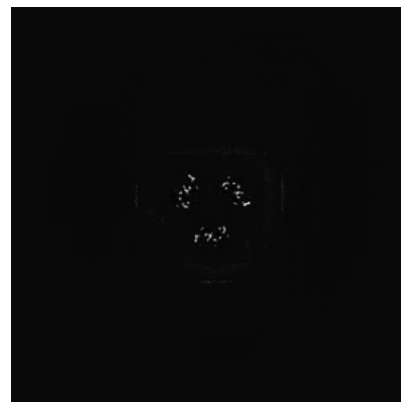
6.2.1 Primary map



X Index: 256

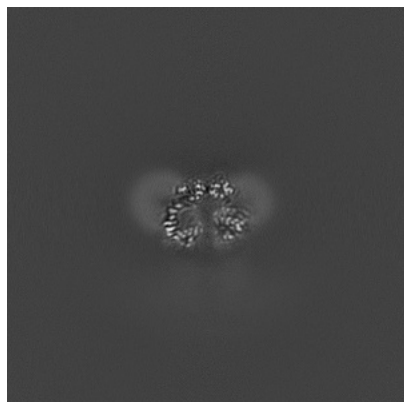


Y Index: 256

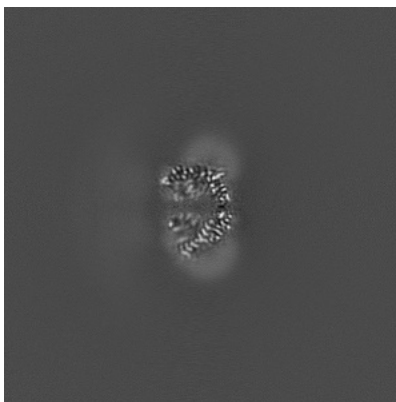


Z Index: 256

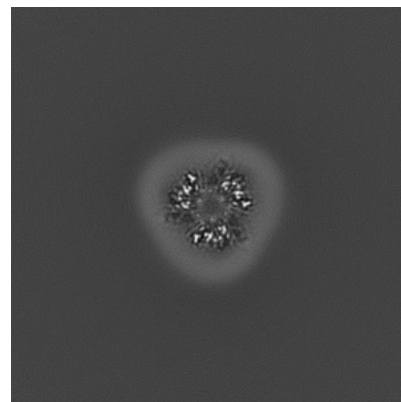
6.2.2 Raw map



X Index: 256



Y Index: 256

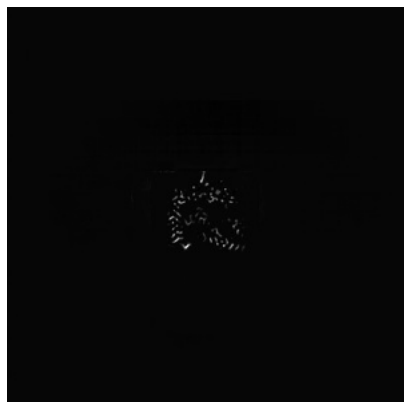


Z Index: 256

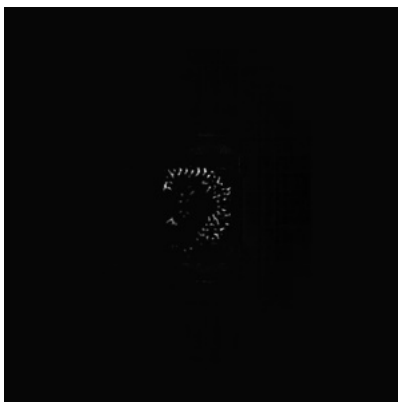
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

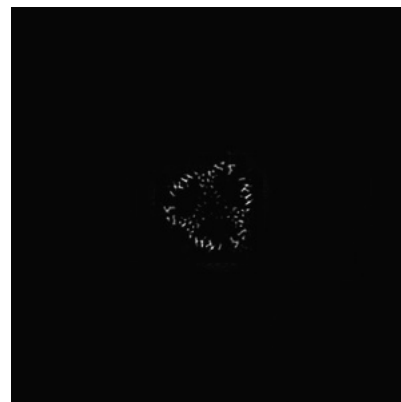
6.3.1 Primary map



X Index: 239

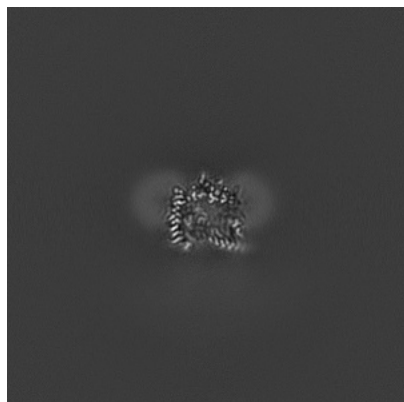


Y Index: 260

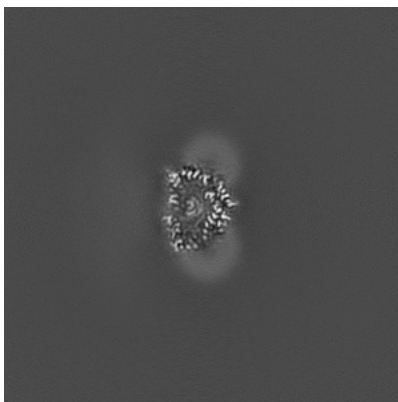


Z Index: 234

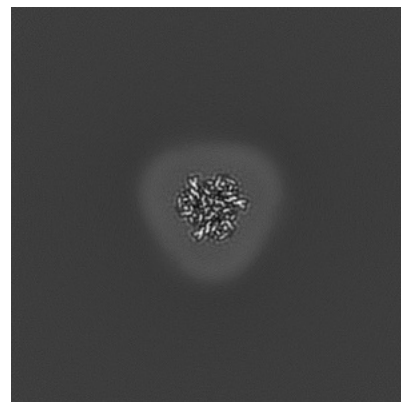
6.3.2 Raw map



X Index: 240



Y Index: 274

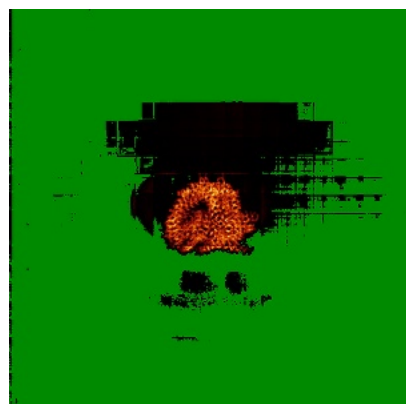


Z Index: 275

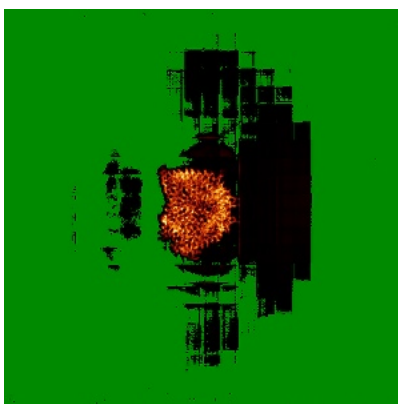
The images above show the largest variance slices of the map in three orthogonal directions.

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

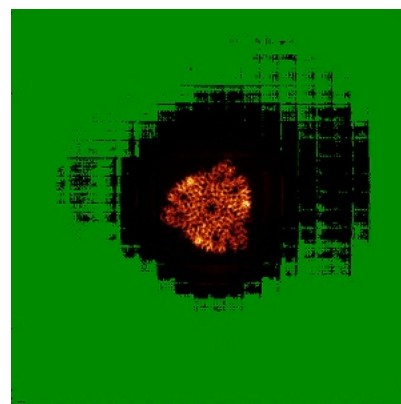
6.4.1 Primary map



X

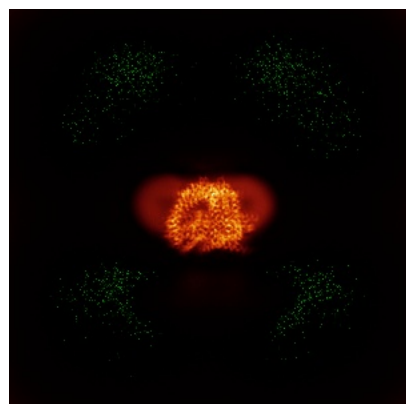


Y

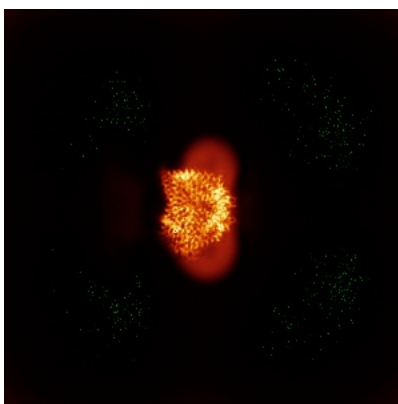


Z

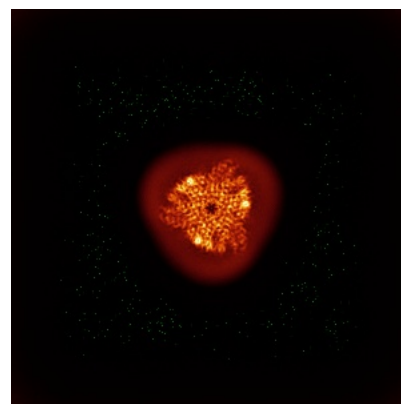
6.4.2 Raw map



X



Y

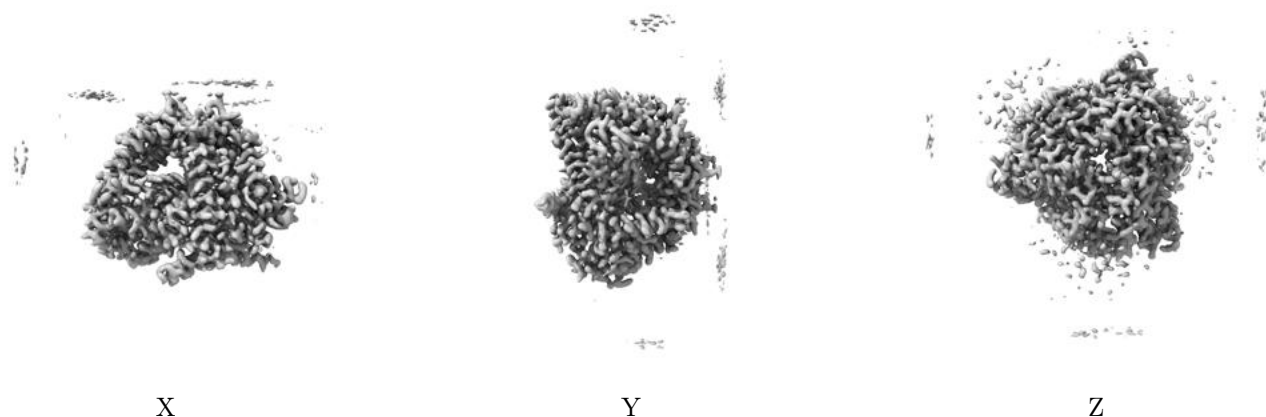


Z

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.

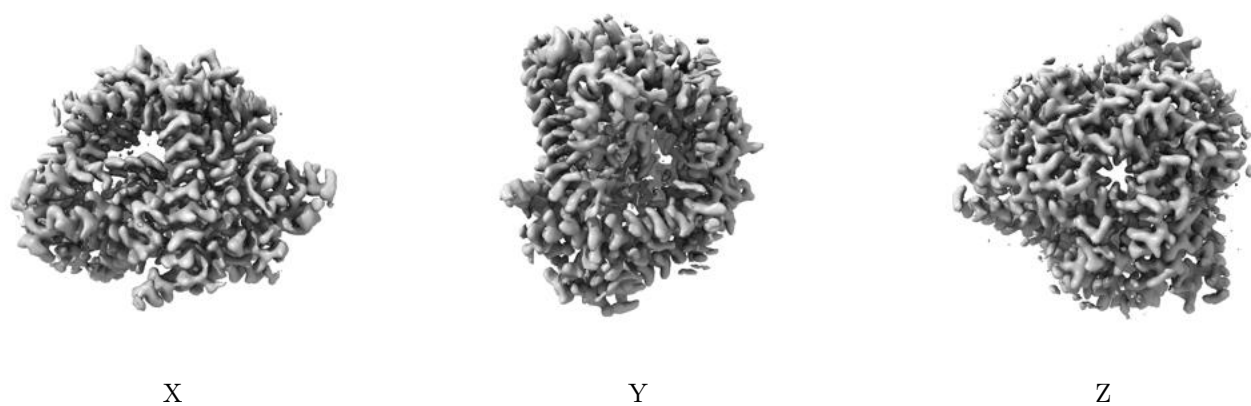
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.1. 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.

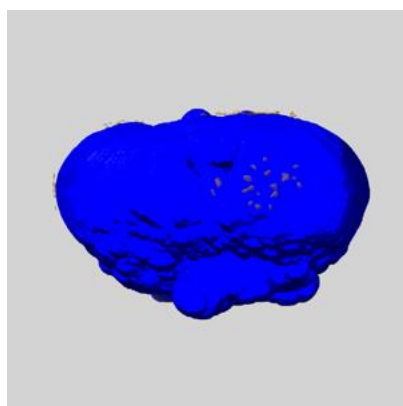
6.6 Mask visualisation [i](#)

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

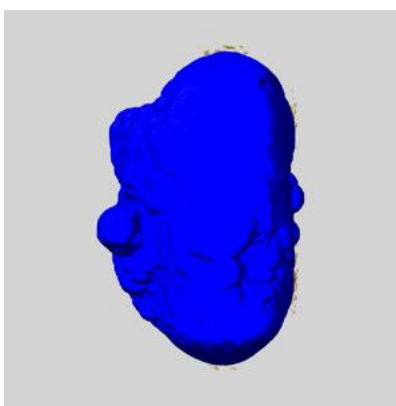
A mask typically either:

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

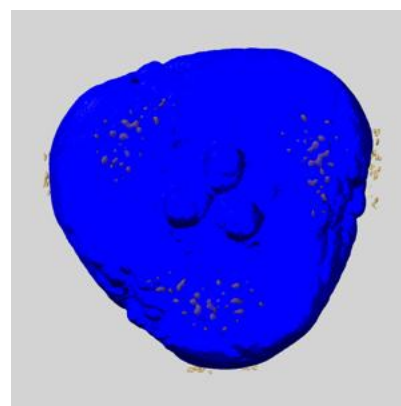
6.6.1 emd_47345_msk_1.map [i](#)



X



Y

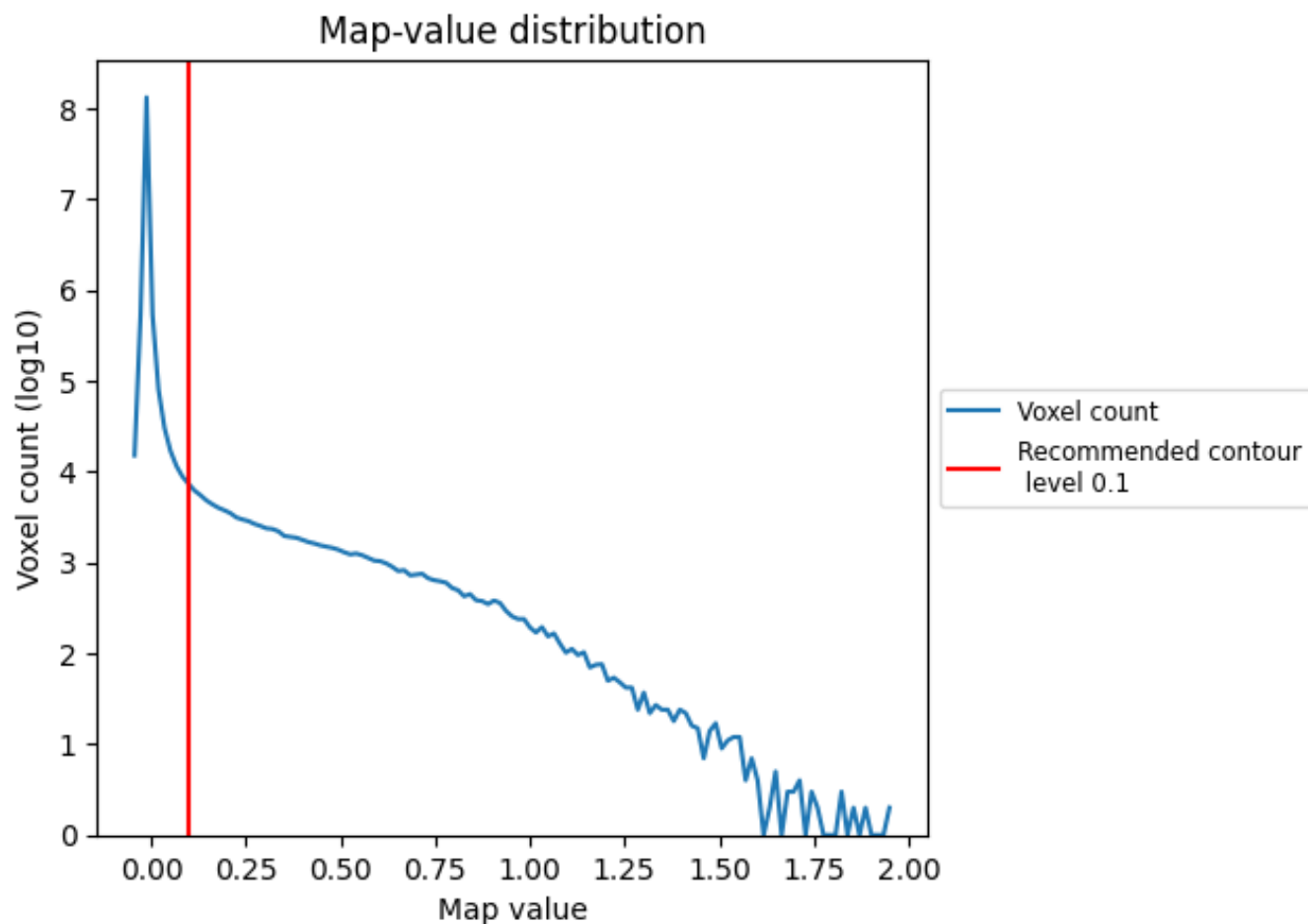


Z

7 Map analysis [i](#)

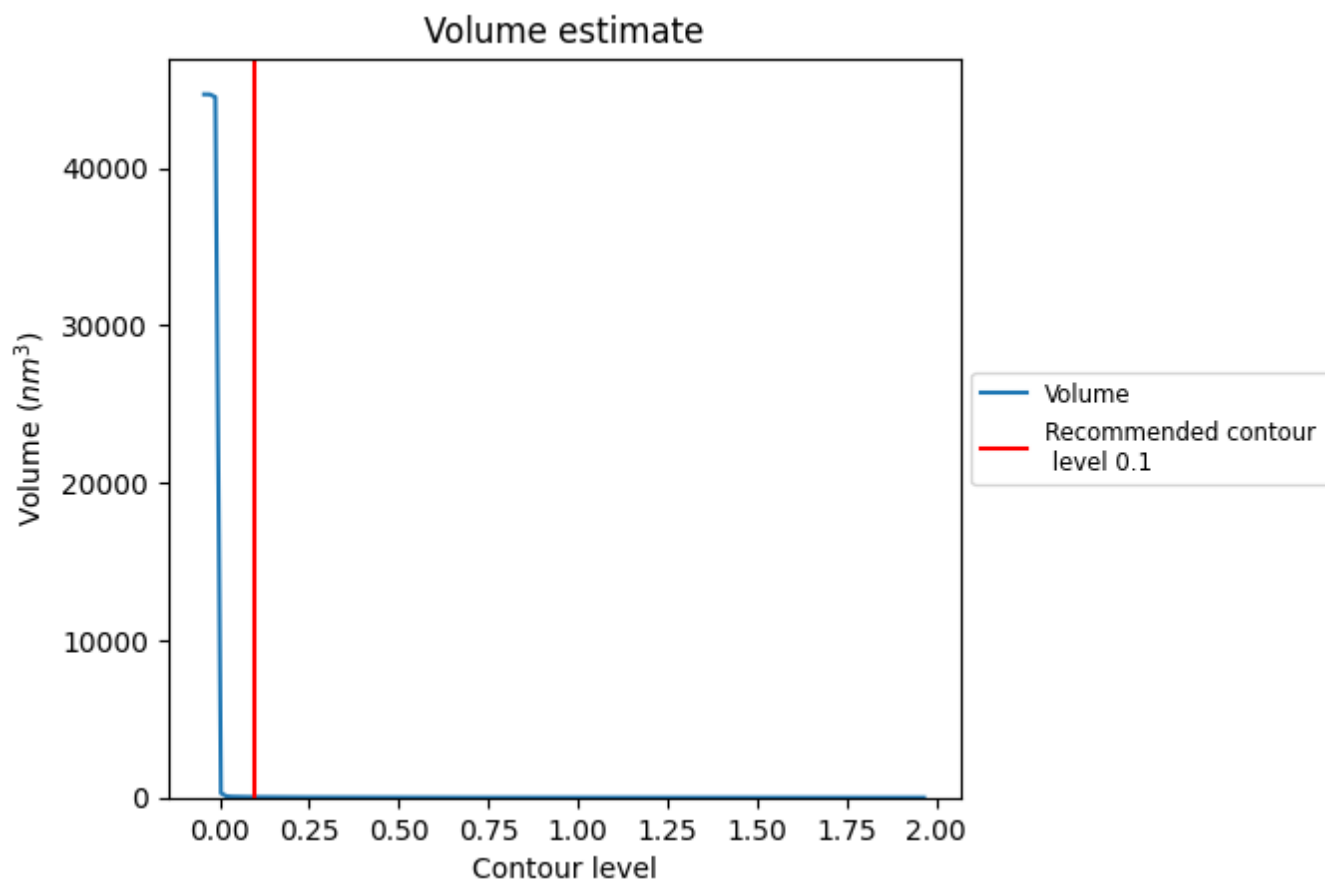
This section contains the results of statistical analysis of the map.

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.

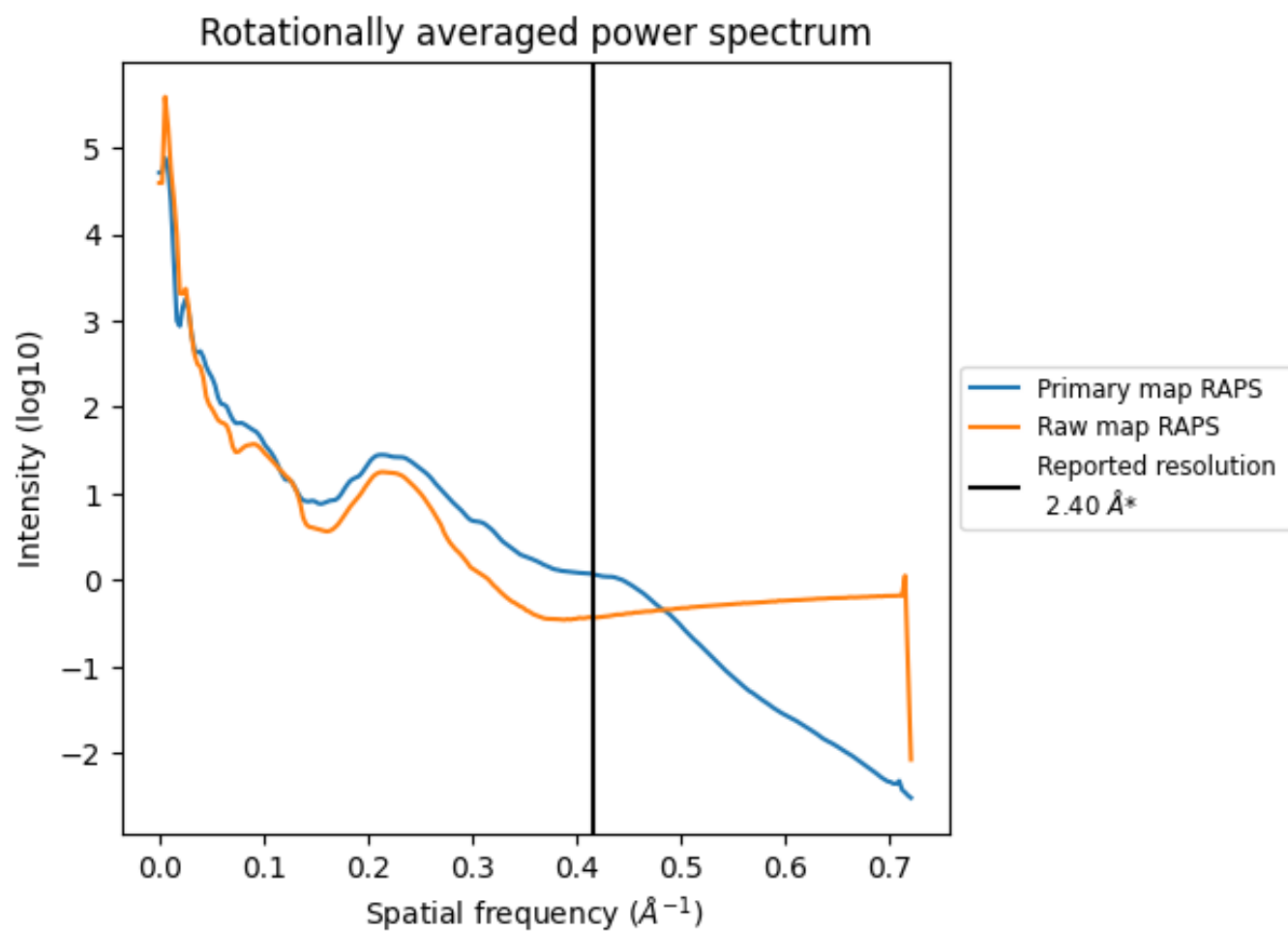
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 33 nm^3 ; this corresponds to an approximate mass of 30 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 ⓘ

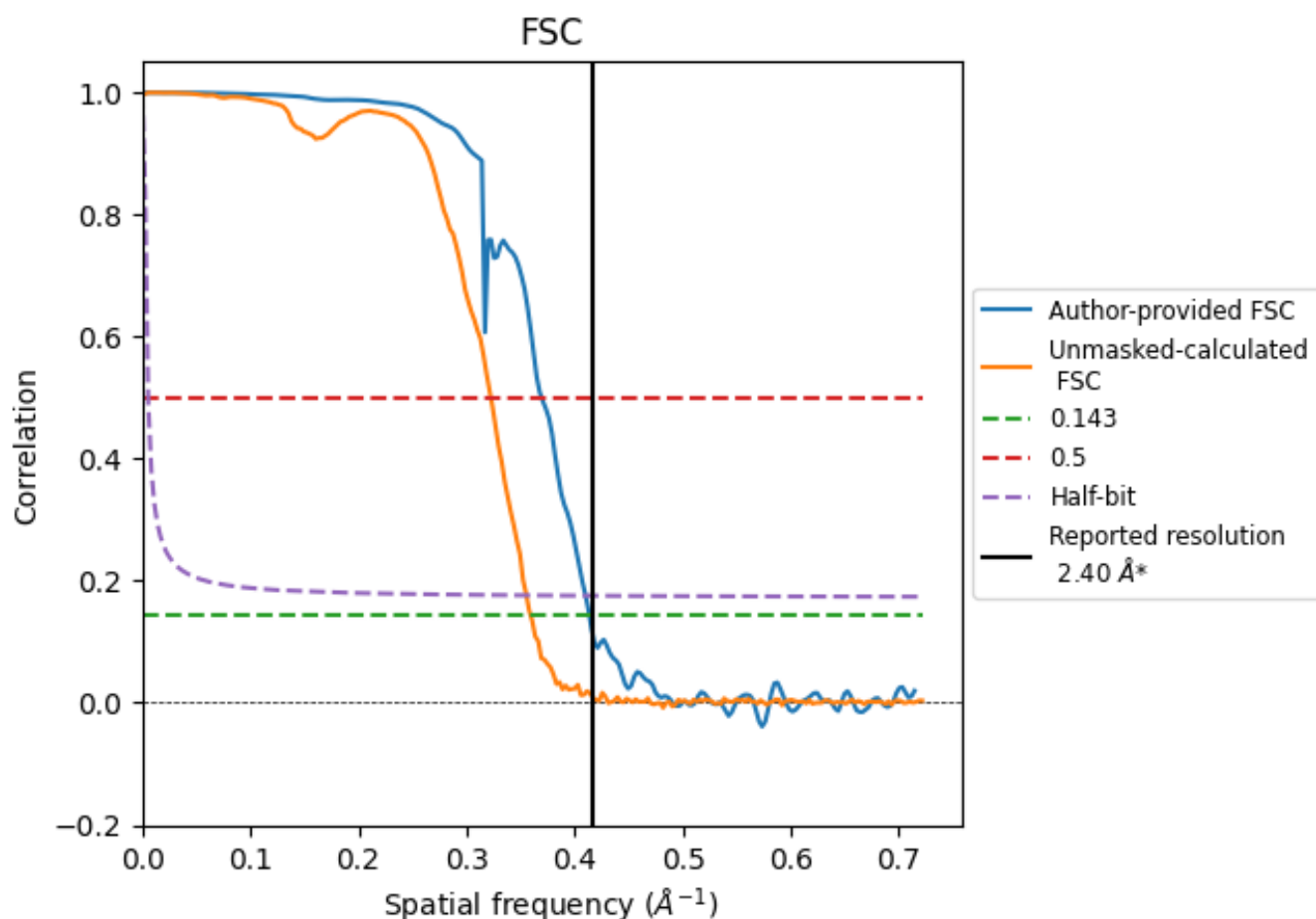


*Reported resolution corresponds to spatial frequency of 0.417 Å⁻¹

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 0.417 \AA^{-1}

8.2 Resolution estimates [i](#)

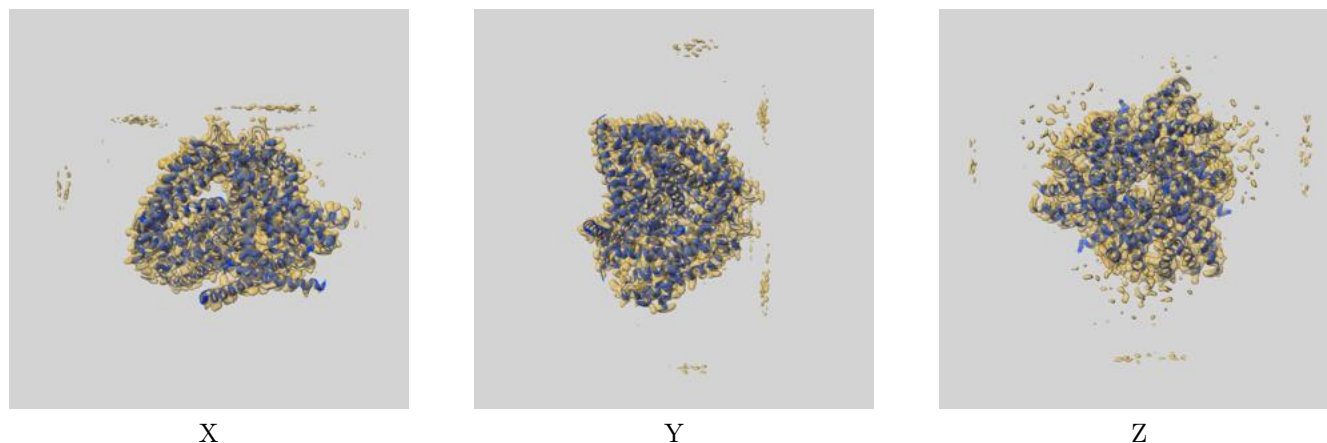
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.40	-	-
Author-provided FSC curve	2.42	2.71	2.44
Unmasked-calculated*	2.78	3.10	2.81

*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 2.78 differs from the reported value 2.4 by more than 10 %

9 Map-model fit [i](#)

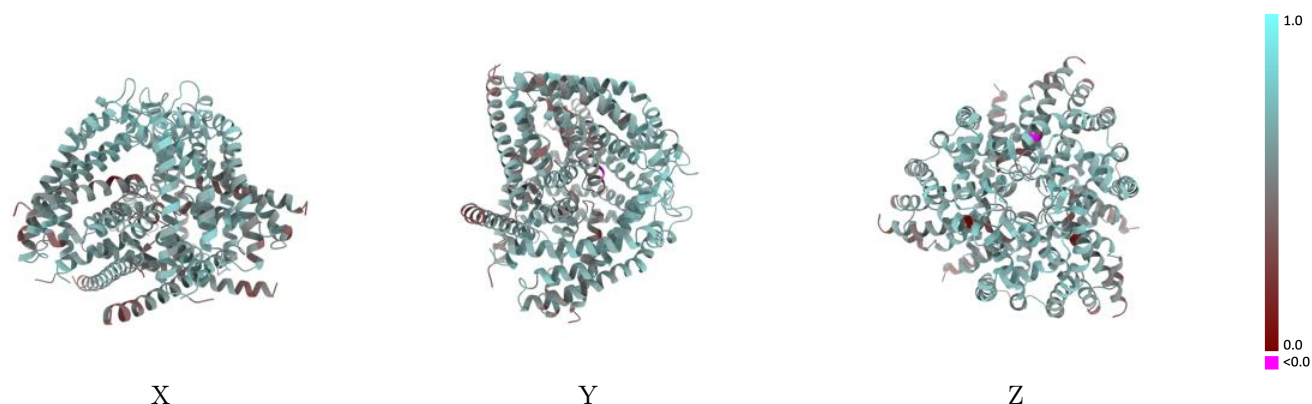
This section contains information regarding the fit between EMDB map EMD-47345 and PDB model 9E01. Per-residue inclusion information can be found in section 3 on page 6.

9.1 Map-model overlay [i](#)



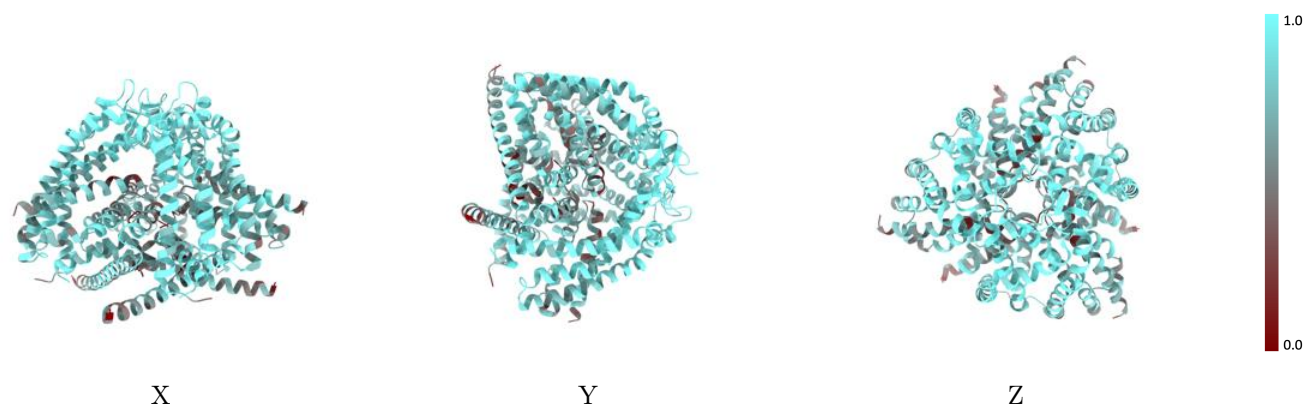
The images above show the 3D surface view of the map at the recommended contour level 0.1 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



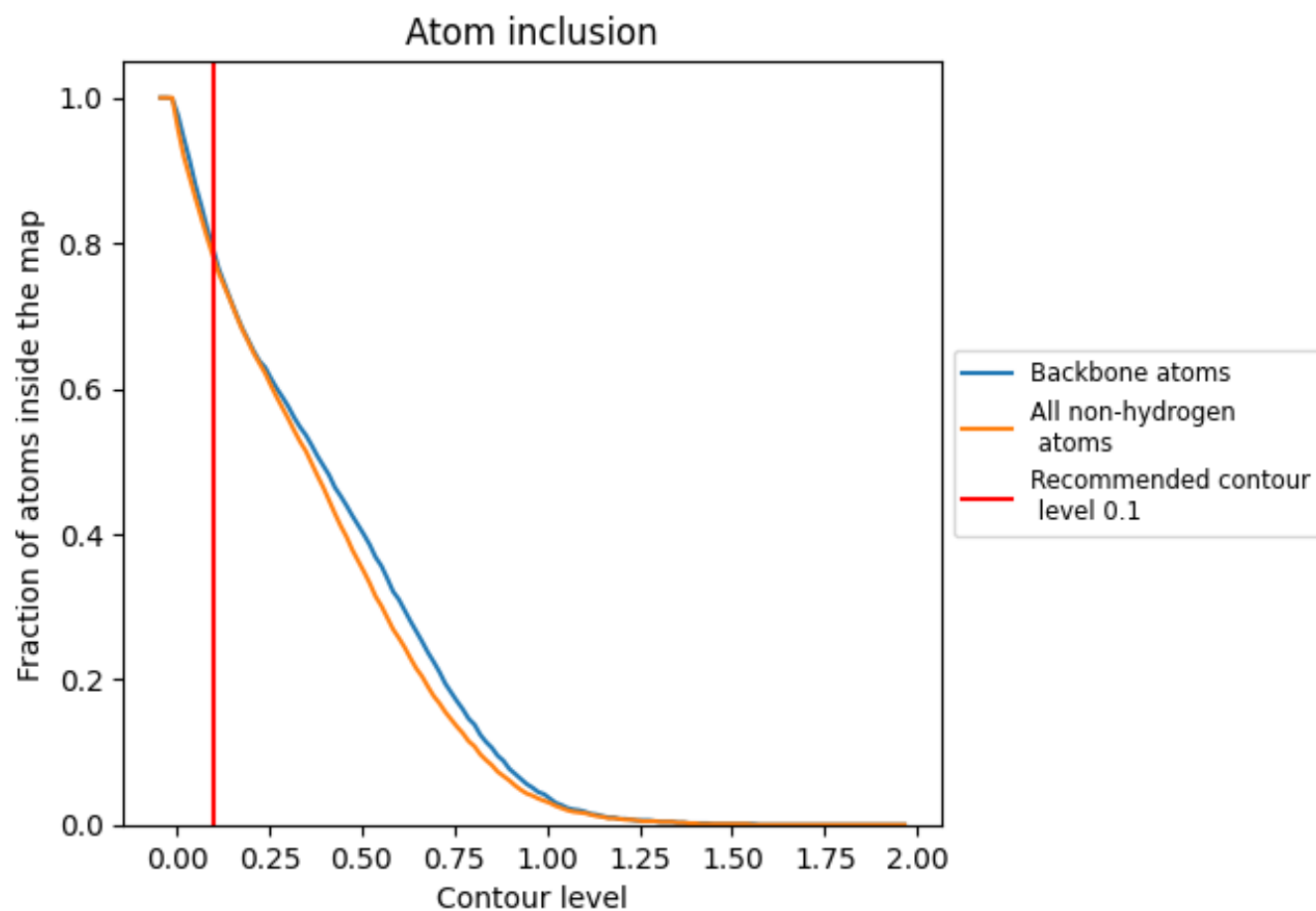
The images above show the model with each residue coloured according its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.1).

9.4 Atom inclusion ⓘ



At the recommended contour level, 79% of all backbone atoms, 78% 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 (0.1) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div>0.7800</div>	<div><div></div>0.5730</div>
A	<div><div></div>0.8260</div>	<div><div></div>0.5900</div>
B	<div><div></div>0.6660</div>	<div><div></div>0.5190</div>
C	<div><div></div>0.8240</div>	<div><div></div>0.5900</div>
D	<div><div></div>0.6740</div>	<div><div></div>0.5140</div>
E	<div><div></div>0.8230</div>	<div><div></div>0.5890</div>
F	<div><div></div>0.6420</div>	<div><div></div>0.5060</div>
G	<div><div></div>0.7110</div>	<div><div></div>0.5790</div>
H	<div><div></div>0.7740</div>	<div><div></div>0.5850</div>
I	<div><div></div>0.7320</div>	<div><div></div>0.5730</div>

1.0

0.0

<0.0