Full wwPDB EM Validation Report (i)

Feb 27, 2024 – 02:02 AM EST

PDB ID	:	6VW2
EMDB ID	:	EMD-21410
Title	:	Cryo-EM structure of human islet amyloid polypeptide (hIAPP, or amylin) fibrils
Authors	:	Cao, Q.; Boyer, D.R.; Sawaya, M.R.; Eisenberg, D.S.
Deposited on	:	2020-02-18
Resolution	:	3.40 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 (i) 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 (i)) were used in the production of this report:

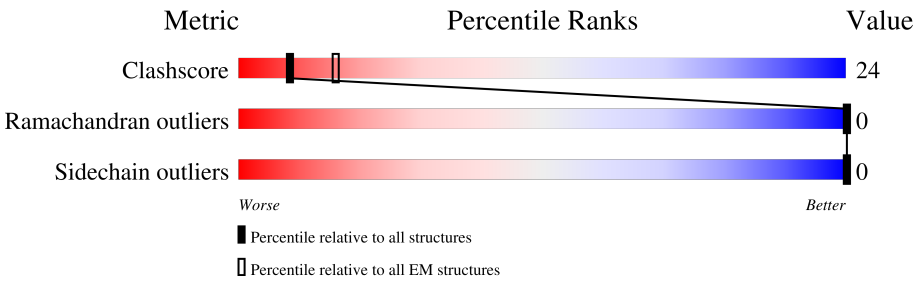
EMDB validation analysis	:	0.0.1.dev70
MolProbity	:	4.02b-467
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ	:	1.9.13
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.40 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	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the EM map (all-atom inclusion < 40%). The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	155	10% 6% 85%
1	B	155	9% 6% 85%
1	C	155	10% 6% 85%
1	D	155	9% 6% 85%
1	E	155	9% 6% 85%
1	F	155	10% 6% 85%
1	G	155	9% 6% 85%
1	H	155	10% 6% 85%

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Mol	Chain	Length	Quality of chain
1	I	155	9% 6% 85%
1	J	155	10% 6% 85%

2 Entry composition (i)

There is only 1 type of molecule in this entry. The entry contains 1800 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 Islet amyloid polypeptide (SUMO-tagged).

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
1	A	24	180	111	31	38	0	0
1	B	24	180	111	31	38	0	0
1	C	24	180	111	31	38	0	0
1	D	24	180	111	31	38	0	0
1	E	24	180	111	31	38	0	0
1	F	24	180	111	31	38	0	0
1	G	24	180	111	31	38	0	0
1	H	24	180	111	31	38	0	0
1	I	24	180	111	31	38	0	0
1	J	24	180	111	31	38	0	0

There are 210 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-117	MET	-	initiating methionine	UNP Q12306
A	-116	GLY	-	expression tag	UNP Q12306
A	-115	SER	-	expression tag	UNP Q12306
A	-114	SER	-	expression tag	UNP Q12306
A	-113	HIS	-	expression tag	UNP Q12306
A	-112	HIS	-	expression tag	UNP Q12306
A	-111	HIS	-	expression tag	UNP Q12306
A	-110	HIS	-	expression tag	UNP Q12306
A	-109	HIS	-	expression tag	UNP Q12306
A	-108	HIS	-	expression tag	UNP Q12306

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Chain	Residue	Modelled	Actual	Comment	Reference
A	-107	GLY	-	expression tag	UNP Q12306
A	-106	SER	-	expression tag	UNP Q12306
A	-105	GLY	-	expression tag	UNP Q12306
A	-104	LEU	-	expression tag	UNP Q12306
A	-103	VAL	-	expression tag	UNP Q12306
A	-102	PRO	-	expression tag	UNP Q12306
A	-101	ARG	-	expression tag	UNP Q12306
A	-100	GLY	-	expression tag	UNP Q12306
A	-99	SER	-	expression tag	UNP Q12306
A	-98	ALA	-	expression tag	UNP Q12306
A	-97	SER	-	expression tag	UNP Q12306
B	-117	MET	-	initiating methionine	UNP Q12306
B	-116	GLY	-	expression tag	UNP Q12306
B	-115	SER	-	expression tag	UNP Q12306
B	-114	SER	-	expression tag	UNP Q12306
B	-113	HIS	-	expression tag	UNP Q12306
B	-112	HIS	-	expression tag	UNP Q12306
B	-111	HIS	-	expression tag	UNP Q12306
B	-110	HIS	-	expression tag	UNP Q12306
B	-109	HIS	-	expression tag	UNP Q12306
B	-108	HIS	-	expression tag	UNP Q12306
B	-107	GLY	-	expression tag	UNP Q12306
B	-106	SER	-	expression tag	UNP Q12306
B	-105	GLY	-	expression tag	UNP Q12306
B	-104	LEU	-	expression tag	UNP Q12306
B	-103	VAL	-	expression tag	UNP Q12306
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B	-100	GLY	-	expression tag	UNP Q12306
B	-99	SER	-	expression tag	UNP Q12306
B	-98	ALA	-	expression tag	UNP Q12306
B	-97	SER	-	expression tag	UNP Q12306
C	-117	MET	-	initiating methionine	UNP Q12306
C	-116	GLY	-	expression tag	UNP Q12306
C	-115	SER	-	expression tag	UNP Q12306
C	-114	SER	-	expression tag	UNP Q12306
C	-113	HIS	-	expression tag	UNP Q12306
C	-112	HIS	-	expression tag	UNP Q12306
C	-111	HIS	-	expression tag	UNP Q12306
C	-110	HIS	-	expression tag	UNP Q12306
C	-109	HIS	-	expression tag	UNP Q12306
C	-108	HIS	-	expression tag	UNP Q12306

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-107	GLY	-	expression tag	UNP Q12306
C	-106	SER	-	expression tag	UNP Q12306
C	-105	GLY	-	expression tag	UNP Q12306
C	-104	LEU	-	expression tag	UNP Q12306
C	-103	VAL	-	expression tag	UNP Q12306
C	-102	PRO	-	expression tag	UNP Q12306
C	-101	ARG	-	expression tag	UNP Q12306
C	-100	GLY	-	expression tag	UNP Q12306
C	-99	SER	-	expression tag	UNP Q12306
C	-98	ALA	-	expression tag	UNP Q12306
C	-97	SER	-	expression tag	UNP Q12306
D	-117	MET	-	initiating methionine	UNP Q12306
D	-116	GLY	-	expression tag	UNP Q12306
D	-115	SER	-	expression tag	UNP Q12306
D	-114	SER	-	expression tag	UNP Q12306
D	-113	HIS	-	expression tag	UNP Q12306
D	-112	HIS	-	expression tag	UNP Q12306
D	-111	HIS	-	expression tag	UNP Q12306
D	-110	HIS	-	expression tag	UNP Q12306
D	-109	HIS	-	expression tag	UNP Q12306
D	-108	HIS	-	expression tag	UNP Q12306
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D	-106	SER	-	expression tag	UNP Q12306
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D	-103	VAL	-	expression tag	UNP Q12306
D	-102	PRO	-	expression tag	UNP Q12306
D	-101	ARG	-	expression tag	UNP Q12306
D	-100	GLY	-	expression tag	UNP Q12306
D	-99	SER	-	expression tag	UNP Q12306
D	-98	ALA	-	expression tag	UNP Q12306
D	-97	SER	-	expression tag	UNP Q12306
E	-117	MET	-	initiating methionine	UNP Q12306
E	-116	GLY	-	expression tag	UNP Q12306
E	-115	SER	-	expression tag	UNP Q12306
E	-114	SER	-	expression tag	UNP Q12306
E	-113	HIS	-	expression tag	UNP Q12306
E	-112	HIS	-	expression tag	UNP Q12306
E	-111	HIS	-	expression tag	UNP Q12306
E	-110	HIS	-	expression tag	UNP Q12306
E	-109	HIS	-	expression tag	UNP Q12306
E	-108	HIS	-	expression tag	UNP Q12306

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E	-107	GLY	-	expression tag	UNP Q12306
E	-106	SER	-	expression tag	UNP Q12306
E	-105	GLY	-	expression tag	UNP Q12306
E	-104	LEU	-	expression tag	UNP Q12306
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E	-101	ARG	-	expression tag	UNP Q12306
E	-100	GLY	-	expression tag	UNP Q12306
E	-99	SER	-	expression tag	UNP Q12306
E	-98	ALA	-	expression tag	UNP Q12306
E	-97	SER	-	expression tag	UNP Q12306
F	-117	MET	-	initiating methionine	UNP Q12306
F	-116	GLY	-	expression tag	UNP Q12306
F	-115	SER	-	expression tag	UNP Q12306
F	-114	SER	-	expression tag	UNP Q12306
F	-113	HIS	-	expression tag	UNP Q12306
F	-112	HIS	-	expression tag	UNP Q12306
F	-111	HIS	-	expression tag	UNP Q12306
F	-110	HIS	-	expression tag	UNP Q12306
F	-109	HIS	-	expression tag	UNP Q12306
F	-108	HIS	-	expression tag	UNP Q12306
F	-107	GLY	-	expression tag	UNP Q12306
F	-106	SER	-	expression tag	UNP Q12306
F	-105	GLY	-	expression tag	UNP Q12306
F	-104	LEU	-	expression tag	UNP Q12306
F	-103	VAL	-	expression tag	UNP Q12306
F	-102	PRO	-	expression tag	UNP Q12306
F	-101	ARG	-	expression tag	UNP Q12306
F	-100	GLY	-	expression tag	UNP Q12306
F	-99	SER	-	expression tag	UNP Q12306
F	-98	ALA	-	expression tag	UNP Q12306
F	-97	SER	-	expression tag	UNP Q12306
G	-117	MET	-	initiating methionine	UNP Q12306
G	-116	GLY	-	expression tag	UNP Q12306
G	-115	SER	-	expression tag	UNP Q12306
G	-114	SER	-	expression tag	UNP Q12306
G	-113	HIS	-	expression tag	UNP Q12306
G	-112	HIS	-	expression tag	UNP Q12306
G	-111	HIS	-	expression tag	UNP Q12306
G	-110	HIS	-	expression tag	UNP Q12306
G	-109	HIS	-	expression tag	UNP Q12306
G	-108	HIS	-	expression tag	UNP Q12306

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Chain	Residue	Modelled	Actual	Comment	Reference
G	-107	GLY	-	expression tag	UNP Q12306
G	-106	SER	-	expression tag	UNP Q12306
G	-105	GLY	-	expression tag	UNP Q12306
G	-104	LEU	-	expression tag	UNP Q12306
G	-103	VAL	-	expression tag	UNP Q12306
G	-102	PRO	-	expression tag	UNP Q12306
G	-101	ARG	-	expression tag	UNP Q12306
G	-100	GLY	-	expression tag	UNP Q12306
G	-99	SER	-	expression tag	UNP Q12306
G	-98	ALA	-	expression tag	UNP Q12306
G	-97	SER	-	expression tag	UNP Q12306
H	-117	MET	-	initiating methionine	UNP Q12306
H	-116	GLY	-	expression tag	UNP Q12306
H	-115	SER	-	expression tag	UNP Q12306
H	-114	SER	-	expression tag	UNP Q12306
H	-113	HIS	-	expression tag	UNP Q12306
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H	-108	HIS	-	expression tag	UNP Q12306
H	-107	GLY	-	expression tag	UNP Q12306
H	-106	SER	-	expression tag	UNP Q12306
H	-105	GLY	-	expression tag	UNP Q12306
H	-104	LEU	-	expression tag	UNP Q12306
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H	-102	PRO	-	expression tag	UNP Q12306
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H	-99	SER	-	expression tag	UNP Q12306
H	-98	ALA	-	expression tag	UNP Q12306
H	-97	SER	-	expression tag	UNP Q12306
I	-117	MET	-	initiating methionine	UNP Q12306
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I	-115	SER	-	expression tag	UNP Q12306
I	-114	SER	-	expression tag	UNP Q12306
I	-113	HIS	-	expression tag	UNP Q12306
I	-112	HIS	-	expression tag	UNP Q12306
I	-111	HIS	-	expression tag	UNP Q12306
I	-110	HIS	-	expression tag	UNP Q12306
I	-109	HIS	-	expression tag	UNP Q12306
I	-108	HIS	-	expression tag	UNP Q12306

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I	-105	GLY	-	expression tag	UNP Q12306
I	-104	LEU	-	expression tag	UNP Q12306
I	-103	VAL	-	expression tag	UNP Q12306
I	-102	PRO	-	expression tag	UNP Q12306
I	-101	ARG	-	expression tag	UNP Q12306
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J	-110	HIS	-	expression tag	UNP Q12306
J	-109	HIS	-	expression tag	UNP Q12306
J	-108	HIS	-	expression tag	UNP Q12306
J	-107	GLY	-	expression tag	UNP Q12306
J	-106	SER	-	expression tag	UNP Q12306
J	-105	GLY	-	expression tag	UNP Q12306
J	-104	LEU	-	expression tag	UNP Q12306
J	-103	VAL	-	expression tag	UNP Q12306
J	-102	PRO	-	expression tag	UNP Q12306
J	-101	ARG	-	expression tag	UNP Q12306
J	-100	GLY	-	expression tag	UNP Q12306
J	-99	SER	-	expression tag	UNP Q12306
J	-98	ALA	-	expression tag	UNP Q12306
J	-97	SER	-	expression tag	UNP Q12306

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: Islet amyloid polypeptide (SUMO-tagged)

Chain A: 10% 6% 85%

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

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

ASN THR ALA THR CYS ALA THR GLN ARG LEU ALA N14 V17 N22 I26 S29 T30 N31 V32 G33 S34 N35 T36 Y37

- Molecule 1: Islet amyloid polypeptide (SUMO-tagged)

Chain B: 9% 6% 85%

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

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

ASN THR ALA THR CYS ALA THR GLN ARG LEU ALA N14 V17 S20 N21 N22 I26 S29 T30 N31 V32 G33 S34 N35 T36 Y37

- Molecule 1: Islet amyloid polypeptide (SUMO-tagged)

Chain C: 10% 6% 85%

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

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

ASN THR ALA THR CYS ALA THR GLN ARG LEU ALA N14 V17 N22 I26 S29 T30 N31 V32 G33 S34 N35 T36 Y37

- Molecule 1: Islet amyloid polypeptide (SUMO-tagged)

Chain D: 9% 6% 85%

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

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

ASN THR ALA THR CYS ALA THR GLN ARG LEU ALA N14 V17 S20 N21 N22 I26 S29 T30 N31 V32 G33 S34 N35 T36 Y37

- Molecule 1: Islet amyloid polypeptide (SUMO-tagged)

Chain E: 9% 6% 85%

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

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

ASN THR ALA THR CYS ALA THR GLN ARG LEU ALA N14 V17 S20 N21 N22 I26 S29 T30 N31 V32 G33 S34 N35 T36 Y37

- Molecule 1: Islet amyloid polypeptide (SUMO-tagged)

Chain F: 10% 6% 85%

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

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

ASN THR ALA THR CYS ALA THR GLN ARG LEU ALA N14 V17 N22 I26 S29 T30 N31 V32 G33 S34 N35 T36 Y37

- Molecule 1: Islet amyloid polypeptide (SUMO-tagged)

Chain G: 9% 6% 85%

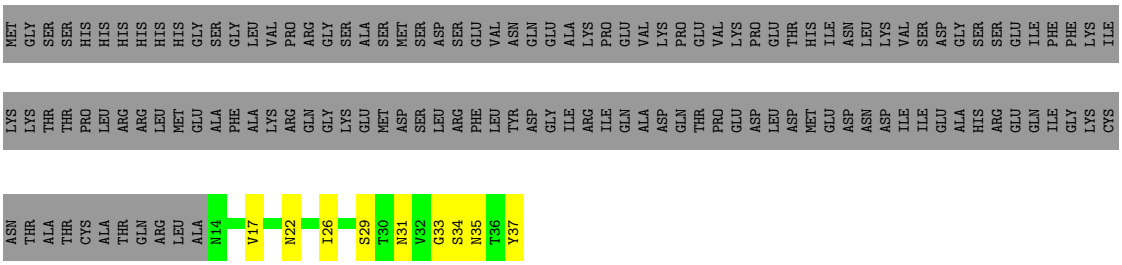
MET GLY SER SER HIS HIS HIS HIS HIS HIS GLY SER GLY LEU VAL PRO ARG GLY SER ALA SER MET SER ASP SER GLU VAL ASN GLN GLU ALA LYS PRO GLU VAL LYS PRO GLU THR HIS ILE ASN LEU LYS VAL SER ASP GLY SER SER GLU ILE PHE PHE LYS ILE

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

ASN THR ALA THR CYS ALA THR GLN ARG LEU ALA N14 V17 S20 N21 N22 I26 S29 T30 N31 V32 G33 S34 N35 T36 Y37

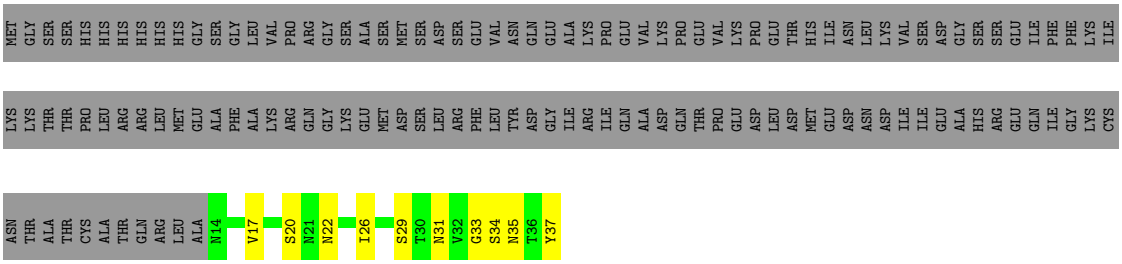
- Molecule 1: Islet amyloid polypeptide (SUMO-tagged)

Chain H: 10% 6% 85%



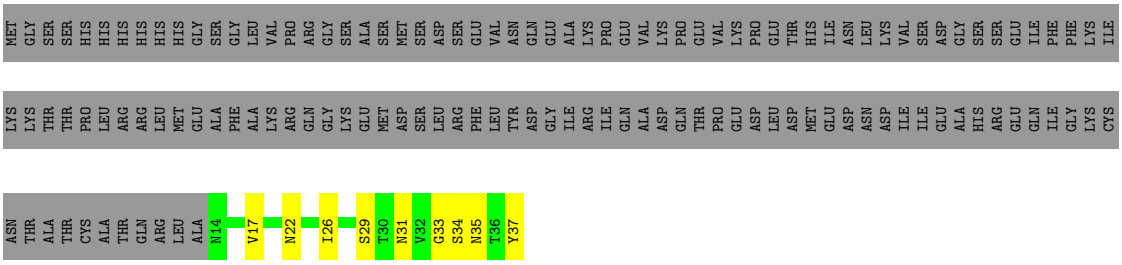
• Molecule 1: Islet amyloid polypeptide (SUMO-tagged)

Chain I: 9% 6% 85%



• Molecule 1: Islet amyloid polypeptide (SUMO-tagged)

Chain J: 10% 6% 85%



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	HELICAL	Depositor
Imposed symmetry	HELICAL, twist=179.42°, rise=2.406 Å, axial sym=C1	Depositor
Number of segments used	25767	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 ($e^-/Å^2$)	44	Depositor
Minimum defocus (nm)	2000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	14.139	Depositor
Minimum map value	-8.656	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	2.0	Depositor
Map size (Å)	136.192, 136.192, 136.192	wwPDB
Map dimensions	128, 128, 128	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.064, 1.064, 1.064	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	0.35	0/183	0.73	0/247
1	B	0.35	0/183	0.73	0/247
1	C	0.35	0/183	0.73	0/247
1	D	0.35	0/183	0.73	0/247
1	E	0.35	0/183	0.73	0/247
1	F	0.35	0/183	0.74	0/247
1	G	0.35	0/183	0.73	0/247
1	H	0.35	0/183	0.73	0/247
1	I	0.35	0/183	0.73	0/247
1	J	0.35	0/183	0.73	0/247
All	All	0.35	0/1830	0.73	0/2470

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	180	0	164	14	0
1	B	180	0	164	16	0
1	C	180	0	164	15	0
1	D	180	0	164	15	0
1	E	180	0	164	16	0
1	F	180	0	164	15	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	G	180	0	164	10	0
1	H	180	0	164	10	0
1	I	180	0	164	10	0
1	J	180	0	164	9	0
All	All	1800	0	1640	83	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 24.

All (83) 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:H:34:SER:HB2	1:H:37:TYR:OH	1.74	0.88
1:D:34:SER:HB2	1:D:37:TYR:OH	1.74	0.87
1:A:34:SER:HB2	1:A:37:TYR:OH	1.74	0.87
1:J:34:SER:HB2	1:J:37:TYR:OH	1.74	0.87
1:E:34:SER:HB2	1:E:37:TYR:OH	1.74	0.87
1:F:34:SER:HB2	1:F:37:TYR:OH	1.74	0.87
1:I:34:SER:HB2	1:I:37:TYR:OH	1.74	0.86
1:B:34:SER:HB2	1:B:37:TYR:OH	1.74	0.86
1:C:34:SER:HB2	1:C:37:TYR:OH	1.74	0.86
1:G:34:SER:HB2	1:G:37:TYR:OH	1.74	0.85
1:B:17:VAL:O	1:F:17:VAL:HB	1.79	0.83
1:A:17:VAL:HB	1:E:17:VAL:O	1.79	0.82
1:B:17:VAL:HB	1:C:17:VAL:O	1.80	0.81
1:C:17:VAL:HB	1:G:17:VAL:O	1.80	0.81
1:D:17:VAL:O	1:H:17:VAL:HB	1.80	0.81
1:E:17:VAL:HB	1:I:17:VAL:O	1.80	0.81
1:F:17:VAL:O	1:J:17:VAL:HB	1.81	0.81
1:A:17:VAL:O	1:D:17:VAL:HB	1.81	0.80
1:A:22:ASN:HA	1:E:22:ASN:O	1.96	0.66
1:D:22:ASN:O	1:H:22:ASN:HA	1.96	0.66
1:E:22:ASN:HA	1:I:22:ASN:O	1.96	0.66
1:B:22:ASN:O	1:F:22:ASN:HA	1.96	0.65
1:A:22:ASN:O	1:D:22:ASN:HA	1.97	0.65
1:B:22:ASN:HA	1:C:22:ASN:O	1.96	0.64
1:C:22:ASN:HA	1:G:22:ASN:O	1.96	0.64
1:F:22:ASN:O	1:J:22:ASN:HA	1.97	0.64
1:C:33:GLY:O	1:C:35:ASN:ND2	2.35	0.60
1:G:33:GLY:O	1:G:35:ASN:ND2	2.35	0.60
1:B:33:GLY:O	1:B:35:ASN:ND2	2.35	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:29:SER:O	1:I:29:SER:HA	2.02	0.60
1:F:33:GLY:O	1:F:35:ASN:ND2	2.35	0.60
1:A:29:SER:O	1:E:29:SER:HA	2.02	0.59
1:A:29:SER:HA	1:D:29:SER:O	2.02	0.59
1:B:29:SER:O	1:C:29:SER:HA	2.02	0.59
1:J:33:GLY:O	1:J:35:ASN:ND2	2.35	0.59
1:C:29:SER:O	1:G:29:SER:HA	2.02	0.59
1:F:29:SER:HA	1:J:29:SER:O	2.02	0.59
1:B:29:SER:HA	1:F:29:SER:O	2.02	0.59
1:D:29:SER:HA	1:H:29:SER:O	2.02	0.59
1:B:26:ILE:HA	1:F:26:ILE:O	2.04	0.57
1:C:26:ILE:O	1:G:26:ILE:HA	2.05	0.57
1:A:33:GLY:O	1:A:35:ASN:ND2	2.35	0.57
1:E:33:GLY:O	1:E:35:ASN:ND2	2.35	0.57
1:A:26:ILE:O	1:E:26:ILE:HA	2.04	0.57
1:B:26:ILE:O	1:C:26:ILE:HA	2.05	0.57
1:D:33:GLY:O	1:D:35:ASN:ND2	2.35	0.57
1:I:33:GLY:O	1:I:35:ASN:ND2	2.35	0.57
1:D:26:ILE:HA	1:H:26:ILE:O	2.05	0.57
1:E:26:ILE:O	1:I:26:ILE:HA	2.05	0.57
1:H:33:GLY:O	1:H:35:ASN:ND2	2.35	0.57
1:A:26:ILE:HA	1:D:26:ILE:O	2.05	0.57
1:F:26:ILE:HA	1:J:26:ILE:O	2.05	0.56
1:C:31:ASN:HD22	1:G:31:ASN:HB2	1.74	0.53
1:D:22:ASN:N	1:D:22:ASN:OD1	2.42	0.53
1:B:31:ASN:HD22	1:C:31:ASN:HB2	1.74	0.53
1:F:31:ASN:HB2	1:J:31:ASN:HD22	1.74	0.53
1:H:22:ASN:N	1:H:22:ASN:OD1	2.42	0.52
1:B:22:ASN:N	1:B:22:ASN:OD1	2.42	0.52
1:B:31:ASN:HB2	1:F:31:ASN:HD22	1.75	0.52
1:F:22:ASN:N	1:F:22:ASN:OD1	2.42	0.52
1:E:31:ASN:HD22	1:I:31:ASN:HB2	1.74	0.52
1:A:31:ASN:HB2	1:D:31:ASN:HD22	1.74	0.52
1:C:22:ASN:N	1:C:22:ASN:OD1	2.42	0.52
1:A:22:ASN:N	1:A:22:ASN:OD1	2.42	0.51
1:D:31:ASN:HB2	1:H:31:ASN:HD22	1.74	0.51
1:J:22:ASN:OD1	1:J:22:ASN:N	2.42	0.51
1:A:31:ASN:HD22	1:E:31:ASN:HB2	1.75	0.51
1:G:22:ASN:N	1:G:22:ASN:OD1	2.42	0.51
1:E:22:ASN:N	1:E:22:ASN:OD1	2.42	0.51
1:I:22:ASN:N	1:I:22:ASN:OD1	2.42	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:20:SER:O	1:G:20:SER:OG	2.29	0.48
1:E:20:SER:O	1:E:20:SER:OG	2.29	0.47
1:I:20:SER:O	1:I:20:SER:OG	2.29	0.46
1:H:34:SER:HB2	1:H:37:TYR:HH	1.75	0.46
1:B:22:ASN:OD1	1:F:22:ASN:HB3	2.20	0.42
1:A:22:ASN:HB3	1:E:22:ASN:OD1	2.20	0.42
1:B:22:ASN:HB3	1:C:22:ASN:OD1	2.21	0.41
1:C:22:ASN:HB3	1:G:22:ASN:OD1	2.21	0.41
1:B:20:SER:O	1:B:20:SER:OG	2.29	0.41
1:D:22:ASN:OD1	1:H:22:ASN:HB3	2.21	0.41
1:E:22:ASN:HB3	1:I:22:ASN:OD1	2.21	0.41
1:D:20:SER:O	1:D:20:SER:OG	2.29	0.40
1:F:22:ASN:OD1	1:J:22:ASN:HB3	2.22	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	22/155 (14%)	16 (73%)	6 (27%)	0	100	100
1	B	22/155 (14%)	16 (73%)	6 (27%)	0	100	100
1	C	22/155 (14%)	16 (73%)	6 (27%)	0	100	100
1	D	22/155 (14%)	16 (73%)	6 (27%)	0	100	100
1	E	22/155 (14%)	16 (73%)	6 (27%)	0	100	100
1	F	22/155 (14%)	16 (73%)	6 (27%)	0	100	100
1	G	22/155 (14%)	16 (73%)	6 (27%)	0	100	100
1	H	22/155 (14%)	16 (73%)	6 (27%)	0	100	100
1	I	22/155 (14%)	16 (73%)	6 (27%)	0	100	100
1	J	22/155 (14%)	16 (73%)	6 (27%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
All	All	220/1550 (14%)	160 (73%)	60 (27%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	21/135 (16%)	21 (100%)	0	100	100
1	B	21/135 (16%)	21 (100%)	0	100	100
1	C	21/135 (16%)	21 (100%)	0	100	100
1	D	21/135 (16%)	21 (100%)	0	100	100
1	E	21/135 (16%)	21 (100%)	0	100	100
1	F	21/135 (16%)	21 (100%)	0	100	100
1	G	21/135 (16%)	21 (100%)	0	100	100
1	H	21/135 (16%)	21 (100%)	0	100	100
1	I	21/135 (16%)	21 (100%)	0	100	100
1	J	21/135 (16%)	21 (100%)	0	100	100
All	All	210/1350 (16%)	210 (100%)	0	100	100

There are no protein residues with a non-rotameric sidechain to report.

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (8) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	31	ASN
1	B	31	ASN
1	C	31	ASN
1	D	31	ASN
1	E	31	ASN
1	F	31	ASN
1	H	31	ASN

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Mol	Chain	Res	Type
1	J	31	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 monosaccharides 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.

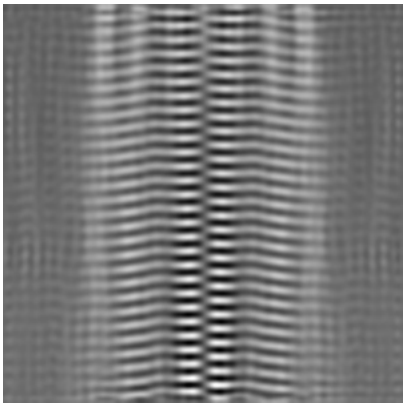
6 Map visualisation (i)

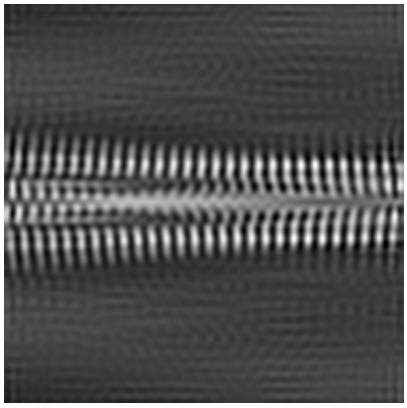
This section contains visualisations of the EMDB entry EMD-21410. These allow visual inspection of the internal detail of the map and identification of artifacts.

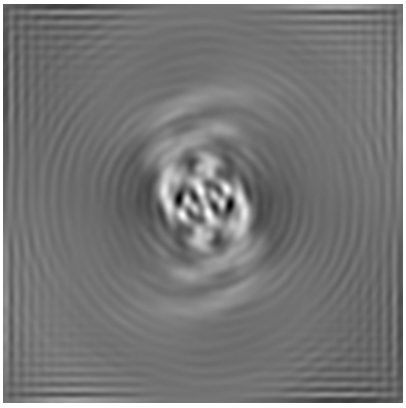
No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections (i)

6.1.1 Primary map


X

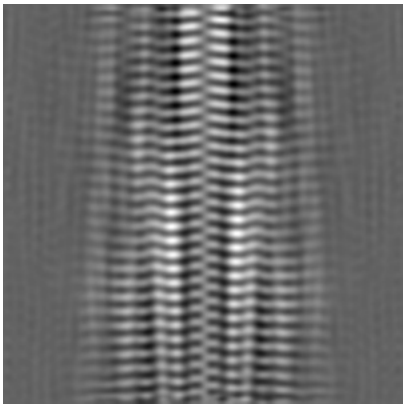

Y

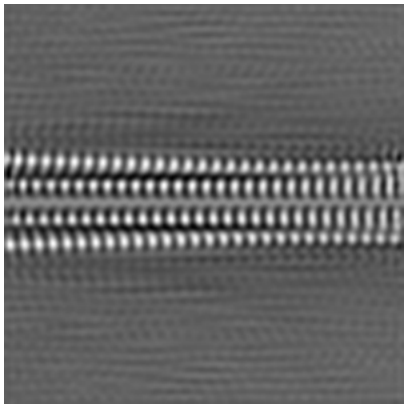

Z

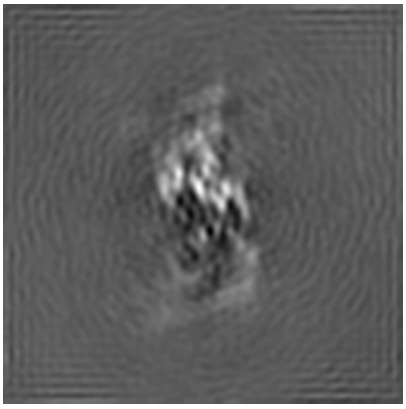
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

6.2.1 Primary map


X Index: 64

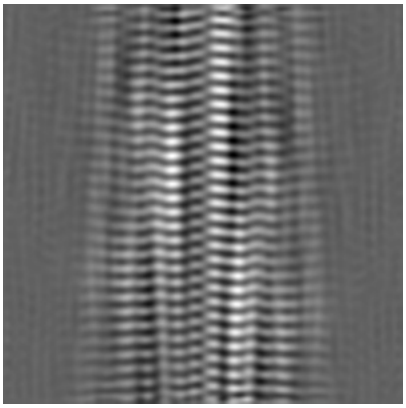

Y Index: 64

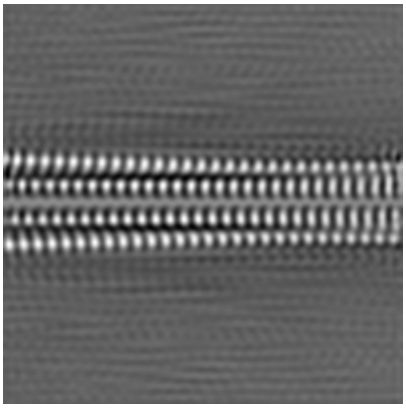

Z Index: 64

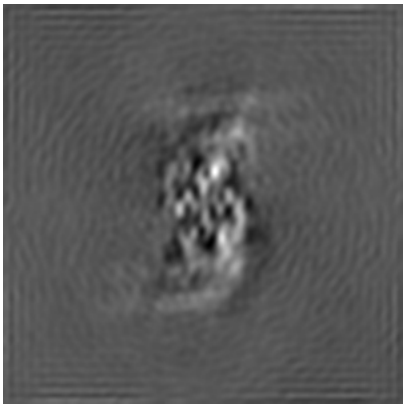
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

6.3.1 Primary map


X Index: 63

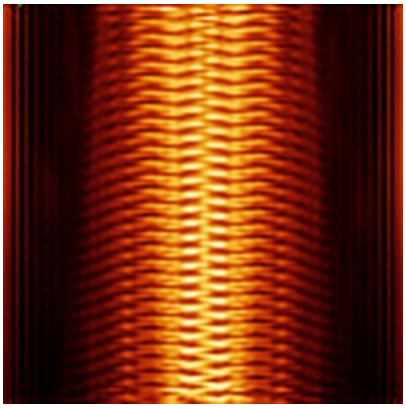

Y Index: 64

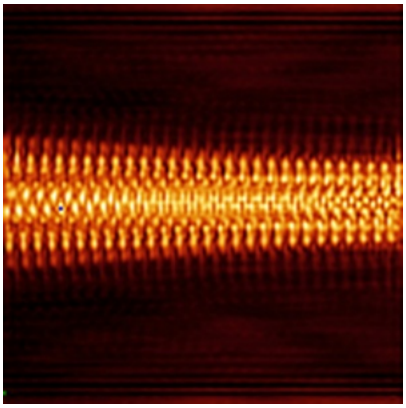

Z Index: 126

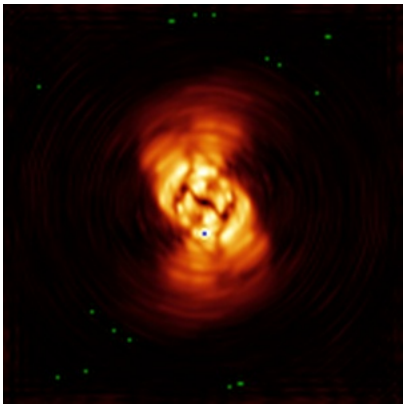
The images above show the largest variance slices of the map in three orthogonal directions.

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

6.4.1 Primary map


X

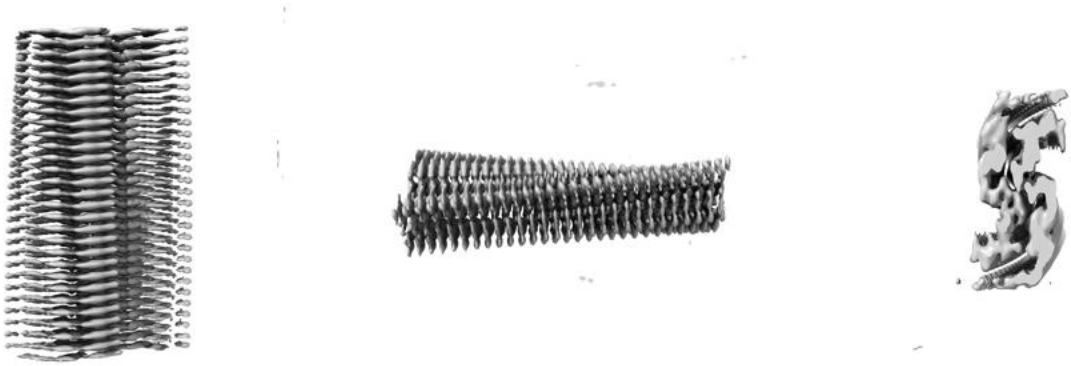

Y


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views (i)

6.5.1 Primary map



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 2.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

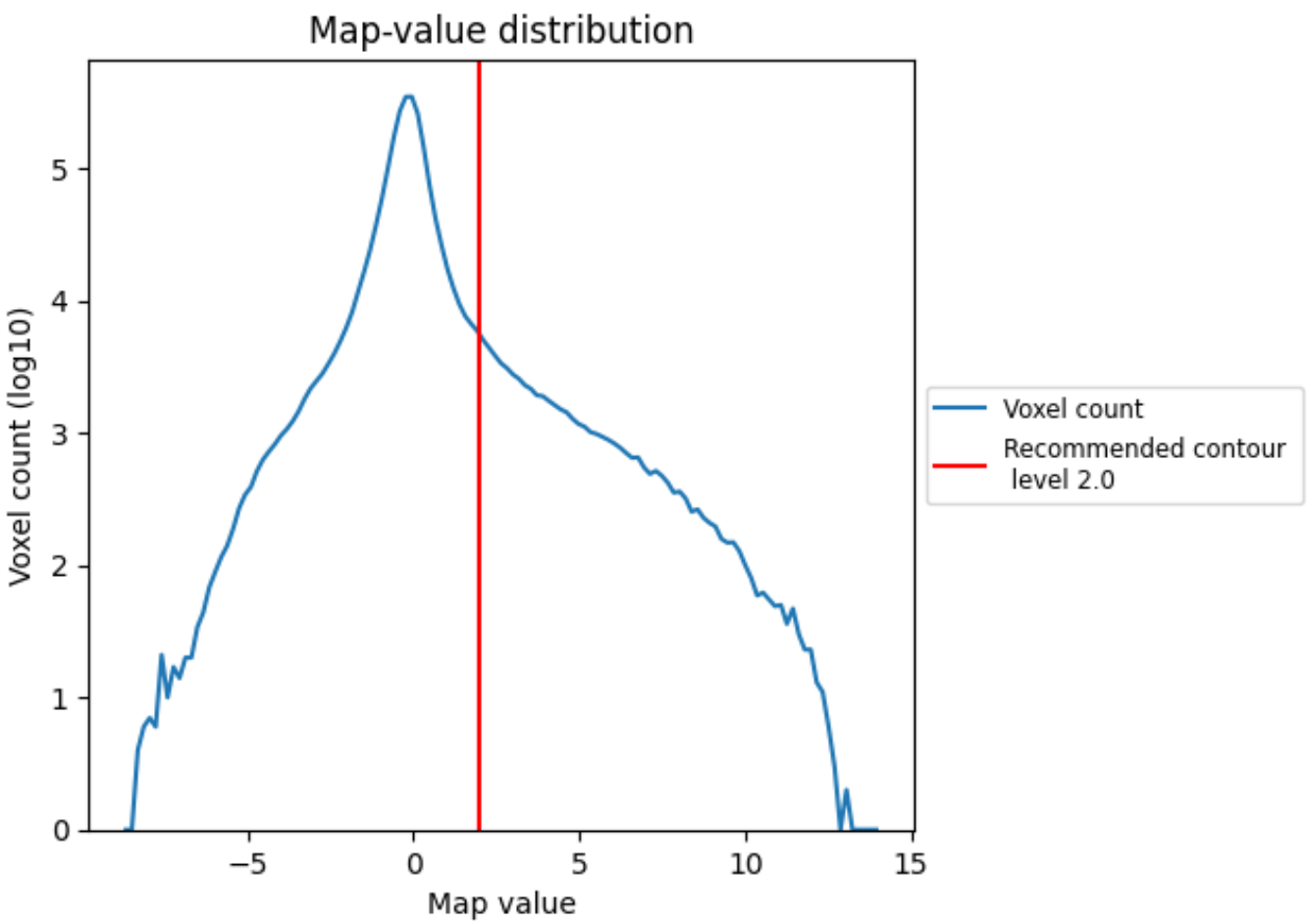
6.6 Mask visualisation (i)

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

7 Map analysis (i)

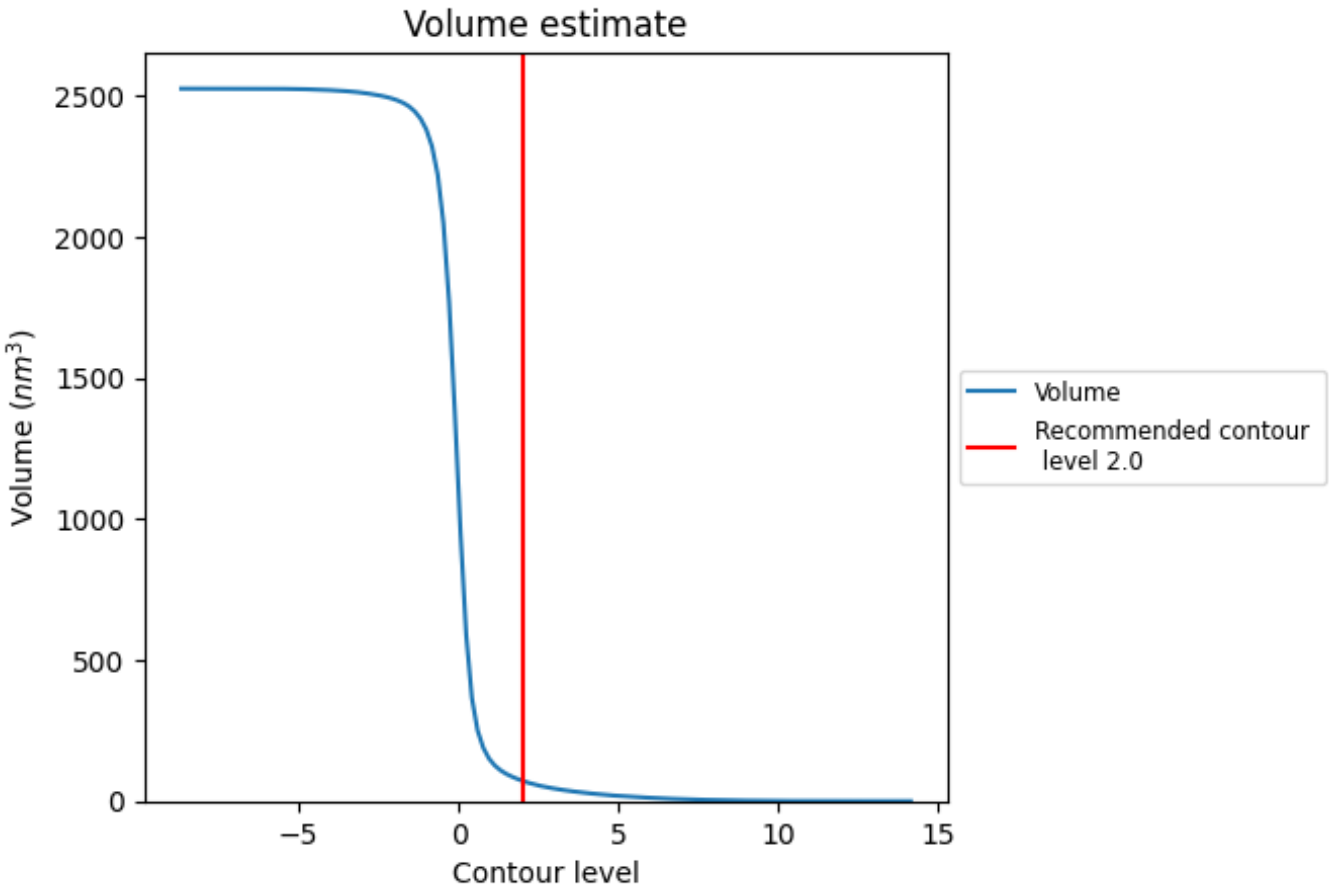
This section contains the results of statistical analysis of the map.

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.

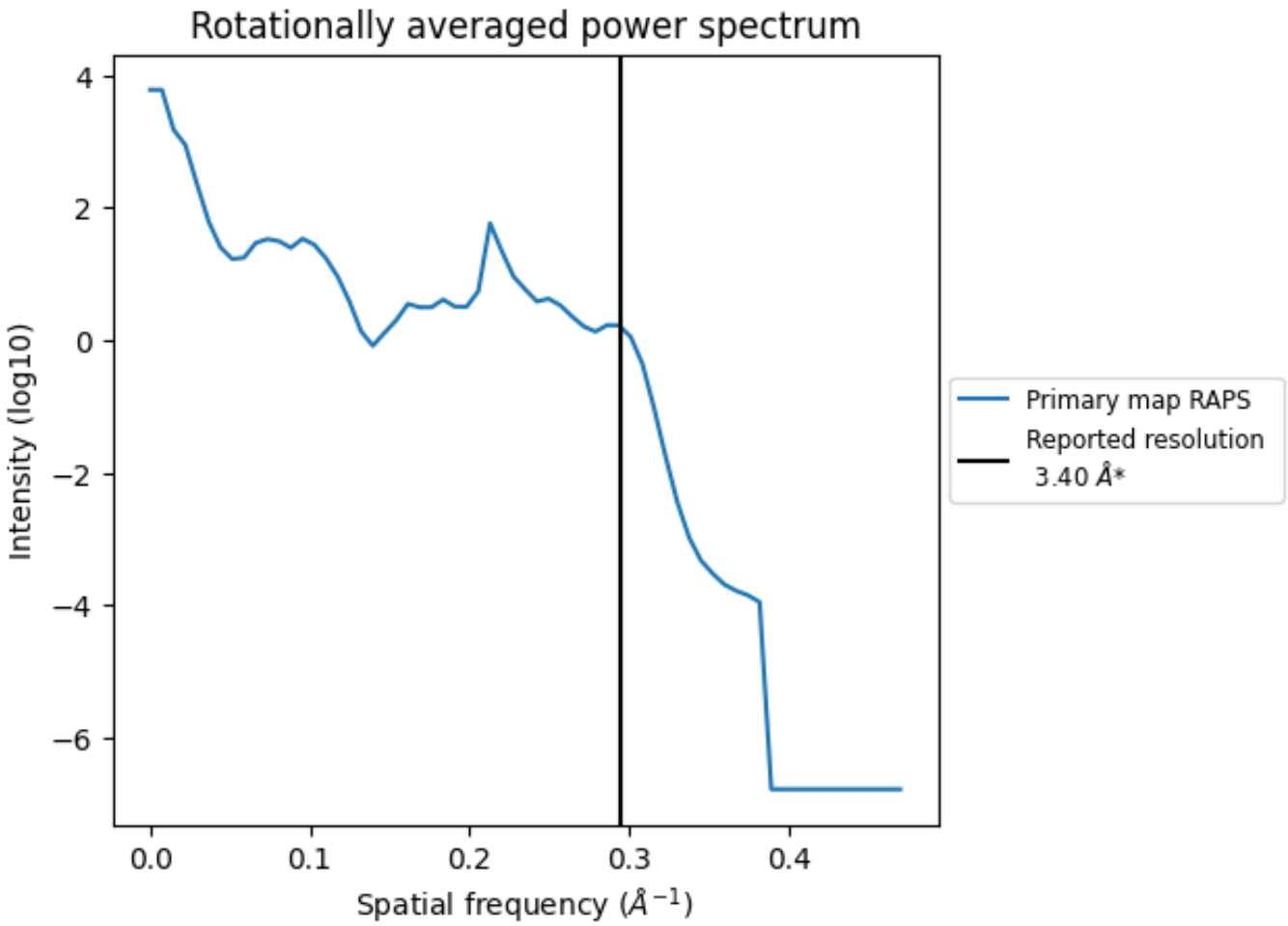
7.2 Volume estimate (i)



The volume at the recommended contour level is 72 nm^3; this corresponds to an approximate mass of 65 kDa.

The volume estimate graph shows how the enclosed volume varies with the contour level. The recommended contour level is shown as a vertical line and the intersection between the line and the curve gives the volume of the enclosed surface at the given level.

7.3 Rotationally averaged power spectrum (i)

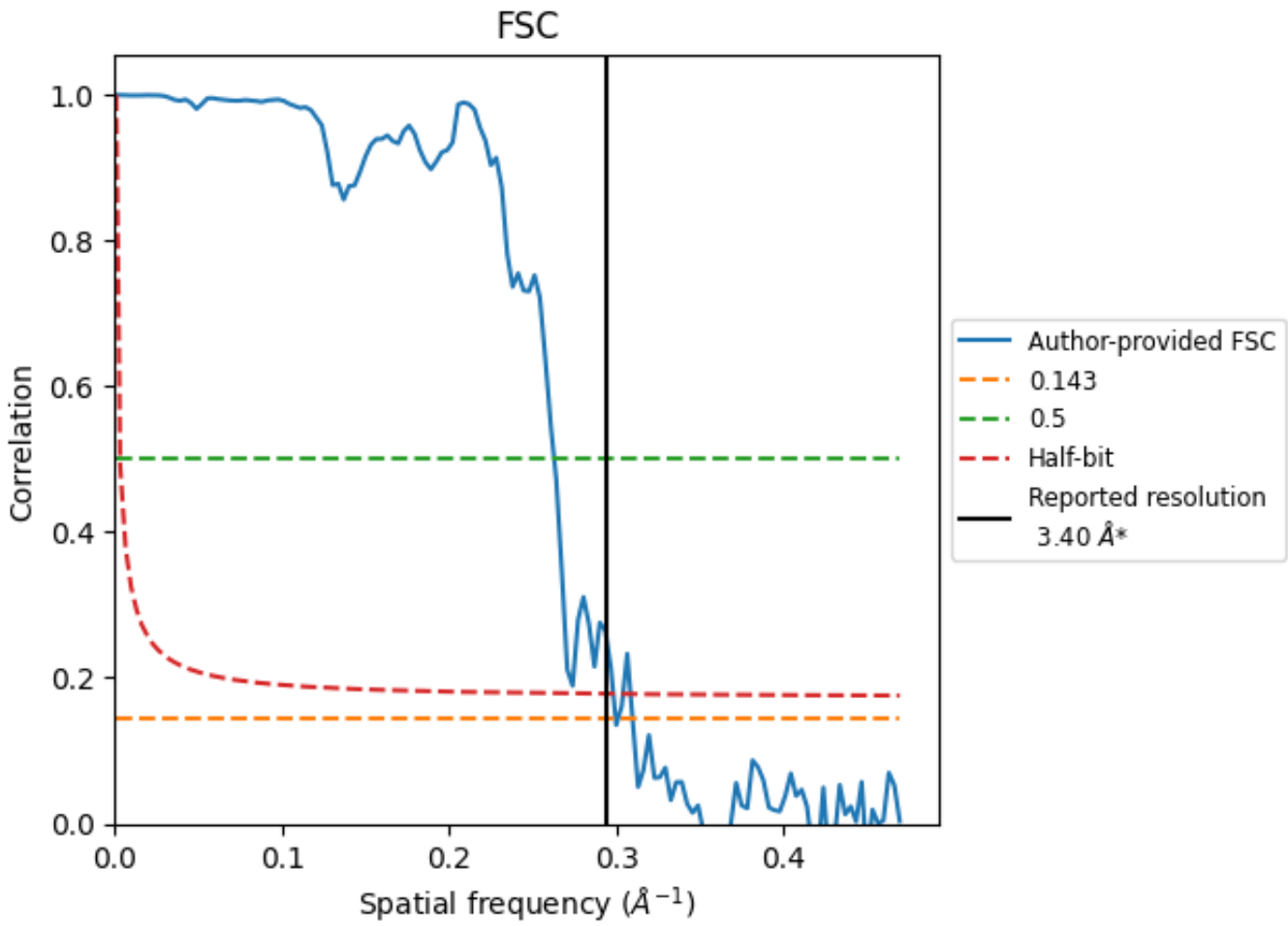


*Reported resolution corresponds to spatial frequency of 0.294 Å^{-1}

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 0.294 Å^{-1}

8.2 Resolution estimates (i)

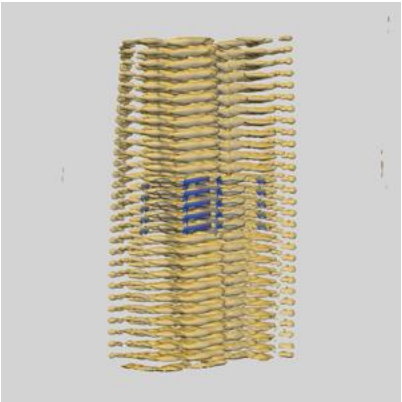
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.40	-	-
Author-provided FSC curve	3.33	3.80	3.35
Unmasked-calculated*	-	-	-

*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-21410 and PDB model 6VW2. Per-residue inclusion information can be found in section 3 on page 10.

9.1 Map-model overlay (i)



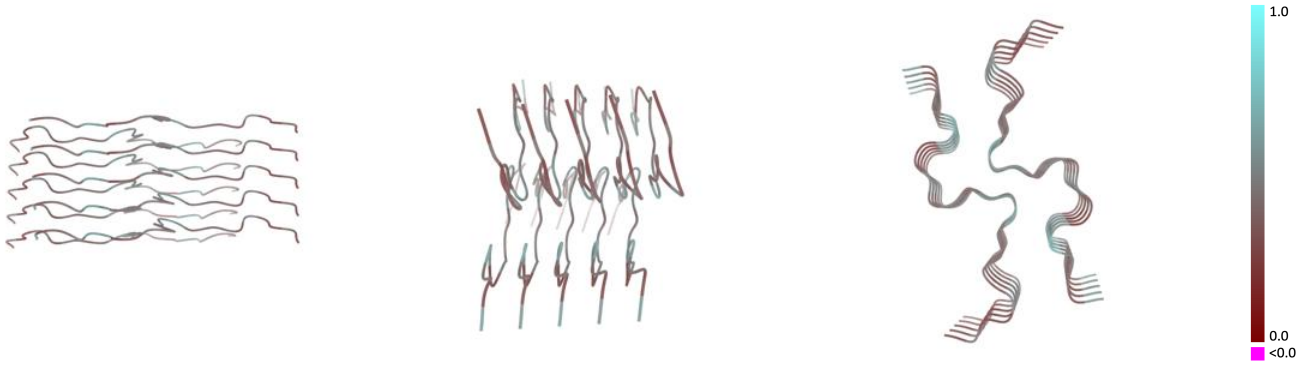
X

Y

Z

The images above show the 3D surface view of the map at the recommended contour level 2.0 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.

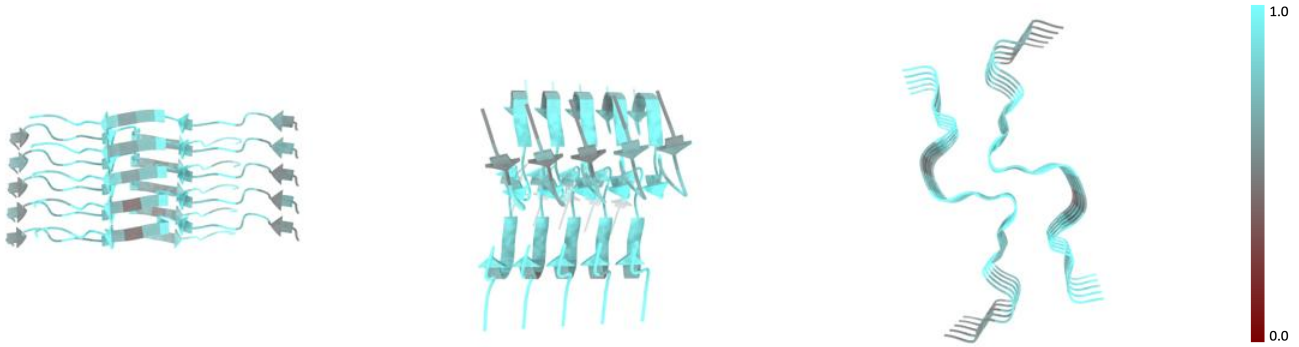
9.2 Q-score mapped to coordinate model (i)



X Y Z

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

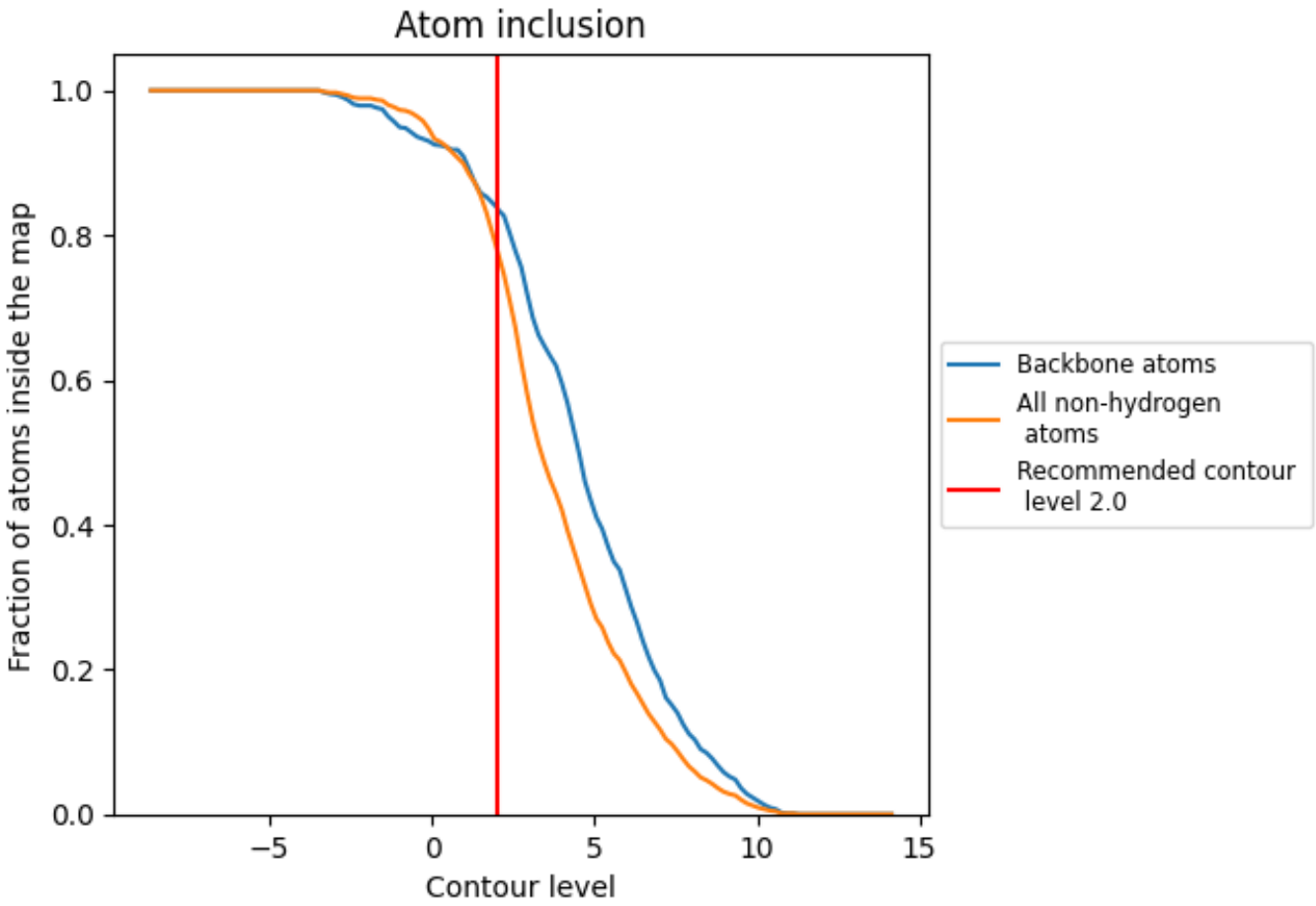
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (2.0).

9.4 Atom inclusion (i)



At the recommended contour level, 84% of all backbone atoms, 78% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary (i)

The table lists the average atom inclusion at the recommended contour level (2.0) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7800	0.4350
A	0.7880	0.4360
B	0.7820	0.4310
C	0.7650	0.4380
D	0.7710	0.4400
E	0.7650	0.4400
F	0.7880	0.4350
G	0.7760	0.4290
H	0.7820	0.4370
I	0.7930	0.4340
J	0.7880	0.4340