



Full wwPDB EM Validation Report ⓘ

Mar 6, 2026 – 05:58 PM UTC

PDB ID : 8F7C / pdb_00008f7c
EMDB ID : EMD-28902
Title : Cryo-EM structure of human pannexin 2
Authors : He, Z.; Yuan, P.
Deposited on : 2022-11-18
Resolution : 3.92 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.dev132
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) : 2.49

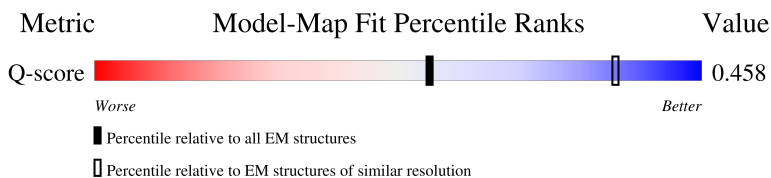
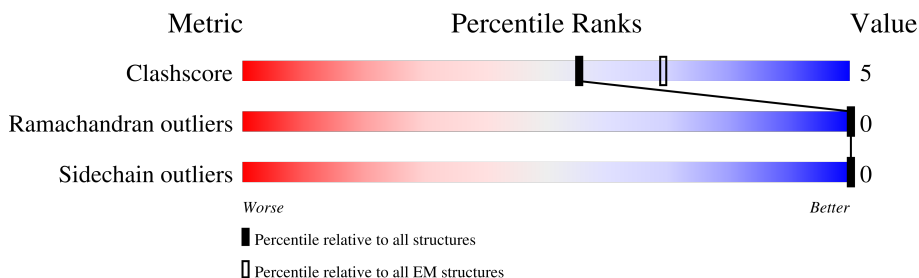
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.92 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.



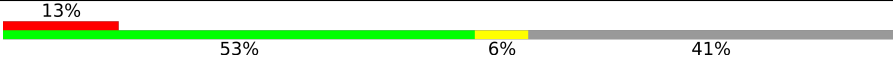
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	7862 (3.42 - 4.42)

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	487	13% 53% 6% 41%
1	B	487	13% 51% 8% 41%
1	C	487	13% 52% 7% 41%
1	D	487	13% 52% 7% 41%

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Mol	Chain	Length	Quality of chain
1	E	487	
1	F	487	
1	G	487	

2 Entry composition

There is only 1 type of molecule in this entry. The entry contains 15925 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 Pannexin-2, Soluble cytochrome b562 fusion.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	287	Total	C	N	O	S	0	0
			2275	1489	388	387	11		
1	B	287	Total	C	N	O	S	0	0
			2275	1489	388	387	11		
1	C	287	Total	C	N	O	S	0	0
			2275	1489	388	387	11		
1	D	287	Total	C	N	O	S	0	0
			2275	1489	388	387	11		
1	E	287	Total	C	N	O	S	0	0
			2275	1489	388	387	11		
1	F	287	Total	C	N	O	S	0	0
			2275	1489	388	387	11		
1	G	287	Total	C	N	O	S	0	0
			2275	1489	388	387	11		

There are 84 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	379	TRP	MET	conflict	UNP P0ABE7
A	474	ILE	HIS	conflict	UNP P0ABE7
A	478	LEU	-	expression tag	UNP P0ABE7
A	479	SER	-	expression tag	UNP P0ABE7
A	480	ASN	-	expression tag	UNP P0ABE7
A	481	SER	-	expression tag	UNP P0ABE7
A	482	LEU	-	expression tag	UNP P0ABE7
A	483	GLU	-	expression tag	UNP P0ABE7
A	484	VAL	-	expression tag	UNP P0ABE7
A	485	LEU	-	expression tag	UNP P0ABE7
A	486	PHE	-	expression tag	UNP P0ABE7
A	487	GLN	-	expression tag	UNP P0ABE7
B	379	TRP	MET	conflict	UNP P0ABE7
B	474	ILE	HIS	conflict	UNP P0ABE7
B	478	LEU	-	expression tag	UNP P0ABE7
B	479	SER	-	expression tag	UNP P0ABE7

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Chain	Residue	Modelled	Actual	Comment	Reference
B	480	ASN	-	expression tag	UNP P0ABE7
B	481	SER	-	expression tag	UNP P0ABE7
B	482	LEU	-	expression tag	UNP P0ABE7
B	483	GLU	-	expression tag	UNP P0ABE7
B	484	VAL	-	expression tag	UNP P0ABE7
B	485	LEU	-	expression tag	UNP P0ABE7
B	486	PHE	-	expression tag	UNP P0ABE7
B	487	GLN	-	expression tag	UNP P0ABE7
C	379	TRP	MET	conflict	UNP P0ABE7
C	474	ILE	HIS	conflict	UNP P0ABE7
C	478	LEU	-	expression tag	UNP P0ABE7
C	479	SER	-	expression tag	UNP P0ABE7
C	480	ASN	-	expression tag	UNP P0ABE7
C	481	SER	-	expression tag	UNP P0ABE7
C	482	LEU	-	expression tag	UNP P0ABE7
C	483	GLU	-	expression tag	UNP P0ABE7
C	484	VAL	-	expression tag	UNP P0ABE7
C	485	LEU	-	expression tag	UNP P0ABE7
C	486	PHE	-	expression tag	UNP P0ABE7
C	487	GLN	-	expression tag	UNP P0ABE7
D	379	TRP	MET	conflict	UNP P0ABE7
D	474	ILE	HIS	conflict	UNP P0ABE7
D	478	LEU	-	expression tag	UNP P0ABE7
D	479	SER	-	expression tag	UNP P0ABE7
D	480	ASN	-	expression tag	UNP P0ABE7
D	481	SER	-	expression tag	UNP P0ABE7
D	482	LEU	-	expression tag	UNP P0ABE7
D	483	GLU	-	expression tag	UNP P0ABE7
D	484	VAL	-	expression tag	UNP P0ABE7
D	485	LEU	-	expression tag	UNP P0ABE7
D	486	PHE	-	expression tag	UNP P0ABE7
D	487	GLN	-	expression tag	UNP P0ABE7
E	379	TRP	MET	conflict	UNP P0ABE7
E	474	ILE	HIS	conflict	UNP P0ABE7
E	478	LEU	-	expression tag	UNP P0ABE7
E	479	SER	-	expression tag	UNP P0ABE7
E	480	ASN	-	expression tag	UNP P0ABE7
E	481	SER	-	expression tag	UNP P0ABE7
E	482	LEU	-	expression tag	UNP P0ABE7
E	483	GLU	-	expression tag	UNP P0ABE7
E	484	VAL	-	expression tag	UNP P0ABE7
E	485	LEU	-	expression tag	UNP P0ABE7

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Chain	Residue	Modelled	Actual	Comment	Reference
E	486	PHE	-	expression tag	UNP P0ABE7
E	487	GLN	-	expression tag	UNP P0ABE7
F	379	TRP	MET	conflict	UNP P0ABE7
F	474	ILE	HIS	conflict	UNP P0ABE7
F	478	LEU	-	expression tag	UNP P0ABE7
F	479	SER	-	expression tag	UNP P0ABE7
F	480	ASN	-	expression tag	UNP P0ABE7
F	481	SER	-	expression tag	UNP P0ABE7
F	482	LEU	-	expression tag	UNP P0ABE7
F	483	GLU	-	expression tag	UNP P0ABE7
F	484	VAL	-	expression tag	UNP P0ABE7
F	485	LEU	-	expression tag	UNP P0ABE7
F	486	PHE	-	expression tag	UNP P0ABE7
F	487	GLN	-	expression tag	UNP P0ABE7
G	379	TRP	MET	conflict	UNP P0ABE7
G	474	ILE	HIS	conflict	UNP P0ABE7
G	478	LEU	-	expression tag	UNP P0ABE7
G	479	SER	-	expression tag	UNP P0ABE7
G	480	ASN	-	expression tag	UNP P0ABE7
G	481	SER	-	expression tag	UNP P0ABE7
G	482	LEU	-	expression tag	UNP P0ABE7
G	483	GLU	-	expression tag	UNP P0ABE7
G	484	VAL	-	expression tag	UNP P0ABE7
G	485	LEU	-	expression tag	UNP P0ABE7
G	486	PHE	-	expression tag	UNP P0ABE7
G	487	GLN	-	expression tag	UNP P0ABE7

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

● Molecule 1: Pannexin-2, Soluble cytochrome b562 fusion



ALA	GLU	GLN	LEU	LEU	THR	THR	ARG	ASN	ALA	ALA	TYR	ILE	GLN	LYS	LYS	THR	PRO	PRO	LYS	GLY	SER	LEU	GLY	ASP	LEU	GLY	VAL	LEU	PHE	GLN																													
THR	LYS	MET	ARG	LYS	ALA	ALA	LEU	ASP	ALA	ALA	TYR	ILE	GLN	LYS	LYS	THR	PRO	PRO	LYS	GLY	SER	LEU	GLY	ASP	LEU	GLY	VAL	LEU	PHE	GLN																													
R334	Q335	W336	R337	R338	S339	Q340	F341	C342	D343	I344	N352	F353	N354	K359	S360	L361	N362	R363	L364	D365	F366	I367	T368	ASN	GLY	SER	PHE	ASP	ILE	VAL	GLY	GLN	ASP	ASN	LYS	LEU	ASP	ALA	ASP	LEU	GLY	THR	ALA	LEU	THR	ASN	GLY	VAL	GLN	ASP	ASN	ALA	ALA	GLN	VAL	LYS	ASP	ALA	LEU
GLU	LYS	SER	PRO	GLU	GLN	N208	L209	F210	E211	K212	E215	R216	R219	R230	L237	K253	A263	S264	P265	D266	GLY	ALA	ALA	GLY	ALA	GLY	ALA	GLY	P273	C279	R288	I289	I290	I310	I314	F315	R316	K317	S318	D323	K324	L325	H326	K327	V328	G329	I330	K331	T332	R333									
A106	G109	V110	D111	A112	F131	A132	A133	I134	V137	L154	L158	Q159	E160	I161	D162	N163	C164	Y165	H166	ARG	ALA	ALA	GLY	ARG	ALA	ALA	PRO	LYS	ILE	GLY	GLN	ILE	GLN	SER	LYS	GLY	PRO	GLY	ILE	THR	GLY	ARG	GLY	LYS	ARG	GLY	ILE	ILE	GLY	ASN	ALA	LYS							
MET	HIS	HIS	LEU	LEU	GLU	GLN	SER	ALA	ASP	MET	ALA	THR	ALA	LEU	LEU	ALA	GLY	GLY	LYS	ARG	GLU	LEU	LEU	ALA	ASP	ASP	LYS	LYS	ALA	G35	A36	L37	A38	A39	L40	L41	L42	Q43	L44	K45	L46	E47	D51	L65	K72	E77	Y80	C81	D105										

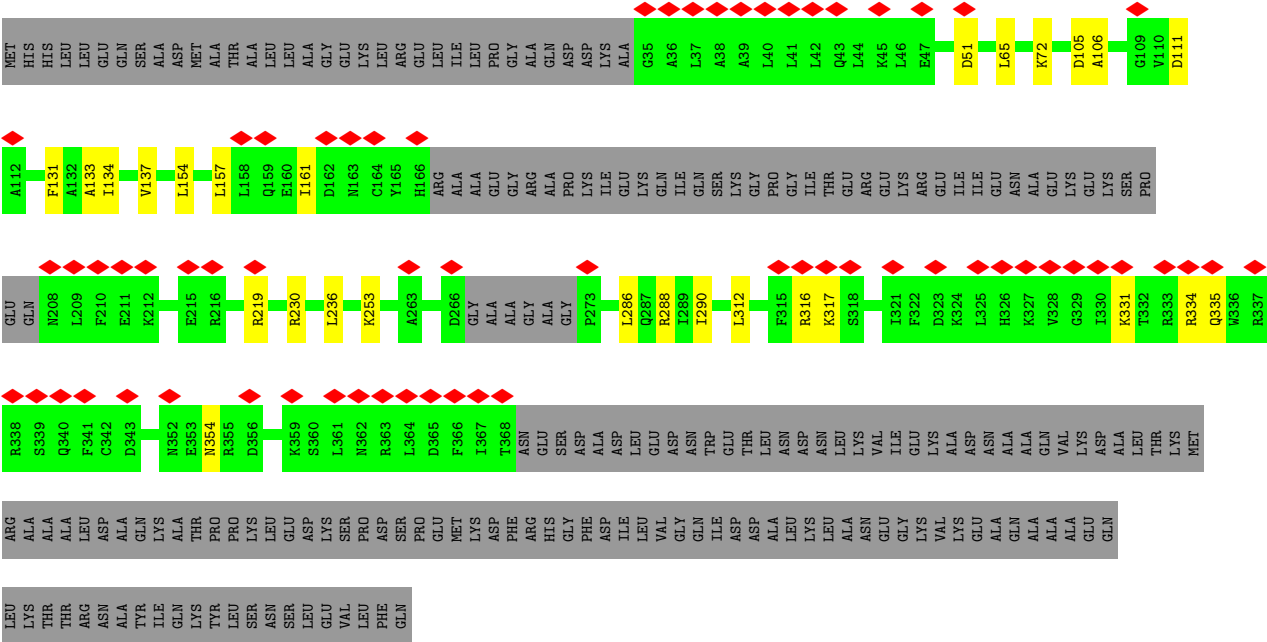
● Molecule 1: Pannexin-2, Soluble cytochrome b562 fusion



ALA	THR	PRO	PRO	LEU	GLU	ASP	LYS	SER	PRO	ASP	SER	PRO	GLU	MET	ASP	PHE	ARG	HIS	GLY	GLY	ALA	ASP	LEU	ASP	LEU	ILE	LEU	VAL	GLY	GLN	ILE	ASP	ASP	ALA	LEU	LYS	LEU	LEU	ALA	ALA	ASN	ILE	GLY	GLY	GLY	LYS	LYS	VAL	GLU	GLU	GLN	ALA	ALA	ALA	GLU	GLN	VAL	LYS	ASP	ASP	ALA	ALA	ALA	THR	THR	THR	ARG	ASN	ALA	TYR	ILE																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																																							
D343		N352		K359	S360	L361	N362	R363	L364	D365	F366	I367	T368	ASN	GLU	SER	ASP	ALA	GLY	ALA	ASP	LEU	ASP	LEU	GLY	ASP	ASP	TRP	GLY	GLN	ILE	THR	LEU	ASN	ASP	ASN	LEU	LEU	LYS	LEU	VAL	ILE	GLY	GLY	GLY	LYS	ALA	ASP	ASN	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	GLY	

GLN
LYS
TYR
LEU
SER
ASN
SER
LEU
GLU
VAL
LEU
PHE
GLN

● Molecule 1: Pannexin-2, Soluble cytochrome b562 fusion



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C7	Depositor
Number of particles used	25191	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	49	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2400	Depositor
Magnification	Not provided	
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	5.438	Depositor
Minimum map value	-2.889	Depositor
Average map value	0.019	Depositor
Map value standard deviation	0.154	Depositor
Recommended contour level	0.86	Depositor
Map size ($\text{\AA}$)	252.0, 252.0, 252.0	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.9, 0.9, 0.9	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	0.10	0/2327	0.28	0/3164
1	B	0.11	0/2327	0.28	0/3164
1	C	0.11	0/2327	0.28	0/3164
1	D	0.11	0/2327	0.29	0/3164
1	E	0.11	0/2327	0.29	0/3164
1	F	0.10	0/2327	0.30	0/3164
1	G	0.10	0/2327	0.24	0/3164
All	All	0.11	0/16289	0.28	0/22148

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	2275	0	2284	32	0
1	B	2275	0	2285	28	0
1	C	2275	0	2285	22	0
1	D	2275	0	2285	21	0
1	E	2275	0	2285	22	0
1	F	2275	0	2285	19	0
1	G	2275	0	2285	15	0
All	All	15925	0	15994	158	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 (158) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:335:GLN:CB	1:A:338:ARG:NH2	1.68	1.56
1:A:335:GLN:HB3	1:A:338:ARG:NH2	1.11	1.38
1:A:335:GLN:HA	1:A:338:ARG:NE	1.45	1.29
1:A:335:GLN:CA	1:A:338:ARG:HE	1.69	1.05
1:B:334:ARG:HD2	1:B:338:ARG:NH1	1.69	1.05
1:A:334:ARG:O	1:A:338:ARG:HG3	1.58	1.03
1:A:335:GLN:HB2	1:A:338:ARG:NH2	1.74	0.98
1:A:335:GLN:HB3	1:A:338:ARG:HH22	1.35	0.92
1:A:335:GLN:CA	1:A:338:ARG:NE	2.32	0.90
1:B:334:ARG:HD2	1:B:338:ARG:HH12	1.33	0.89
1:A:335:GLN:CA	1:A:338:ARG:NH2	2.36	0.88
1:A:81:CYS:HB3	1:A:279:CYS:HA	1.60	0.84
1:B:334:ARG:CD	1:B:338:ARG:NH1	2.42	0.82
1:A:335:GLN:HA	1:A:338:ARG:CZ	2.11	0.80
1:C:81:CYS:HB3	1:C:279:CYS:HA	1.65	0.77
1:A:335:GLN:HA	1:A:338:ARG:HE	0.74	0.76
1:F:81:CYS:HB3	1:F:279:CYS:HA	1.66	0.76
1:E:81:CYS:HB3	1:E:279:CYS:HA	1.68	0.76
1:D:77:GLU:HB2	1:D:80:TYR:HE1	1.50	0.75
1:F:77:GLU:HB2	1:F:80:TYR:HE1	1.50	0.74
1:A:335:GLN:CA	1:A:338:ARG:CZ	2.66	0.71
1:E:77:GLU:HB2	1:E:80:TYR:HE1	1.53	0.71
1:B:81:CYS:HB3	1:B:279:CYS:HA	1.72	0.71
1:B:77:GLU:HB2	1:B:80:TYR:HE1	1.57	0.70
1:D:81:CYS:HB3	1:D:279:CYS:HA	1.74	0.70
1:B:334:ARG:CD	1:B:338:ARG:HH12	2.03	0.66
1:C:77:GLU:HB2	1:C:80:TYR:HE1	1.61	0.66
1:D:253:LYS:HA	1:D:288:ARG:HH22	1.61	0.65
1:F:253:LYS:HA	1:F:288:ARG:HH22	1.60	0.65
1:F:333:ARG:HB3	1:F:337:ARG:HH22	1.63	0.64
1:B:252:GLN:O	1:B:288:ARG:NH2	2.31	0.64
1:E:253:LYS:HA	1:E:288:ARG:HH22	1.64	0.62
1:G:253:LYS:HA	1:G:288:ARG:HH22	1.64	0.62
1:D:329:GLY:O	1:D:333:ARG:NH1	2.33	0.61
1:D:65:LEU:HD23	1:D:290:ILE:HG23	1.83	0.60
1:B:334:ARG:HD2	1:B:338:ARG:CZ	2.31	0.60
1:C:65:LEU:HD23	1:C:290:ILE:HG23	1.83	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:253:LYS:HA	1:C:288:ARG:HH22	1.67	0.60
1:F:120:GLU:OE2	1:F:252:GLN:NE2	2.33	0.60
1:A:65:LEU:HD23	1:A:290:ILE:HG23	1.85	0.59
1:A:335:GLN:CB	1:A:338:ARG:CZ	2.70	0.59
1:C:252:GLN:O	1:C:288:ARG:NH2	2.35	0.59
1:A:253:LYS:HA	1:A:288:ARG:HH22	1.67	0.59
1:F:330:ILE:O	1:F:334:ARG:NH1	2.36	0.58
1:C:120:GLU:OE1	1:C:252:GLN:NE2	2.36	0.58
1:E:65:LEU:HD23	1:E:290:ILE:HG23	1.85	0.58
1:A:77:GLU:HB2	1:A:80:TYR:HE1	1.69	0.57
1:G:219:ARG:NH1	1:G:335:GLN:OE1	2.36	0.57
1:F:65:LEU:HD23	1:F:290:ILE:HG23	1.85	0.57
1:A:252:GLN:O	1:A:288:ARG:NH2	2.37	0.57
1:B:253:LYS:HA	1:B:288:ARG:HH22	1.68	0.57
1:A:131:PHE:HA	1:A:134:ILE:HG22	1.86	0.57
1:E:331:LYS:HA	1:E:334:ARG:HH12	1.70	0.56
1:G:65:LEU:HD23	1:G:290:ILE:HG23	1.85	0.56
1:D:131:PHE:HA	1:D:134:ILE:HG22	1.86	0.56
1:B:77:GLU:HB2	1:B:80:TYR:CE1	2.40	0.56
1:A:219:ARG:NH1	1:A:335:GLN:OE1	2.39	0.56
1:E:51:ASP:OD1	1:E:230:ARG:NH1	2.39	0.56
1:D:333:ARG:HB3	1:D:337:ARG:HH22	1.71	0.55
1:D:51:ASP:OD1	1:D:230:ARG:NH1	2.39	0.55
1:E:131:PHE:HA	1:E:134:ILE:HG22	1.88	0.55
1:B:157:LEU:O	1:B:161:ILE:HG12	2.06	0.55
1:C:131:PHE:HA	1:C:134:ILE:HG22	1.87	0.55
1:G:331:LYS:HA	1:G:334:ARG:HH12	1.72	0.54
1:E:342:CYS:SG	1:E:344:ILE:HG22	2.48	0.54
1:B:65:LEU:HD23	1:B:290:ILE:HG23	1.89	0.54
1:A:77:GLU:HB2	1:A:80:TYR:CE1	2.43	0.54
1:G:131:PHE:HA	1:G:134:ILE:HG22	1.90	0.54
1:A:120:GLU:OE1	1:A:252:GLN:NE2	2.42	0.53
1:B:131:PHE:HA	1:B:134:ILE:HG22	1.88	0.53
1:E:81:CYS:CB	1:E:279:CYS:HA	2.38	0.53
1:B:219:ARG:NH1	1:B:335:GLN:OE1	2.41	0.53
1:B:51:ASP:OD1	1:B:230:ARG:NH1	2.41	0.52
1:F:154:LEU:HG	1:F:158:LEU:HD13	1.92	0.52
1:E:154:LEU:HD21	1:E:354:ASN:HD22	1.75	0.52
1:B:81:CYS:CB	1:B:279:CYS:HA	2.39	0.52
1:A:51:ASP:OD1	1:A:230:ARG:NH1	2.42	0.52
1:C:51:ASP:OD1	1:C:230:ARG:NH1	2.42	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:252:GLN:O	1:F:288:ARG:NH2	2.42	0.52
1:A:157:LEU:O	1:A:161:ILE:HG12	2.10	0.51
1:G:51:ASP:OD2	1:G:230:ARG:NH1	2.43	0.51
1:D:153:GLU:OE2	1:D:213:TYR:OH	2.24	0.51
1:D:157:LEU:O	1:D:161:ILE:HG12	2.11	0.51
1:C:157:LEU:O	1:C:161:ILE:HG12	2.10	0.51
1:F:51:ASP:OD1	1:F:230:ARG:NH1	2.43	0.50
1:D:81:CYS:CB	1:D:279:CYS:HA	2.40	0.50
1:C:331:LYS:HA	1:C:334:ARG:NH1	2.28	0.49
1:G:157:LEU:O	1:G:161:ILE:HG12	2.12	0.49
1:A:133:ALA:O	1:A:137:VAL:HG23	2.12	0.49
1:A:81:CYS:CB	1:A:279:CYS:HA	2.38	0.49
1:B:133:ALA:O	1:B:137:VAL:HG23	2.13	0.48
1:G:133:ALA:O	1:G:137:VAL:HG23	2.13	0.48
1:C:77:GLU:HB2	1:C:80:TYR:CE1	2.44	0.48
1:B:342:CYS:SG	1:B:344:ILE:HG22	2.52	0.48
1:D:133:ALA:O	1:D:137:VAL:HG23	2.13	0.48
1:E:105:ASP:OD1	1:E:106:ALA:N	2.44	0.48
1:E:77:GLU:HB2	1:E:80:TYR:CE1	2.42	0.48
1:G:105:ASP:OD1	1:G:106:ALA:N	2.44	0.48
1:B:149:ARG:O	1:B:153:GLU:HG2	2.14	0.48
1:E:133:ALA:O	1:E:137:VAL:HG23	2.14	0.48
1:C:133:ALA:O	1:C:137:VAL:HG23	2.14	0.47
1:G:72:LYS:HG2	1:G:286:LEU:HD22	1.96	0.47
1:D:312:LEU:HD21	1:D:349:MET:HE1	1.96	0.47
1:C:331:LYS:HA	1:C:334:ARG:HH12	1.80	0.47
1:B:154:LEU:HD21	1:B:354:ASN:HD22	1.79	0.47
1:D:252:GLN:O	1:D:288:ARG:NH2	2.48	0.46
1:F:330:ILE:HB	1:F:334:ARG:HH12	1.81	0.45
1:D:328:VAL:HG23	1:D:333:ARG:NH1	2.32	0.45
1:F:105:ASP:OD1	1:F:106:ALA:N	2.45	0.45
1:D:331:LYS:HA	1:D:334:ARG:HH22	1.82	0.45
1:E:219:ARG:NE	1:E:219:ARG:HA	2.32	0.45
1:F:81:CYS:CB	1:F:279:CYS:HA	2.40	0.45
1:C:154:LEU:HD21	1:C:354:ASN:HD22	1.81	0.45
1:B:317:LYS:HA	1:B:317:LYS:HE2	1.99	0.45
1:D:111:ASP:N	1:D:111:ASP:OD1	2.50	0.45
1:A:111:ASP:N	1:A:111:ASP:OD1	2.50	0.45
1:B:72:LYS:HG2	1:B:286:LEU:HD22	1.99	0.45
1:B:111:ASP:OD1	1:B:111:ASP:N	2.50	0.44
1:C:111:ASP:N	1:C:111:ASP:OD1	2.50	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:72:LYS:HG2	1:A:286:LEU:HD22	1.99	0.44
1:A:317:LYS:HA	1:A:317:LYS:HE2	2.00	0.44
1:A:334:ARG:O	1:A:338:ARG:CG	2.46	0.44
1:F:324:LYS:HD3	1:F:324:LYS:HA	1.77	0.44
1:G:312:LEU:HD12	1:G:316:ARG:HE	1.83	0.44
1:G:317:LYS:HA	1:G:317:LYS:HE2	1.99	0.44
1:C:81:CYS:CB	1:C:279:CYS:HA	2.41	0.44
1:C:317:LYS:HA	1:C:317:LYS:HE2	2.00	0.44
1:G:111:ASP:N	1:G:111:ASP:OD1	2.50	0.44
1:G:236:LEU:HD23	1:G:236:LEU:HA	1.86	0.44
1:D:82:TYR:CE2	1:E:265:PRO:HG2	2.53	0.43
1:F:111:ASP:OD1	1:F:111:ASP:N	2.50	0.43
1:F:317:LYS:HA	1:F:317:LYS:HE2	1.99	0.43
1:B:129:LEU:HD12	1:B:129:LEU:HA	1.90	0.43
1:C:333:ARG:O	1:C:337:ARG:NH1	2.52	0.43
1:B:79:ILE:HD11	1:B:279:CYS:HB3	2.01	0.43
1:D:256:GLU:HG2	1:D:280:LYS:HD2	2.01	0.43
1:A:312:LEU:HD12	1:A:316:ARG:HE	1.83	0.43
1:B:328:VAL:HB	1:B:333:ARG:HH22	1.84	0.43
1:E:310:ILE:HG23	1:E:314:ILE:HD12	2.01	0.42
1:E:317:LYS:HA	1:E:317:LYS:HE2	2.00	0.42
1:D:317:LYS:HE2	1:D:317:LYS:HA	2.00	0.42
1:B:236:LEU:HD23	1:B:236:LEU:HA	1.86	0.42
1:B:312:LEU:HD12	1:B:316:ARG:HE	1.85	0.42
1:G:154:LEU:HD21	1:G:354:ASN:HD22	1.84	0.42
1:A:129:LEU:HD12	1:A:129:LEU:HA	1.90	0.42
1:C:312:LEU:HD12	1:C:316:ARG:HE	1.85	0.42
1:E:72:LYS:HD3	1:E:72:LYS:C	2.45	0.42
1:F:312:LEU:HD12	1:F:316:ARG:HE	1.85	0.41
1:C:236:LEU:HD23	1:C:236:LEU:HA	1.87	0.41
1:E:331:LYS:HA	1:E:334:ARG:NH1	2.33	0.41
1:F:264:SER:N	1:F:265:PRO:HD3	2.36	0.41
1:C:310:ILE:HG23	1:C:314:ILE:HD12	2.02	0.41
1:C:115:TRP:HA	1:C:116:PRO:HD3	1.98	0.41
1:E:237:LEU:HD23	1:E:237:LEU:HA	1.91	0.41
1:F:46:LEU:HG	1:F:46:LEU:O	2.21	0.40
1:E:111:ASP:OD1	1:E:111:ASP:N	2.50	0.40
1:E:330:ILE:O	1:E:334:ARG:NH2	2.54	0.40
1:D:105:ASP:OD1	1:D:106:ALA:N	2.43	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	281/487 (58%)	272 (97%)	9 (3%)	0	100	100
1	B	281/487 (58%)	271 (96%)	10 (4%)	0	100	100
1	C	281/487 (58%)	273 (97%)	8 (3%)	0	100	100
1	D	281/487 (58%)	270 (96%)	11 (4%)	0	100	100
1	E	281/487 (58%)	271 (96%)	10 (4%)	0	100	100
1	F	281/487 (58%)	273 (97%)	8 (3%)	0	100	100
1	G	281/487 (58%)	274 (98%)	7 (2%)	0	100	100
All	All	1967/3409 (58%)	1904 (97%)	63 (3%)	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	239/414 (58%)	239 (100%)	0	100	100
1	B	239/414 (58%)	239 (100%)	0	100	100
1	C	239/414 (58%)	239 (100%)	0	100	100
1	D	239/414 (58%)	239 (100%)	0	100	100
1	E	239/414 (58%)	239 (100%)	0	100	100
1	F	239/414 (58%)	239 (100%)	0	100	100
1	G	239/414 (58%)	239 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	1673/2898 (58%)	1673 (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 (26) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	121	HIS
1	A	352	ASN
1	B	121	HIS
1	B	252	GLN
1	B	352	ASN
1	B	354	ASN
1	B	357	HIS
1	C	121	HIS
1	C	155	ASN
1	C	252	GLN
1	C	352	ASN
1	D	121	HIS
1	D	155	ASN
1	D	352	ASN
1	E	121	HIS
1	E	155	ASN
1	E	285	GLN
1	E	352	ASN
1	E	354	ASN
1	F	121	HIS
1	F	285	GLN
1	F	352	ASN
1	F	357	HIS
1	G	285	GLN
1	G	352	ASN
1	G	354	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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.

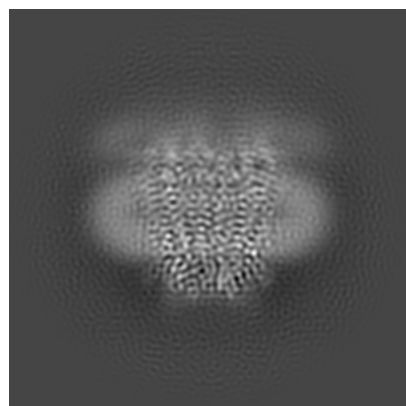
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-28902. These allow visual inspection of the internal detail of the map and identification of artifacts.

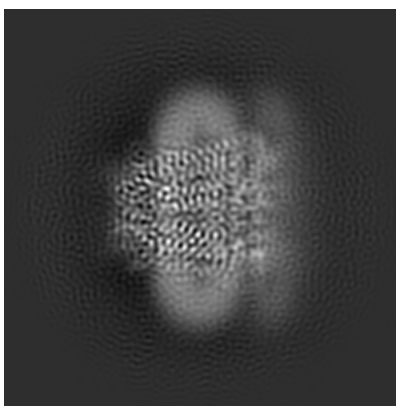
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

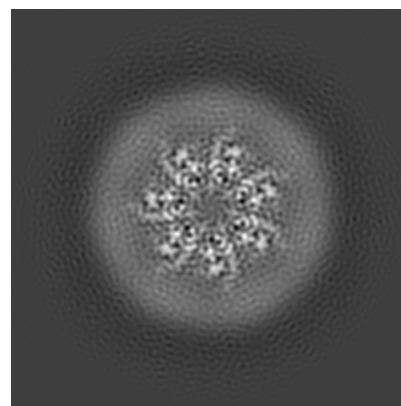
6.1.1 Primary map



X

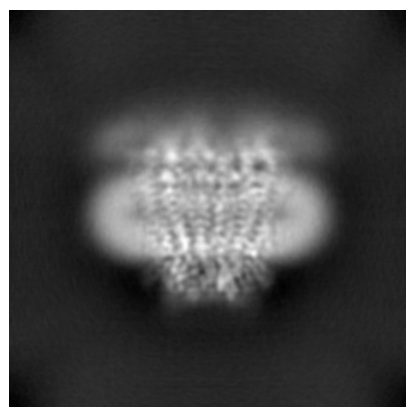


Y

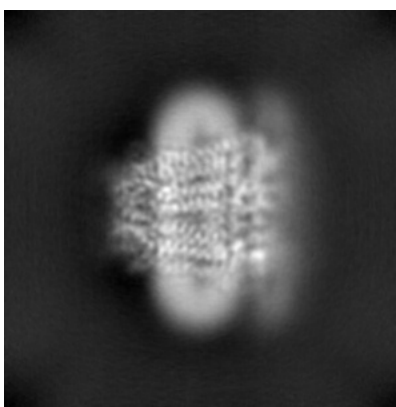


Z

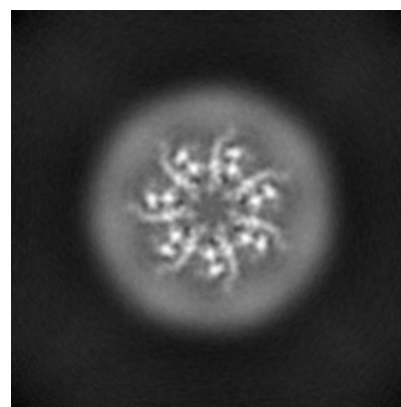
6.1.2 Raw map



X



Y

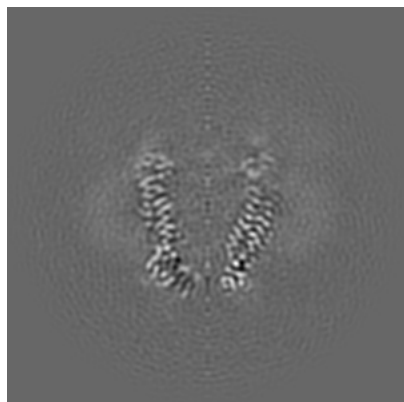


Z

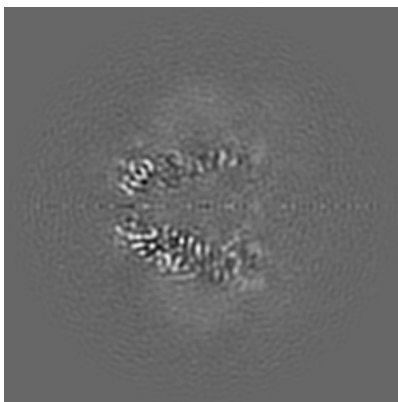
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

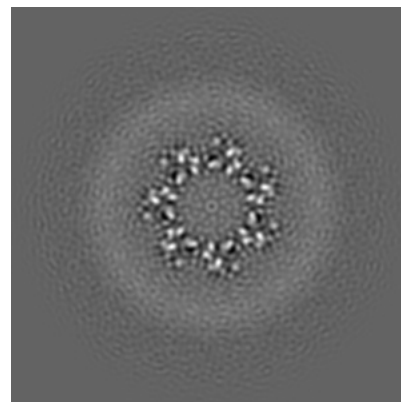
6.2.1 Primary map



X Index: 140

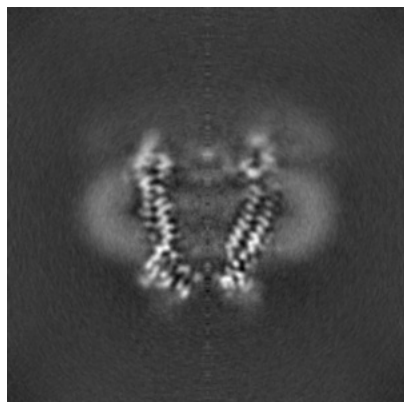


Y Index: 140

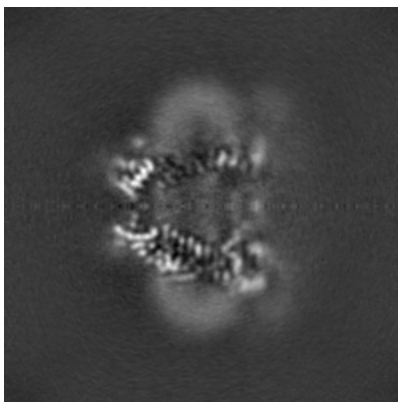


Z Index: 140

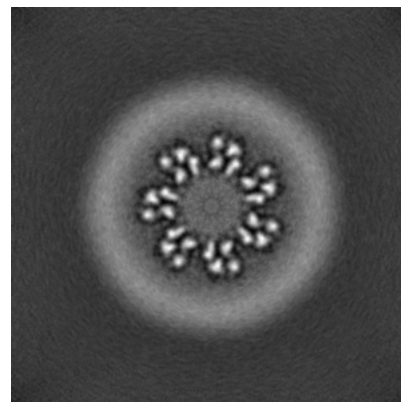
6.2.2 Raw map



X Index: 140



Y Index: 140

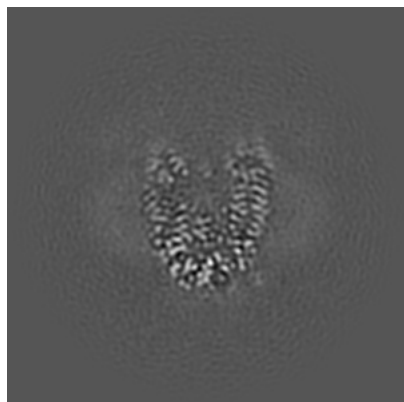


Z Index: 140

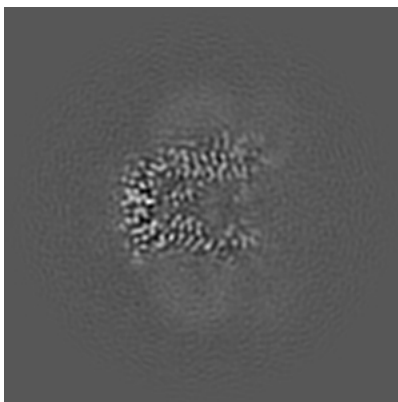
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

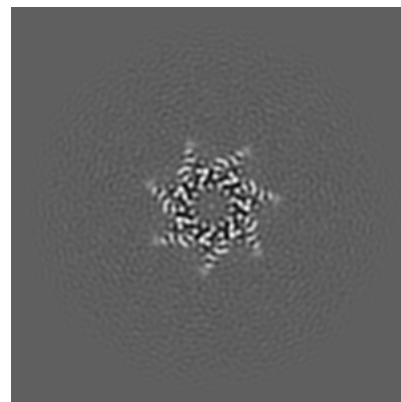
6.3.1 Primary map



X Index: 118

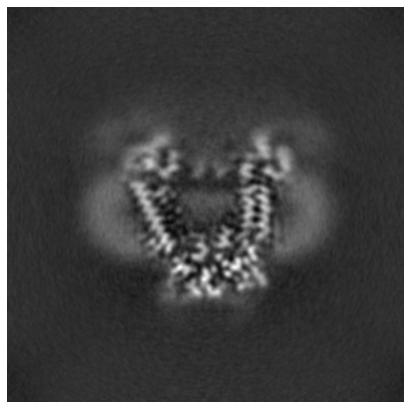


Y Index: 119

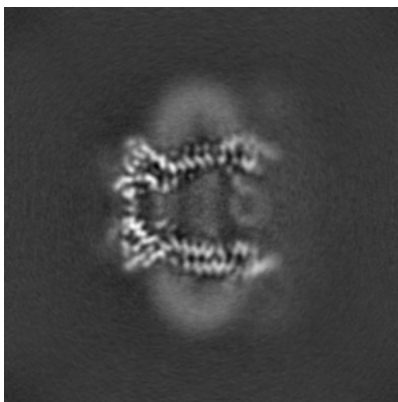


Z Index: 97

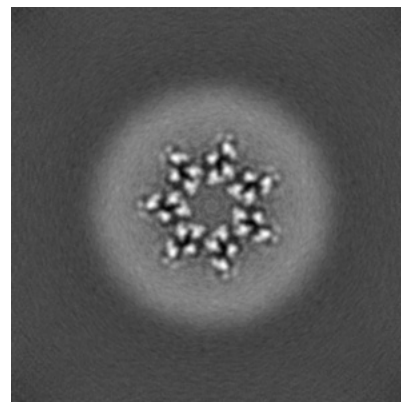
6.3.2 Raw map



X Index: 155



Y Index: 148

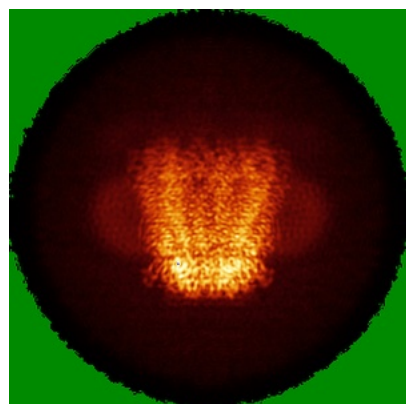


Z Index: 119

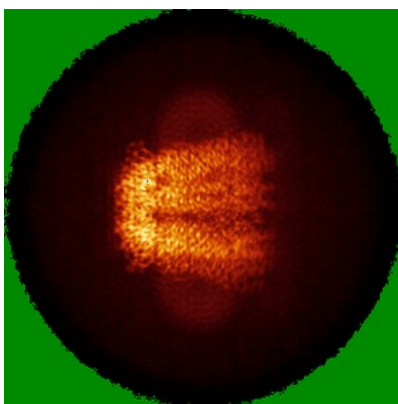
The images above show the largest variance slices of the map in three orthogonal directions.

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

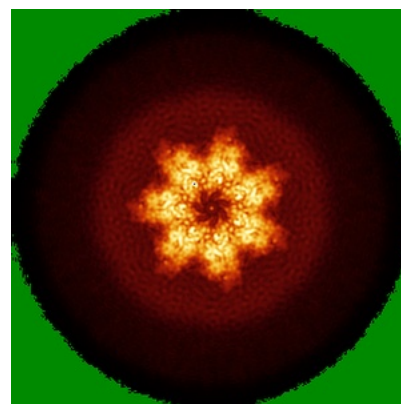
6.4.1 Primary map



X

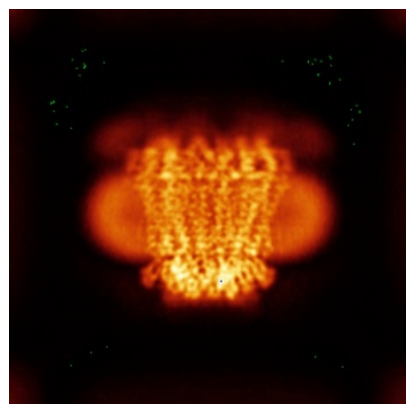


Y

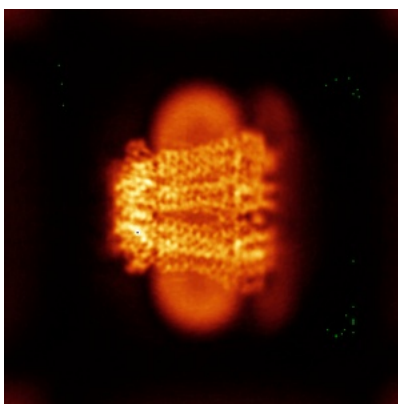


Z

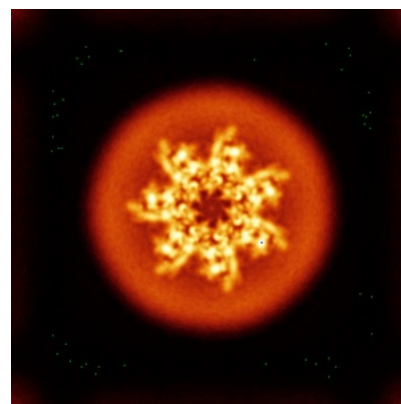
6.4.2 Raw map



X



Y

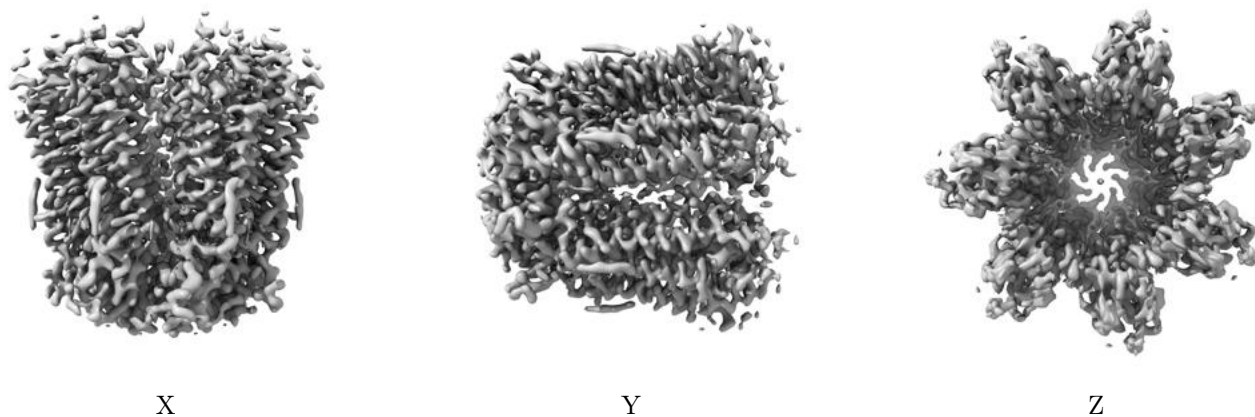


Z

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.

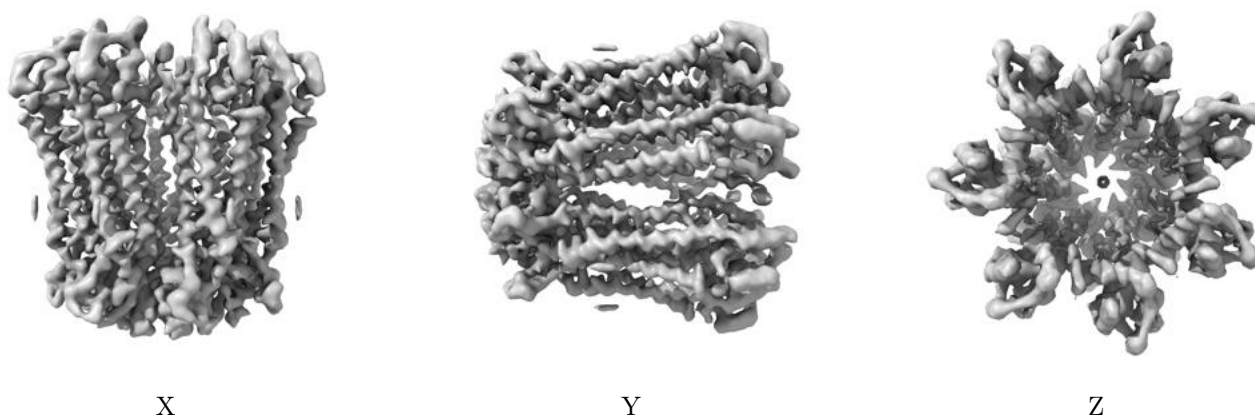
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.86. 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.

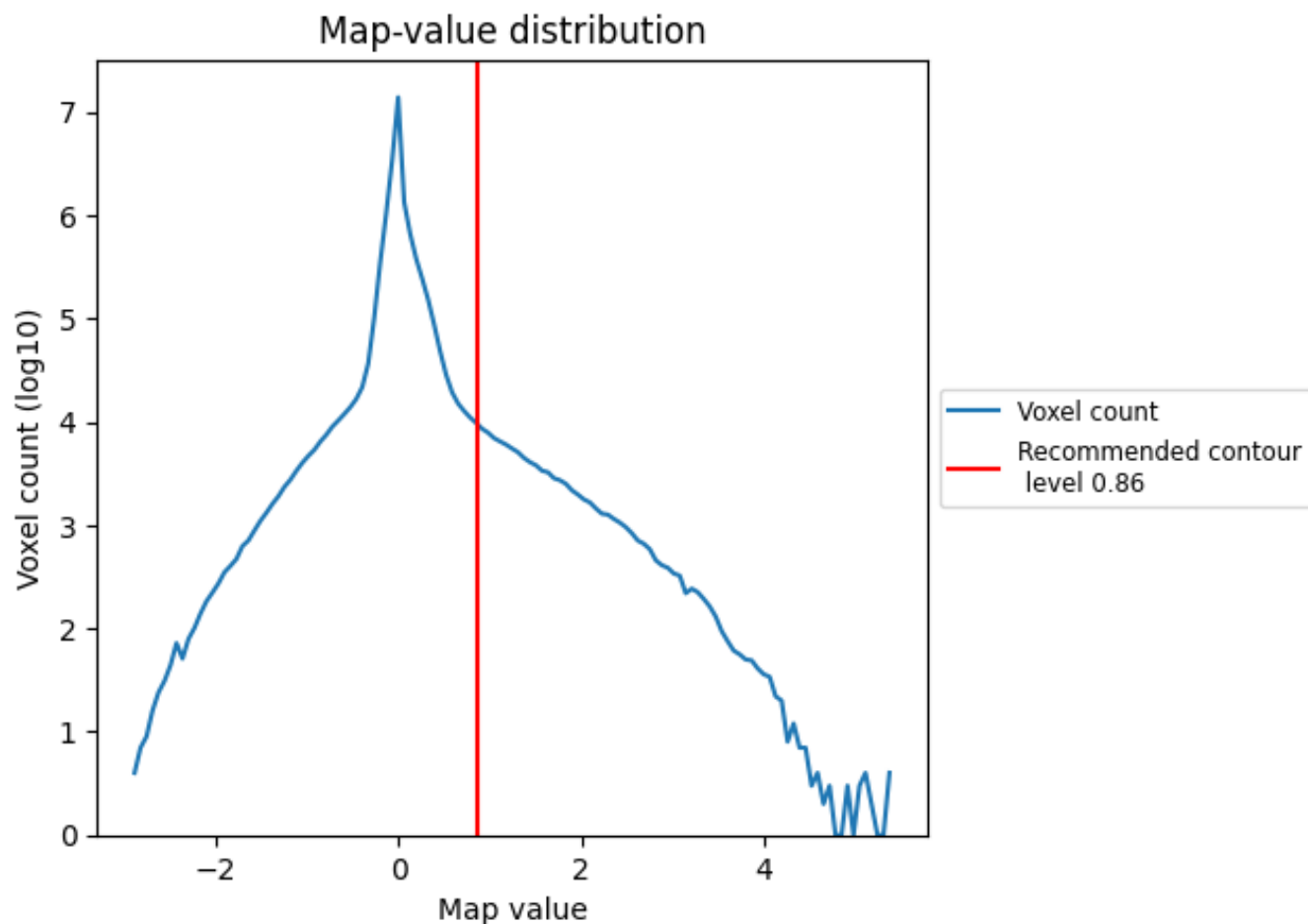
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

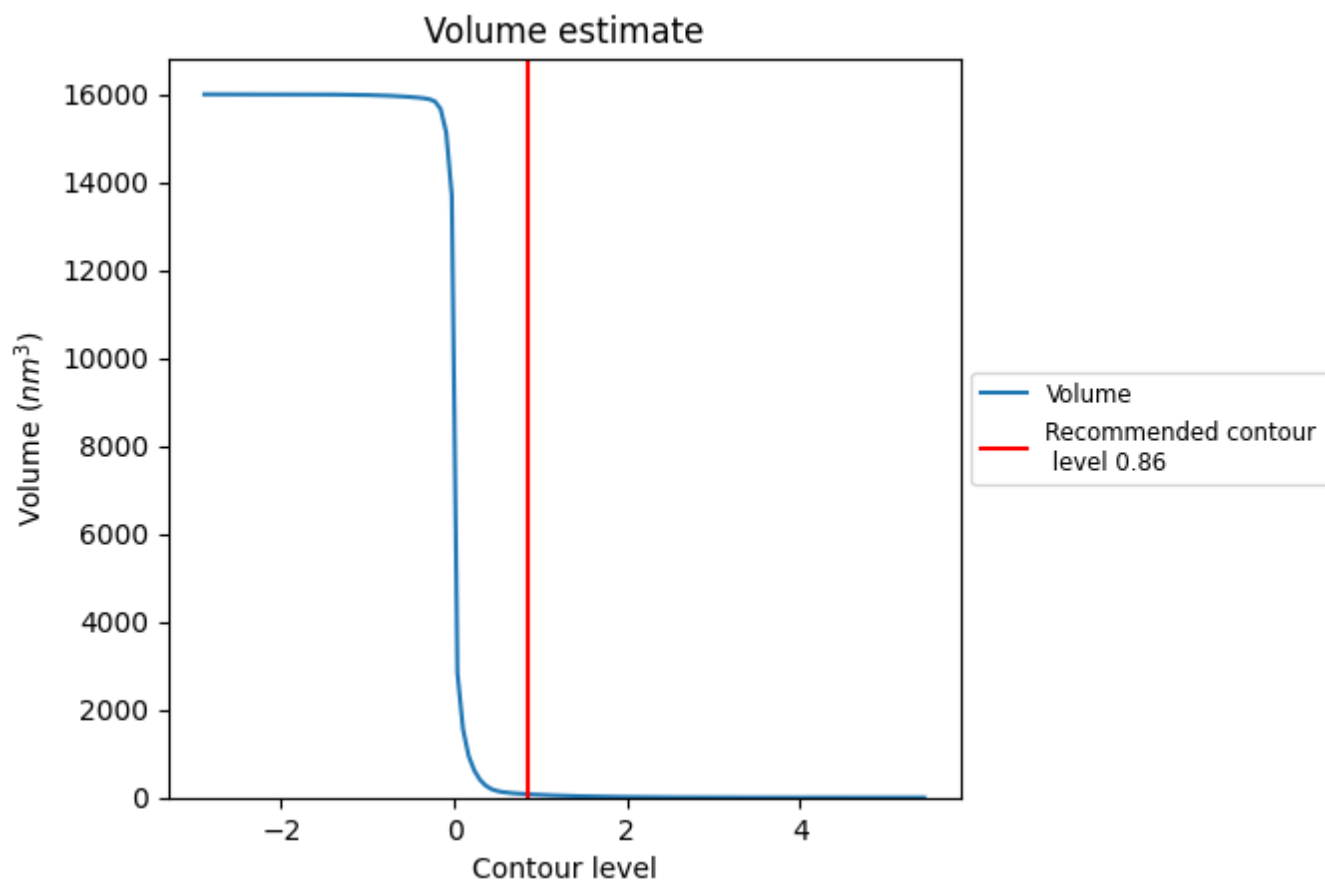
This section contains the results of statistical analysis of the map.

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.

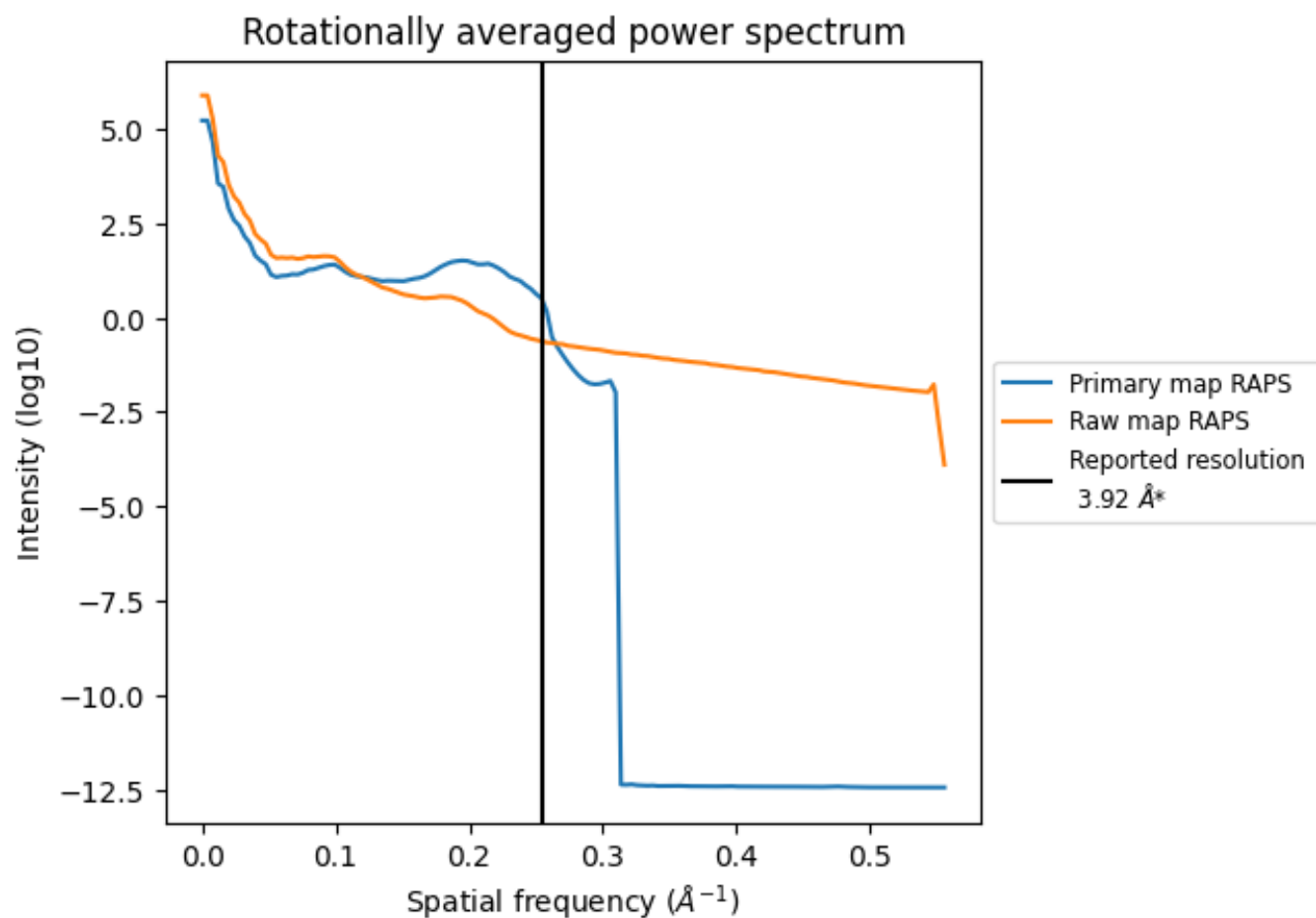
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 75 nm³; this corresponds to an approximate mass of 68 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 ⓘ

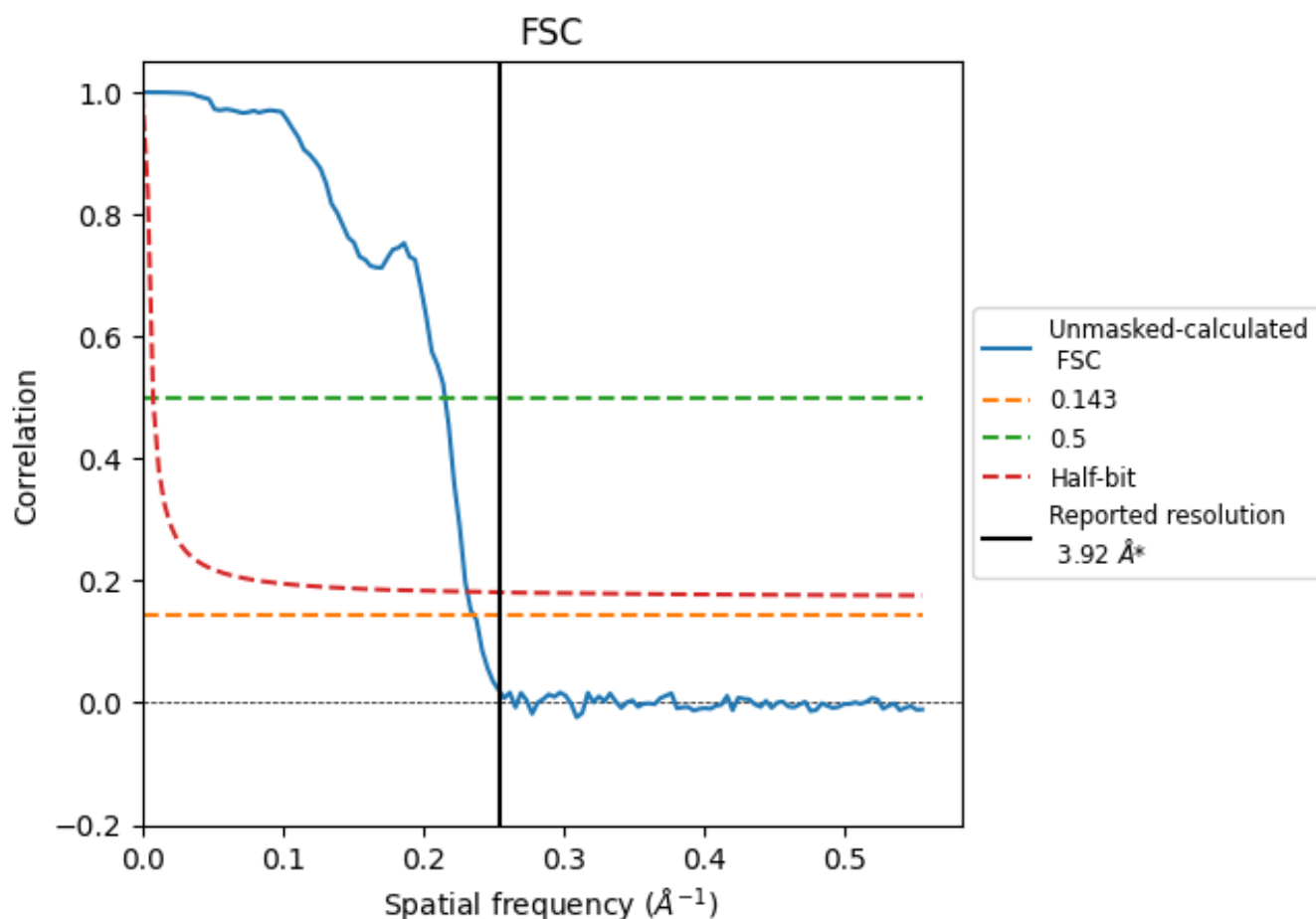


*Reported resolution corresponds to spatial frequency of 0.255 Å⁻¹

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 0.255 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

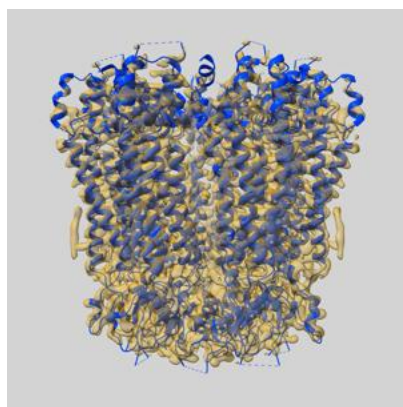
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.92	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	4.23	4.64	4.32

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

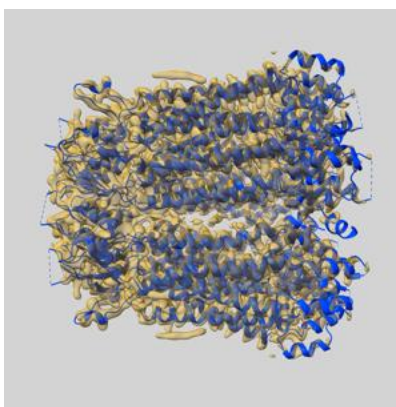
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-28902 and PDB model 8F7C. Per-residue inclusion information can be found in section [3](#) on page [7](#).

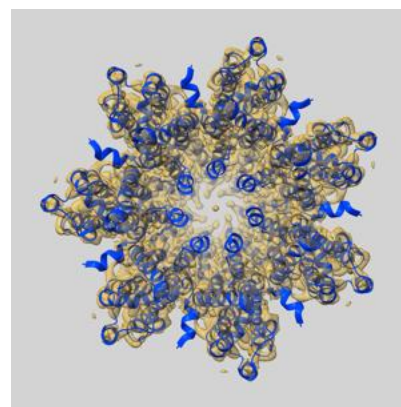
9.1 Map-model overlay [i](#)



X



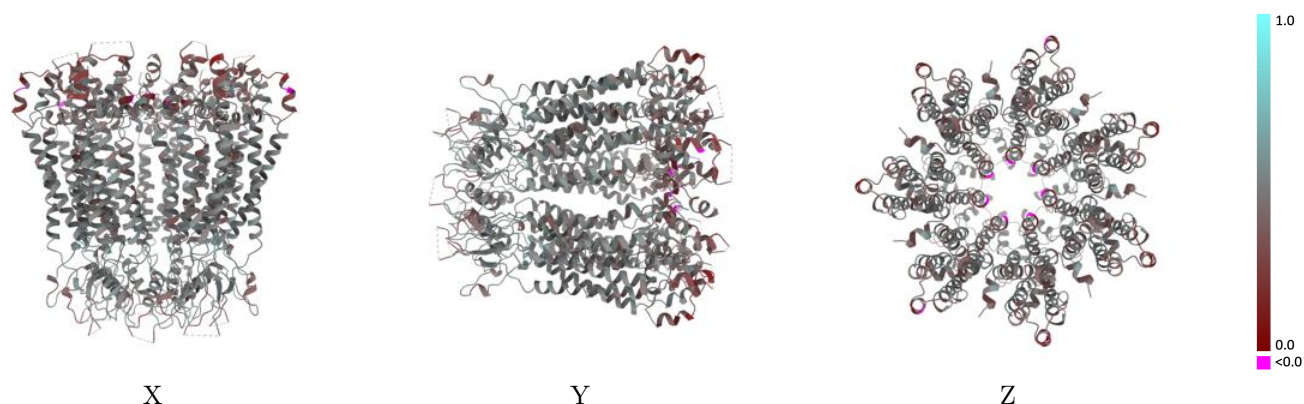
Y



Z

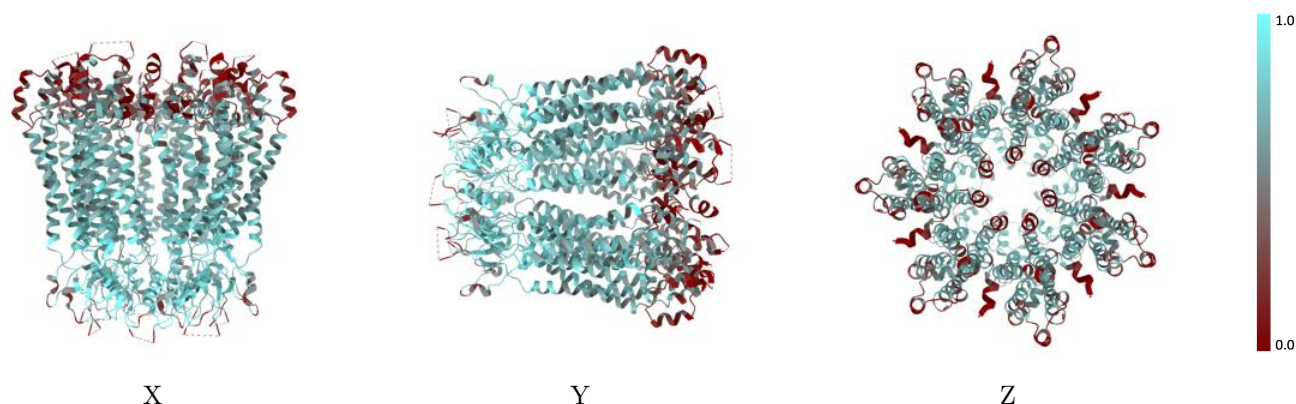
The images above show the 3D surface view of the map at the recommended contour level 0.86 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



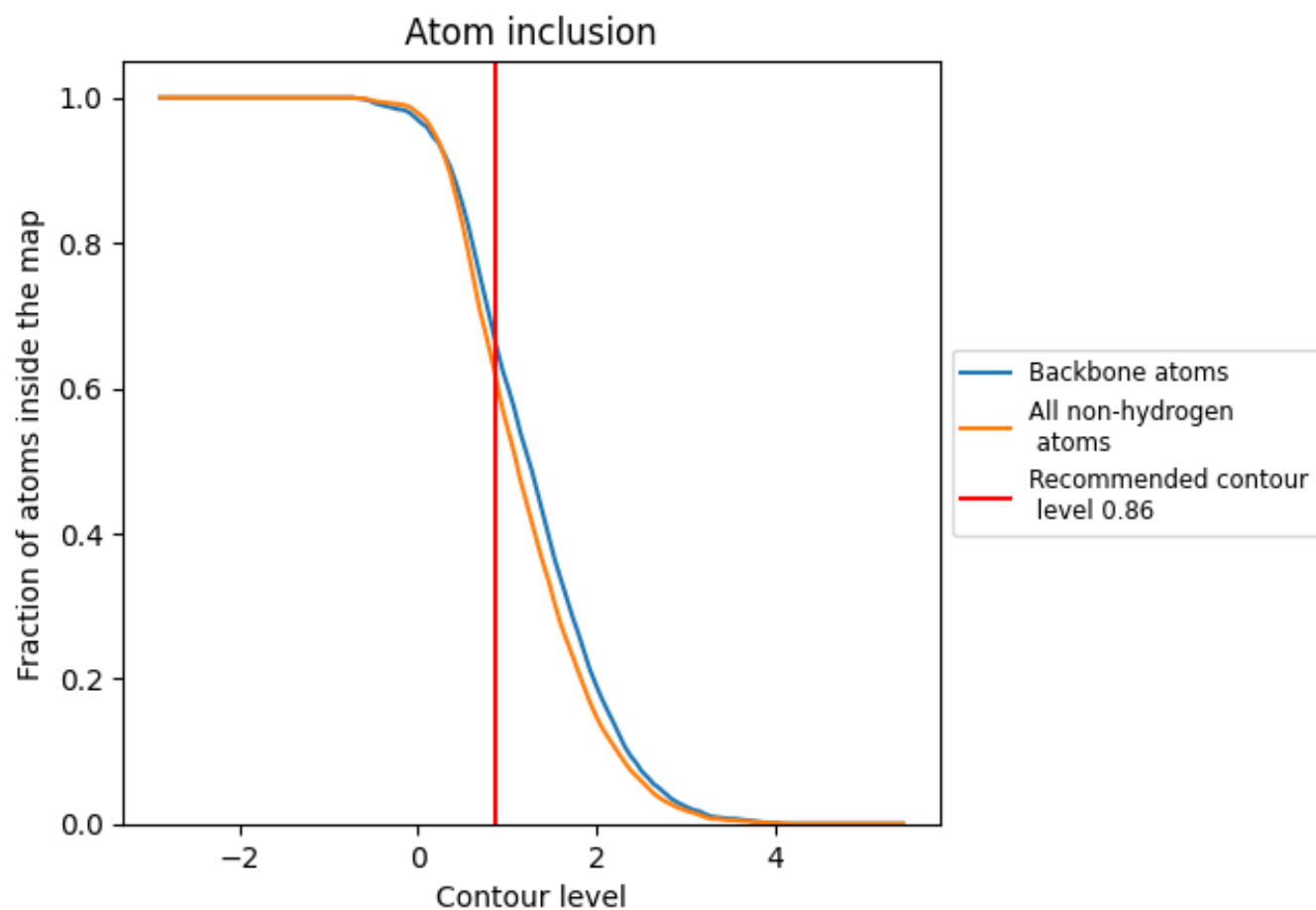
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions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

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



The images above show the model with each residue coloured according to its atom inclusion. This shows to what extent they are inside the map at the recommended contour level (0.86).

9.4 Atom inclusion [i](#)



At the recommended contour level, 67% of all backbone atoms, 62% 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 (0.86) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6220	0.4580
A	0.6210	0.4580
B	0.6210	0.4570
C	0.6180	0.4600
D	0.6210	0.4580
E	0.6250	0.4570
F	0.6280	0.4620
G	0.6210	0.4580

1.0

0.0

<0.0