



wwPDB EM Validation Summary Report ⓘ

Mar 8, 2026 – 10:46 PM UTC

PDB ID : 8CTG / pdb_00008ctg
EMDB ID : EMD-26989
Title : Extracellular architecture of an engineered canonical Wnt signaling ternary complex
Authors : Tsutsumi, N.; Jude, K.M.; Garcia, K.C.
Deposited on : 2022-05-14
Resolution : 3.80 Å(reported)

This is a wwPDB EM Validation Summary 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
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
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

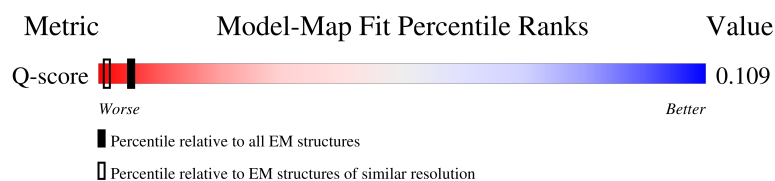
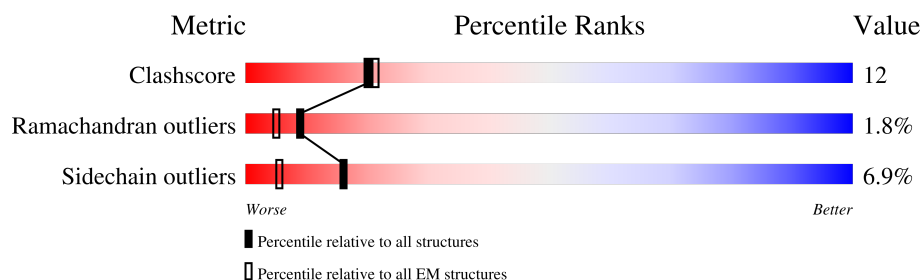
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 3.80 Å.

Percentile scores (ranging between 0-100) for global validation metrics of the entry are shown in the following graphic. The table shows the number of entries on which the scores are based.



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	10198 (3.30 - 4.30)

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	132	
2	B	327	
3	C	649	

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5100 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 Frizzled-8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	118	618	372	118	118	10	0	0

There are 9 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	151	ALA	-	expression tag	UNP Q61091
A	152	ALA	-	expression tag	UNP Q61091
A	153	ALA	-	expression tag	UNP Q61091
A	154	LEU	-	expression tag	UNP Q61091
A	155	GLU	-	expression tag	UNP Q61091
A	156	VAL	-	expression tag	UNP Q61091
A	157	LEU	-	expression tag	UNP Q61091
A	158	PHE	-	expression tag	UNP Q61091
A	159	GLN	-	expression tag	UNP Q61091

- Molecule 2 is a protein called Protein Wnt-8.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	316	1585	930	316	317	22	0	0

There are 25 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	21	GLY	-	expression tag	UNP P28026
B	227	PRO	ARG	conflict	UNP P28026
B	229	THR	GLY	conflict	UNP P28026
B	232	VAL	ALA	conflict	UNP P28026
B	233	ASN	ASP	conflict	UNP P28026
B	234	SER	ASN	conflict	UNP P28026
B	236	ARG	GLY	conflict	UNP P28026
B	330	GLN	-	expression tag	UNP P28026

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Chain	Residue	Modelled	Actual	Comment	Reference
B	331	VAL	-	expression tag	UNP P28026
B	332	VAL	-	expression tag	UNP P28026
B	333	ILE	-	expression tag	UNP P28026
B	334	LYS	-	expression tag	UNP P28026
B	335	HIS	-	expression tag	UNP P28026
B	336	PHE	-	expression tag	UNP P28026
B	337	CYS	-	expression tag	UNP P28026
B	338	ALA	-	expression tag	UNP P28026
B	339	ARG	-	expression tag	UNP P28026
B	340	HIS	-	expression tag	UNP P28026
B	341	HIS	-	expression tag	UNP P28026
B	342	HIS	-	expression tag	UNP P28026
B	343	HIS	-	expression tag	UNP P28026
B	344	HIS	-	expression tag	UNP P28026
B	345	HIS	-	expression tag	UNP P28026
B	346	HIS	-	expression tag	UNP P28026
B	347	HIS	-	expression tag	UNP P28026

- Molecule 3 is a protein called Low-density lipoprotein receptor-related protein 6.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	572	Total	C	N	O	S	0	0
			2880	1723	572	572	13		

There are 39 discrepancies between the modelled and reference sequences:

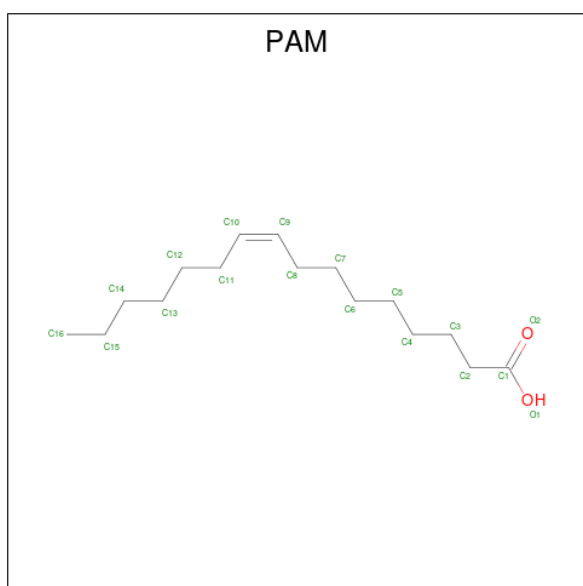
Chain	Residue	Modelled	Actual	Comment	Reference
C	18	GLY	-	expression tag	UNP O75581
C	19	SER	-	expression tag	UNP O75581
C	630	SER	-	expression tag	UNP O75581
C	631	ARG	-	expression tag	UNP O75581
C	632	GLY	-	expression tag	UNP O75581
C	633	LEU	-	expression tag	UNP O75581
C	634	GLU	-	expression tag	UNP O75581
C	635	VAL	-	expression tag	UNP O75581
C	636	LEU	-	expression tag	UNP O75581
C	637	PHE	-	expression tag	UNP O75581
C	638	GLN	-	expression tag	UNP O75581
C	639	GLY	-	expression tag	UNP O75581
C	640	PRO	-	expression tag	UNP O75581
C	641	GLY	-	expression tag	UNP O75581
C	642	ALA	-	expression tag	UNP O75581

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Chain	Residue	Modelled	Actual	Comment	Reference
C	643	ALA	-	expression tag	UNP O75581
C	644	GLY	-	expression tag	UNP O75581
C	645	LEU	-	expression tag	UNP O75581
C	646	ASN	-	expression tag	UNP O75581
C	647	ASP	-	expression tag	UNP O75581
C	648	ILE	-	expression tag	UNP O75581
C	649	PHE	-	expression tag	UNP O75581
C	650	GLU	-	expression tag	UNP O75581
C	651	ALA	-	expression tag	UNP O75581
C	652	GLN	-	expression tag	UNP O75581
C	653	LYS	-	expression tag	UNP O75581
C	654	ILE	-	expression tag	UNP O75581
C	655	GLU	-	expression tag	UNP O75581
C	656	TRP	-	expression tag	UNP O75581
C	657	HIS	-	expression tag	UNP O75581
C	658	GLU	-	expression tag	UNP O75581
C	659	HIS	-	expression tag	UNP O75581
C	660	HIS	-	expression tag	UNP O75581
C	661	HIS	-	expression tag	UNP O75581
C	662	HIS	-	expression tag	UNP O75581
C	663	HIS	-	expression tag	UNP O75581
C	664	HIS	-	expression tag	UNP O75581
C	665	HIS	-	expression tag	UNP O75581
C	666	HIS	-	expression tag	UNP O75581

- Molecule 4 is PALMITOLEIC ACID (CCD ID: PAM) (formula: $C_{16}H_{30}O_2$).

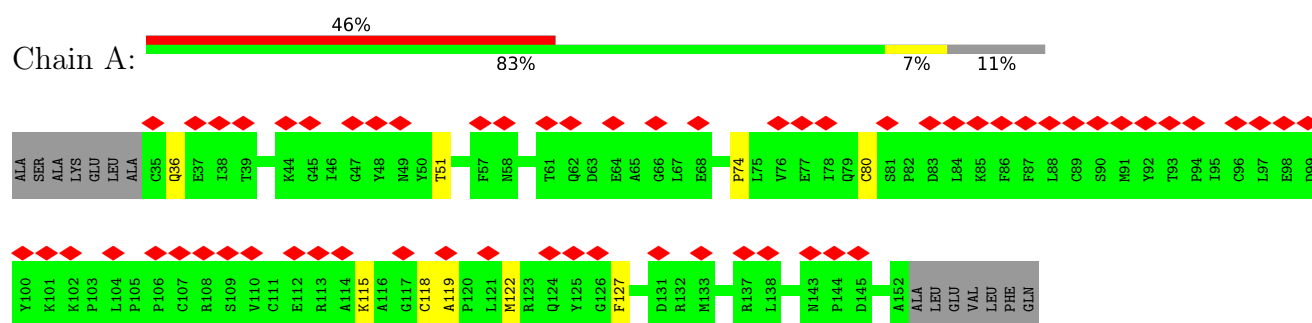


Mol	Chain	Residues	Atoms			AltConf
			Total	C	O	
4	B	1	17	16	1	0

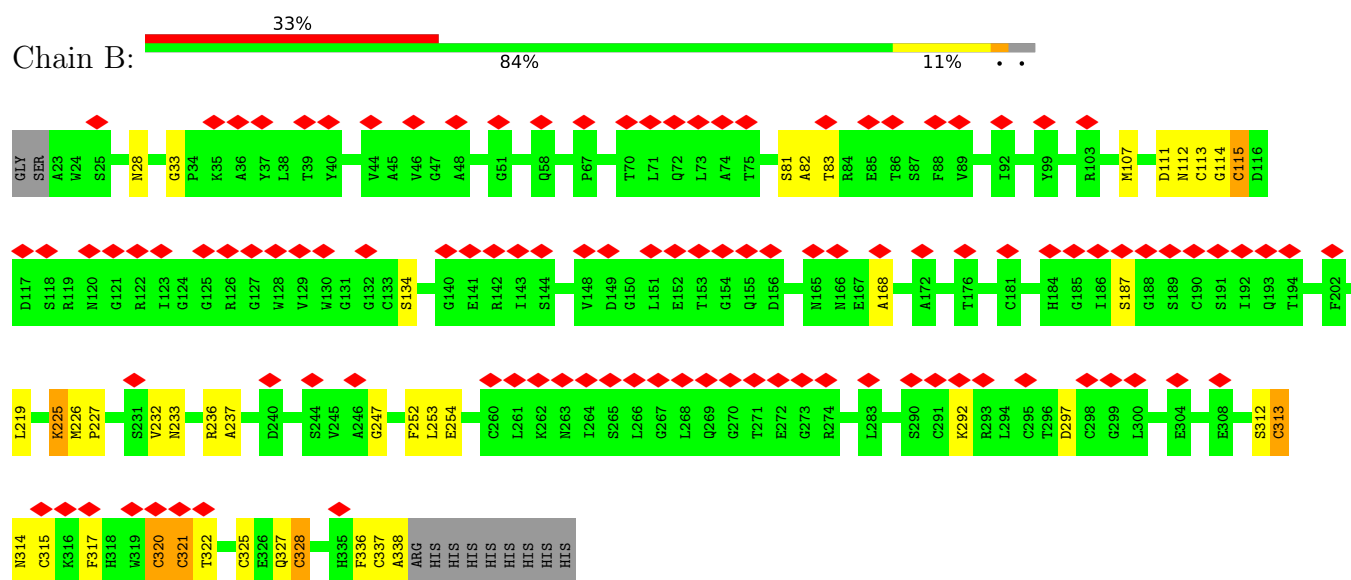
3 Residue-property plots

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

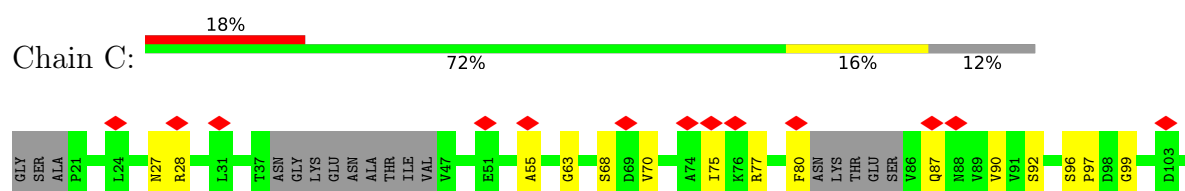
• Molecule 1: Frizzled-8



• Molecule 2: Protein Wnt-8



• Molecule 3: Low-density lipoprotein receptor-related protein 6



GLY	ALA	ALA	GLY	LEU	ASN	ASP	ILE	PHE	GLU	ALA	ALA	GLN	LYS	ILE	GLU	TRP	HIS	GLU	HIS	HIS	HIS	HIS	HIS	HIS	HIS	HIS	GLY	PRO																						
Y546	W547	T548	D549	R552	R553	S562	E563	E564	R565	E566	V567	I568	P573	G577	I587	G588	A593	E594	S600	C603	R606	R611	P615	ILE	GLY	PHE	GLU	LEU	ILE	SER	ASP	MET	LYS	THR	CYS	ILE	VAL	SER	ARG	GLY	LEU	GLU	VAL	LEU	PHE	GLN	GLY	PRO		
D421	T422	D425	E428	V429	T430	K438	D444	L445	E446	F447	P448	Y459	T463	D464	I468	P469	E472	V485	N486	W491	L495	A509	D512	K513	I514	E515	V516	M517	D520	G521	T522	G523	R524	R525	V526	D530	P533	H534	I535	F536	G537	F538								
G291	G292	C293	K303	P304	Q307	G310	P311	T312	G313	C323	G326	E329	L332	R336	R340	P347	D348	E357	H361	I374	Y375	D379	R385	R386	G393	S394	Q395	A400	Q401	H404	P405	D406	G407	I408	A409	A414	Y418	W419	T420											
A201	K202	L203	N204	I206	H207	K208	S209	W210	L211	D212	N215	R216	Q217	ALA	VAL	VAL	LYS	GLY	SER	LEU	PRO	HIS	P227	F228	A229	L232	D235	S246	C250	T254	G255	E256	G257	H262	F266	S267	P268	M269	D270	A273	F274	R278	N281	A282	C286					
Y110	W111	T112	D113	S114	R118	I119	E120	V121	S127	K130	V131	L132	D138	Q139	P140	R141	A142	I143	A144	L145	D146	G150	F151	M152	Y153	W154	T155	D156	W157	G158	E159	W160	P161	R165	N178	W183	P184	R185	G186	L187	T188	L189	D190	Q194	K195	L196	Y197	W198	A199	D200

4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	82235	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	55	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	58680	Depositor
Image detector	GATAN K3 (6k x 4k)	Depositor
Maximum map value	5.127	Depositor
Minimum map value	-0.001	Depositor
Average map value	0.002	Depositor
Map value standard deviation	0.035	Depositor
Recommended contour level	0.2	Depositor
Map size (Å)	299.9392, 299.9392, 299.9392	wwPDB
Map dimensions	352, 352, 352	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.8521, 0.8521, 0.8521	Depositor

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: PAM

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

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	$\# Z > 5$	RMSZ	$\# Z > 5$
1	A	0.23	0/629	0.55	0/881
2	B	0.21	0/1588	0.61	0/2200
3	C	0.22	0/2901	0.67	0/4042
All	All	0.22	0/5118	0.64	0/7123

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	618	0	342	6	0
2	B	1585	0	789	26	0
3	C	2880	0	1444	65	0
4	B	17	0	29	1	0
All	All	5100	0	2604	95	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 12.

The worst 5 of 95 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)
3:C:196:LEU:O	3:C:209:SER:N	2.11	0.82
2:B:312:SER:HA	2:B:327:GLN:HA	1.62	0.79
2:B:219:LEU:H	2:B:253:LEU:HA	1.49	0.78
3:C:422:THR:HA	3:C:448:PRO:HD2	1.64	0.78
3:C:374:ILE:O	3:C:386:ARG:HA	1.84	0.77

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	116/132 (88%)	113 (97%)	3 (3%)	0	100	100
2	B	314/327 (96%)	275 (88%)	32 (10%)	7 (2%)	5	30
3	C	564/649 (87%)	492 (87%)	61 (11%)	11 (2%)	6	32
All	All	994/1108 (90%)	880 (88%)	96 (10%)	18 (2%)	9	33

5 of 18 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	B	115	CYS
2	B	236	ARG
3	C	303	LYS
3	C	562	SER
2	B	237	ALA

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	22/115 (19%)	22 (100%)	0	100	100
2	B	27/278 (10%)	21 (78%)	6 (22%)	1	5
3	C	38/557 (7%)	38 (100%)	0	100	100
All	All	87/950 (9%)	81 (93%)	6 (7%)	16	40

5 of 6 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
2	B	320	CYS
2	B	321	CYS
2	B	328	CYS
2	B	227	PRO
2	B	187	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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](#)

1 ligand is modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. 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 $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	PAM	B	401	2	16,16,17	1.08	1 (6%)	15,15,17	0.68	1 (6%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	PAM	B	401	2	-	4/14/14/15	-

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	B	401	PAM	O1-C1	-4.21	1.20	1.42

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
4	B	401	PAM	O1-C1-C2	2.05	124.47	111.44

There are no chirality outliers.

All (4) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	B	401	PAM	C5-C6-C7-C8
4	B	401	PAM	C9-C10-C11-C12
4	B	401	PAM	C7-C8-C9-C10
4	B	401	PAM	C2-C3-C4-C5

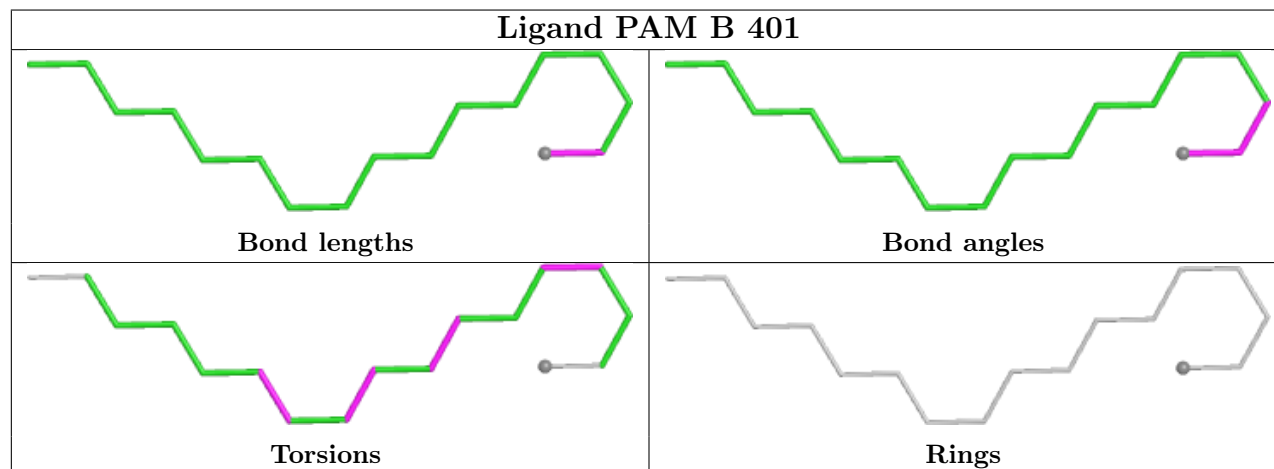
There are no ring outliers.

1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	B	401	PAM	1	0

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.



5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

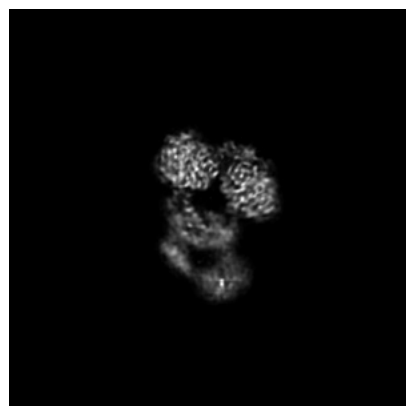
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-26989. These allow visual inspection of the internal detail of the map and identification of artifacts.

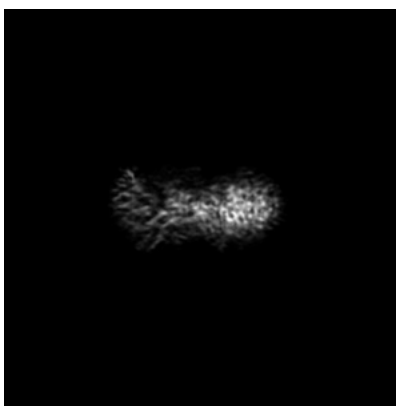
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

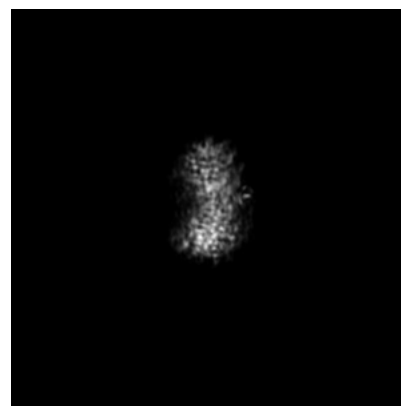
6.1.1 Primary map



X

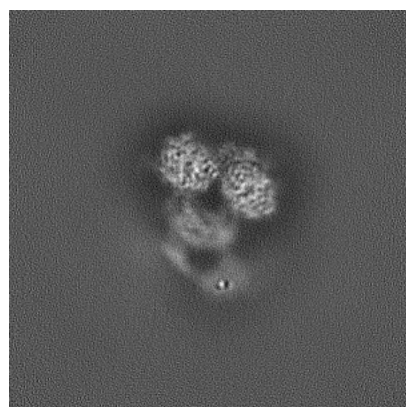


Y

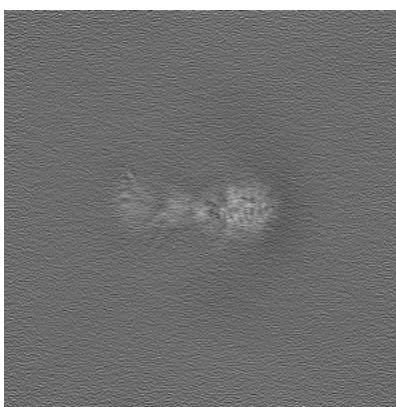


Z

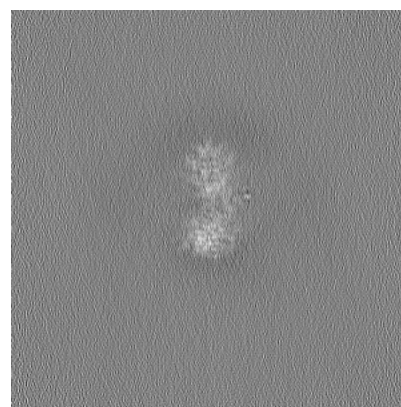
6.1.2 Raw map



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: 176

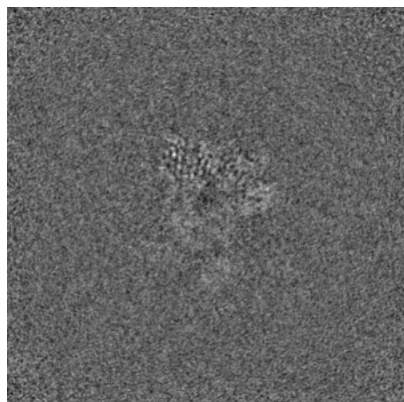


Y Index: 176



Z Index: 176

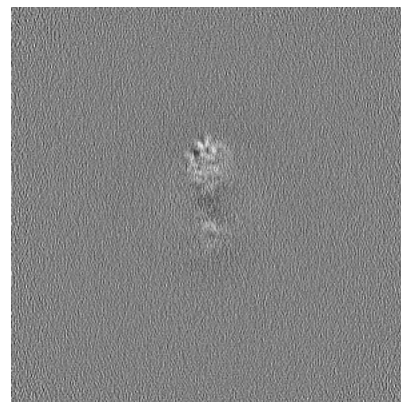
6.2.2 Raw map



X Index: 176



Y Index: 176



Z Index: 176

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: 178

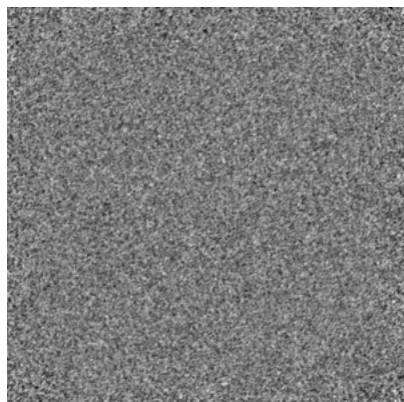


Y Index: 146

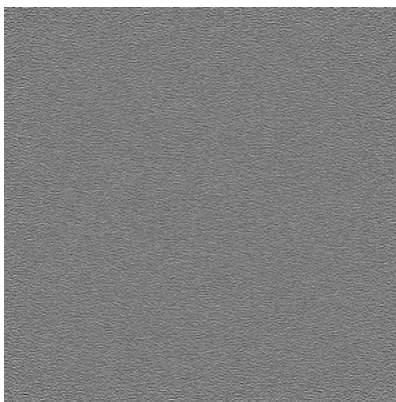


Z Index: 205

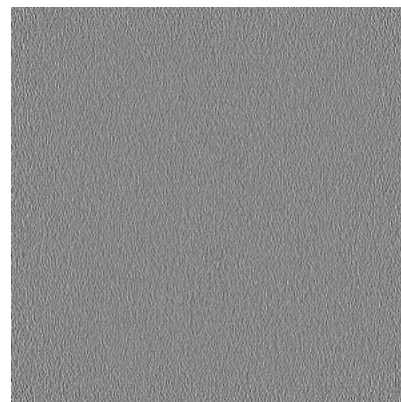
6.3.2 Raw map



X Index: 0



Y Index: 0



Z Index: 0

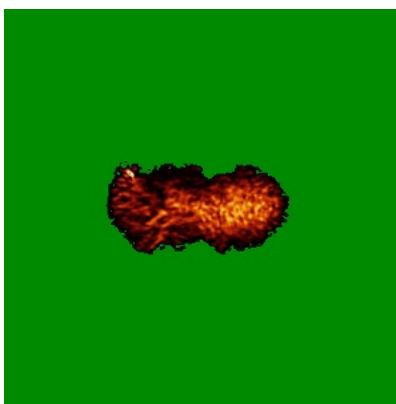
The images above show the largest variance slices of the map in three orthogonal directions.

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

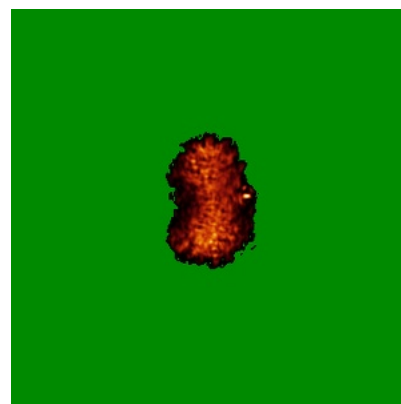
6.4.1 Primary map



X

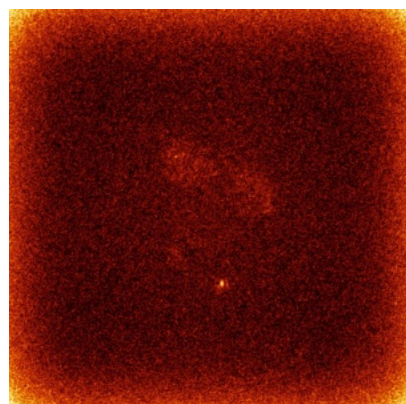


Y

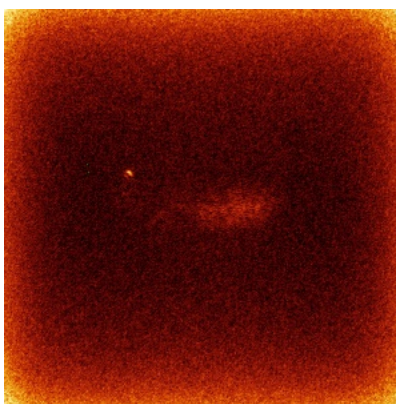


Z

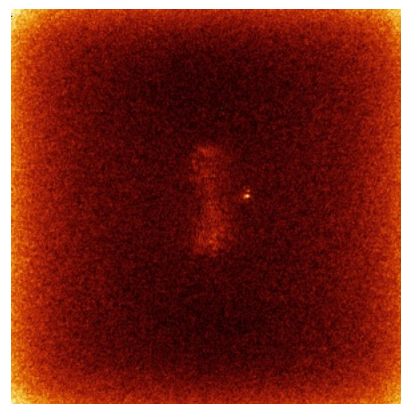
6.4.2 Raw map



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](#)

This section was not generated.

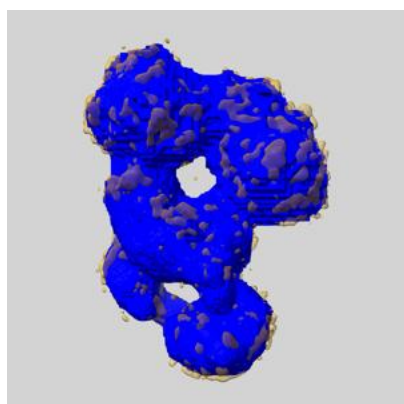
6.6 Mask visualisation [i](#)

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

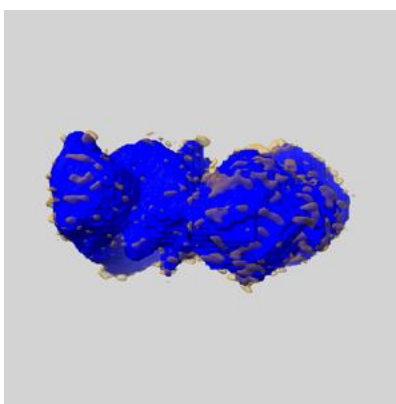
A mask typically either:

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

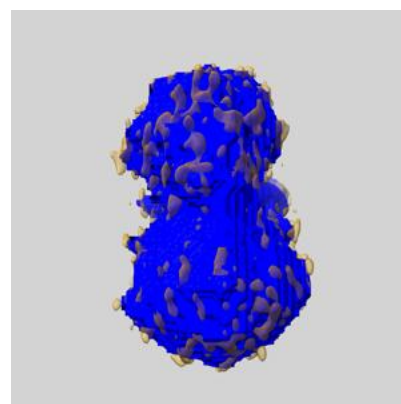
6.6.1 emd_26989_msk_1.map [i](#)



X

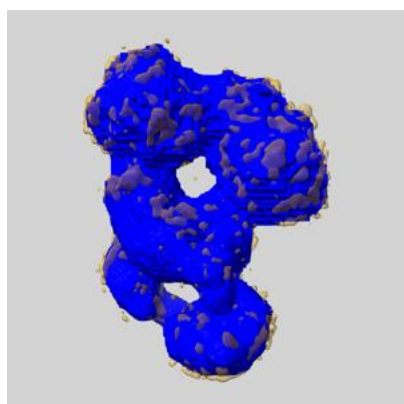


Y

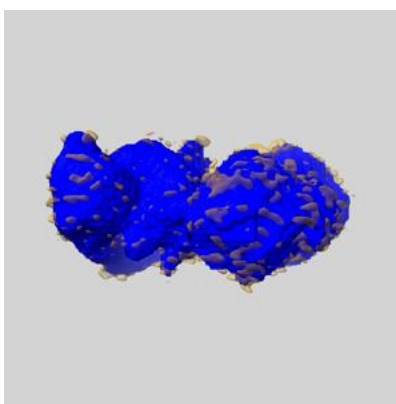


Z

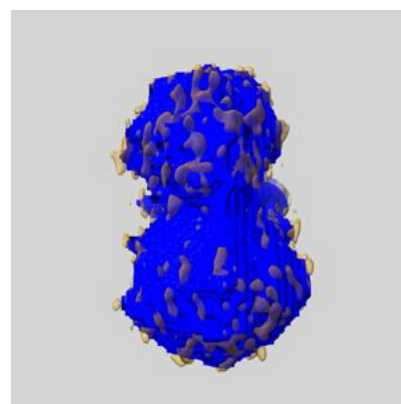
6.6.2 emd_26989_msk_2.map [i](#)



X



Y

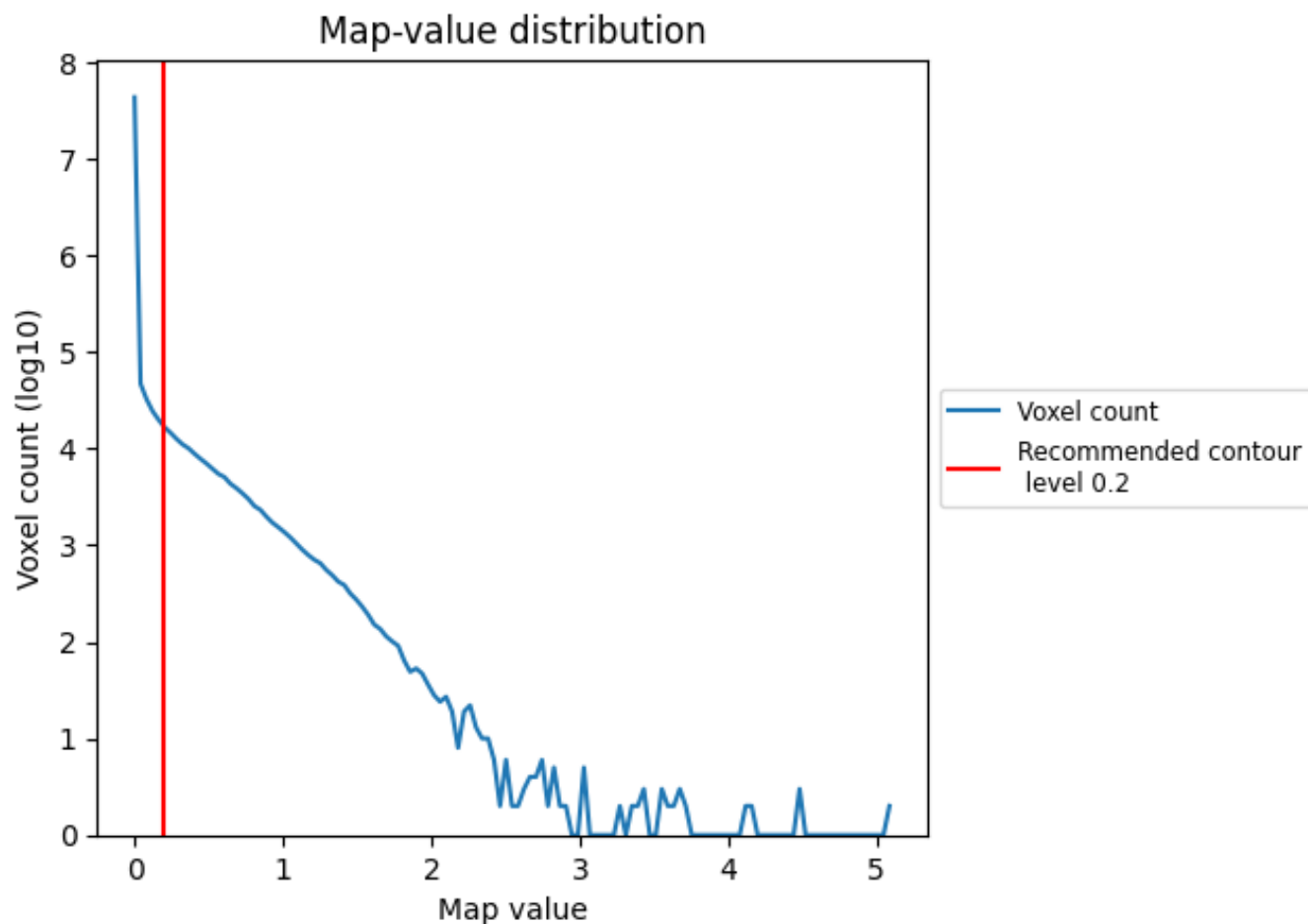


Z

7 Map analysis [i](#)

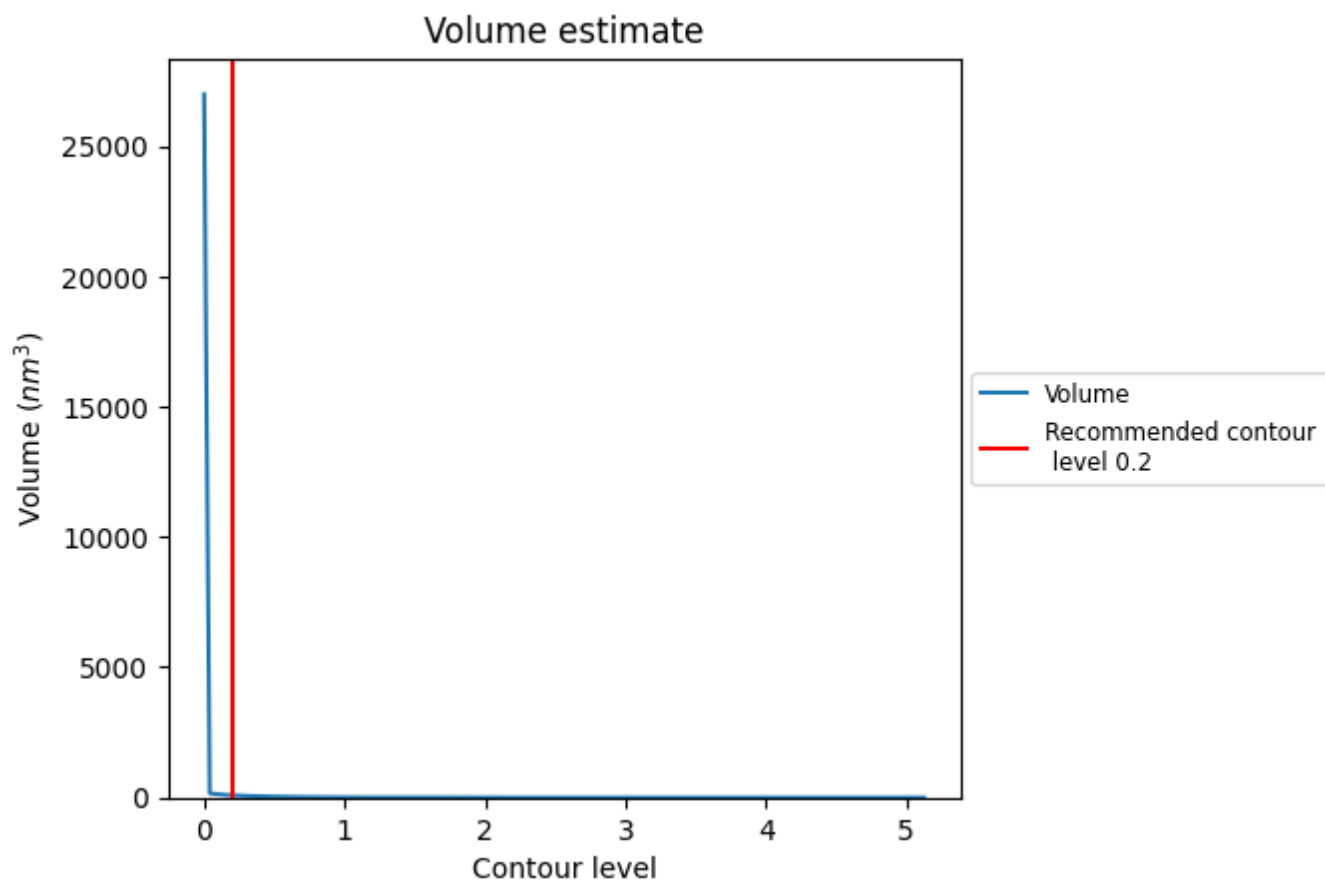
This section contains the results of statistical analysis of the map.

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.

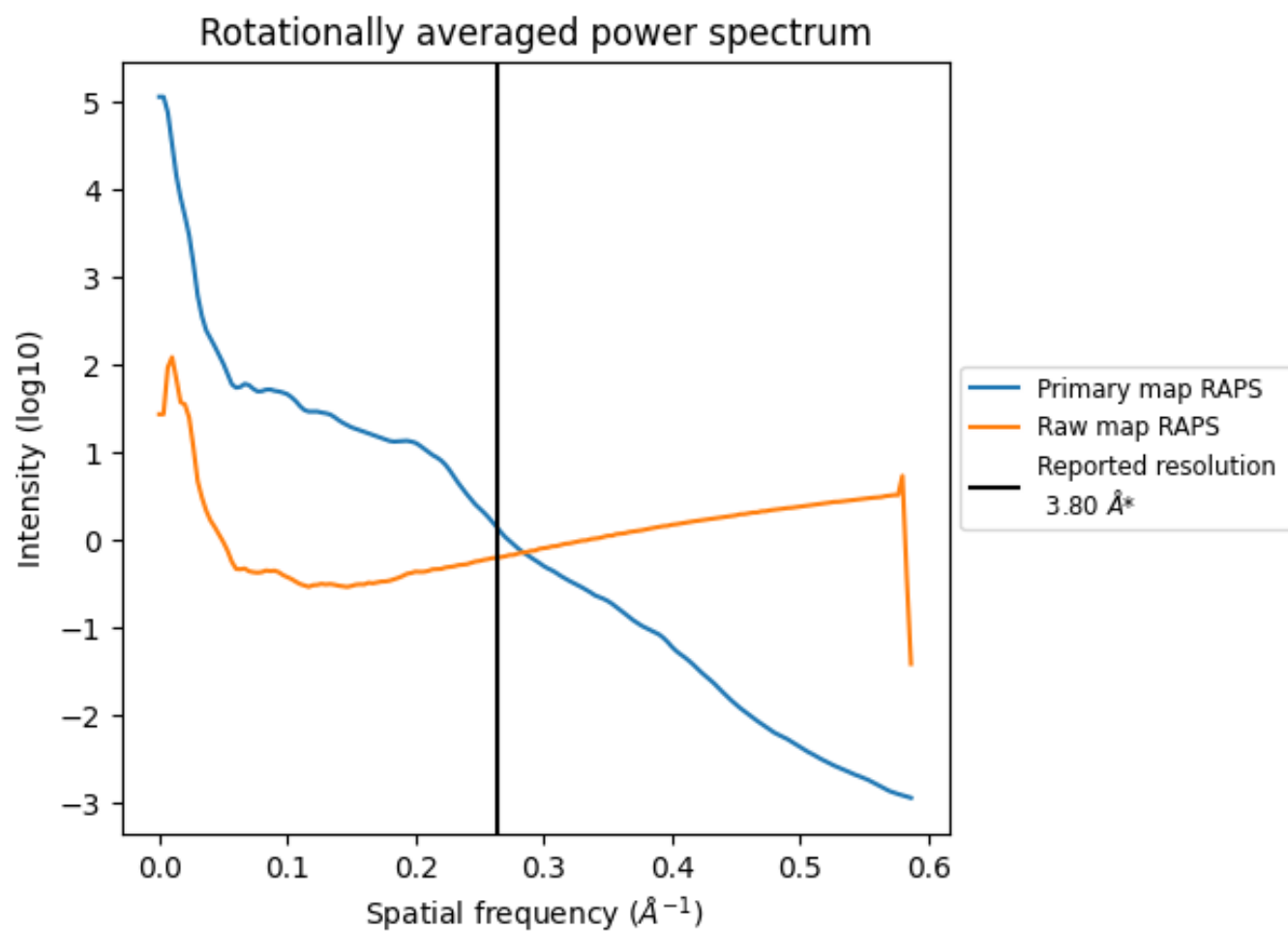
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 87 nm^3 ; this corresponds to an approximate mass of 79 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ

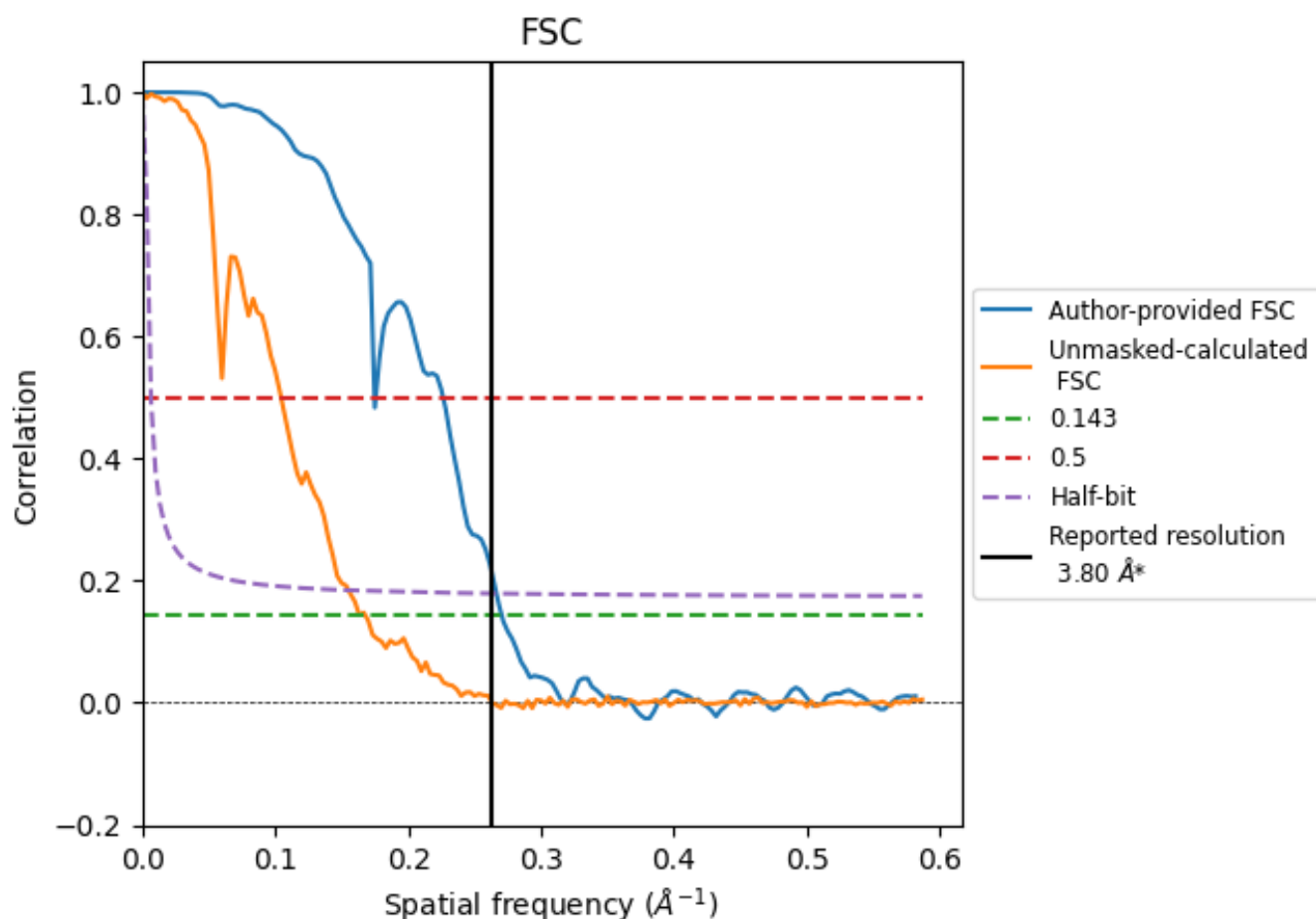


*Reported resolution corresponds to spatial frequency of 0.263 Å⁻¹

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 $0.263 \text{ \AA}^{-1}$

8.2 Resolution estimates [i](#)

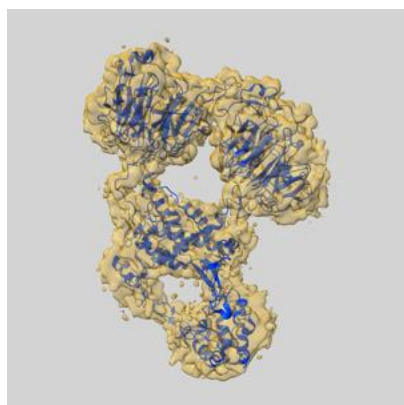
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.80	-	-
Author-provided FSC curve	3.70	5.72	3.75
Unmasked-calculated*	5.96	9.58	6.44

*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 5.96 differs from the reported value 3.8 by more than 10 %

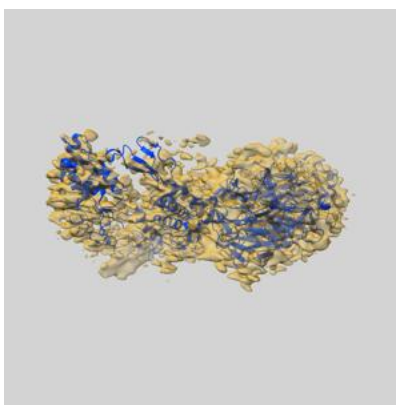
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-26989 and PDB model 8CTG. Per-residue inclusion information can be found in [section 3](#) on [page 7](#).

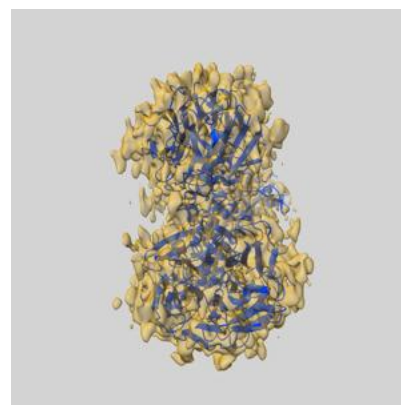
9.1 Map-model overlay [i](#)



X



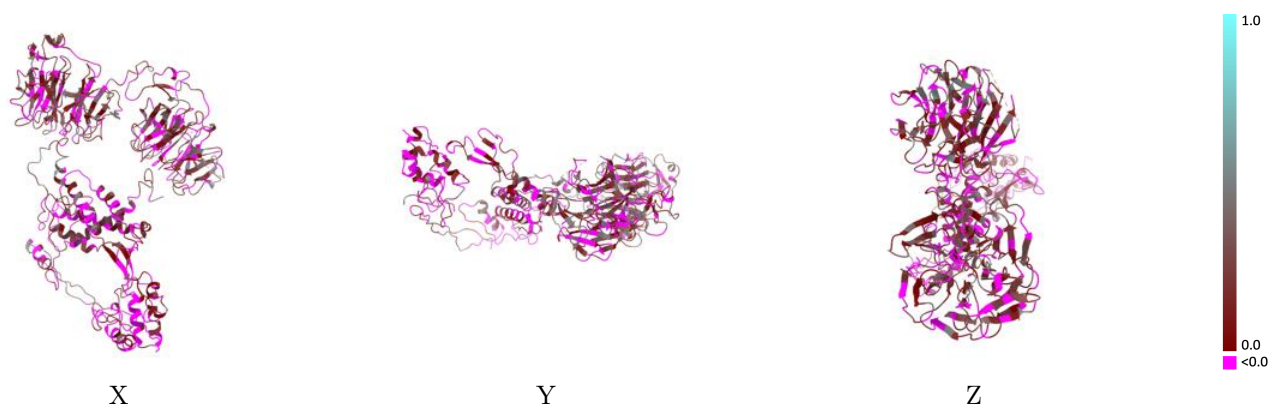
Y



Z

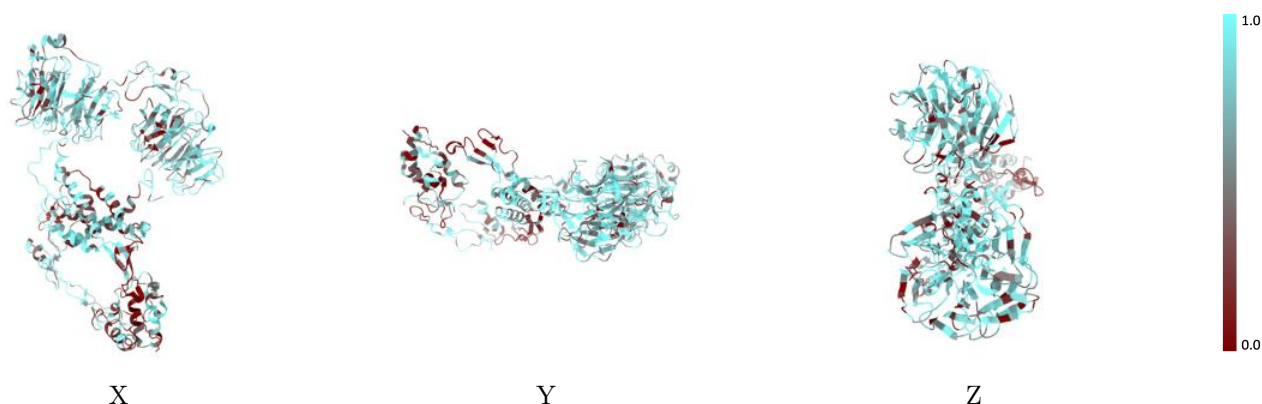
The images above show the 3D surface view of the map at the recommended contour level 0.2 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



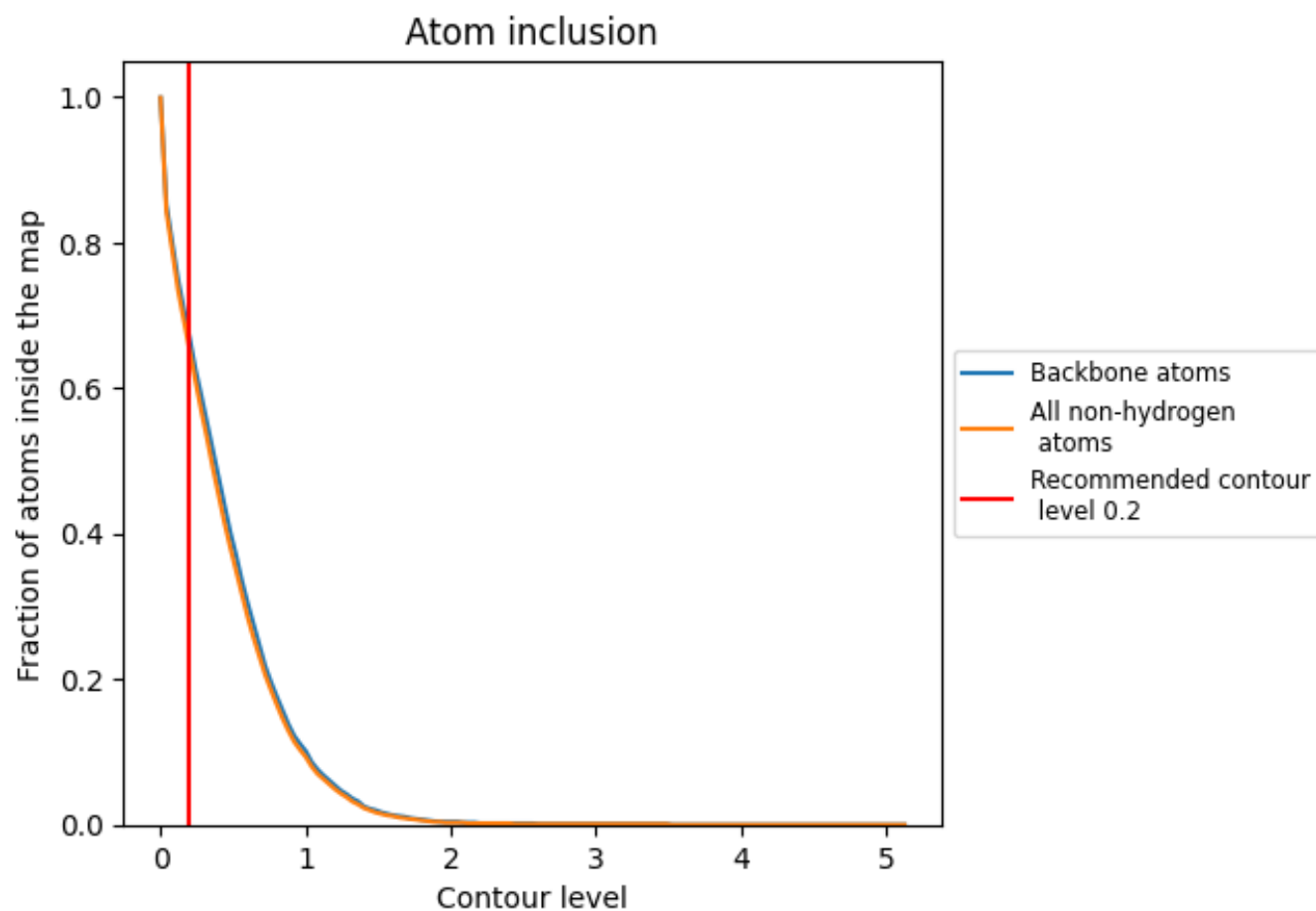
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.2).

9.4 Atom inclusion ⓘ



At the recommended contour level, 67% of all backbone atoms, 65% 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 (0.2) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.6500	0.1090
A	0.4840	0.0110
B	0.6110	0.0770
C	0.7080	0.1470

