



wwPDB EM Validation Summary Report ⓘ

Aug 27, 2026 – 05:39 am BST

PDB ID : 9G03 / pdb_00009g03
EMDB ID : EMD-50909
Title : Structure of carbon monoxide dehydrogenase/acetyl-CoA synthase (CODH/ACS) in complex with ferredoxin (Clostridium autoethanogenum)
Authors : Yin, M.D.; Lemaire, O.N.; Wagner, T.; Murphy, B.J.
Deposited on : 2024-07-06
Resolution : 2.10 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.dev133
Mogul : 1.8.4, CSD as541be (2020)
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) : 2.50

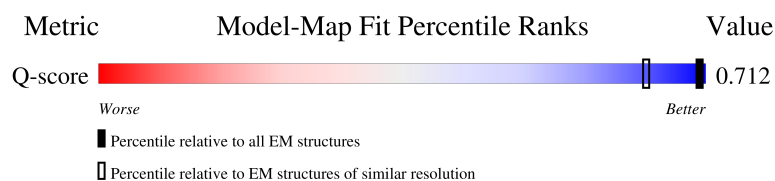
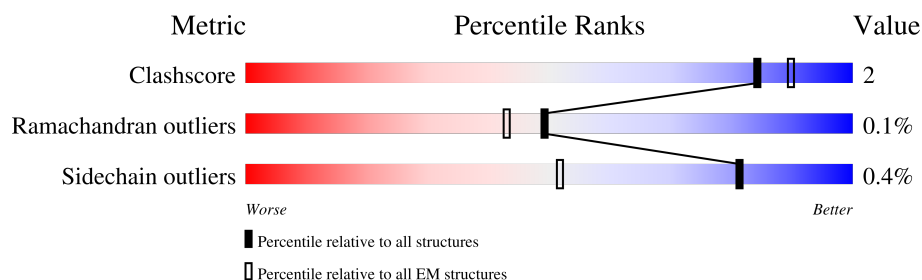
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

The reported resolution of this entry is 2.10 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	2317 (1.60 - 2.60)

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

Mol	Chain	Length	Quality of chain
1	A	708	
1	D	708	
2	B	630	
2	C	630	

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Mol	Chain	Length	Quality of chain
3	E	57	9% 74% 18% 9%

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 14176 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 CO-methylating acetyl-CoA synthase.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	290	Total	C	N	O	S	0	0
			2154	1391	347	409	7		
1	D	291	Total	C	N	O	S	0	0
			2163	1397	349	410	7		

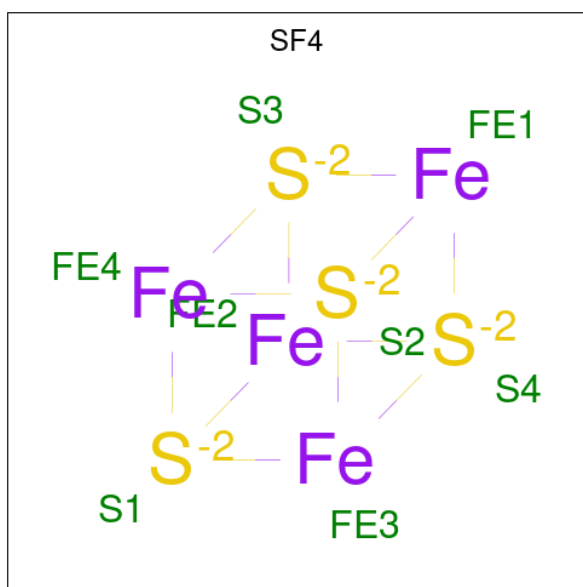
- Molecule 2 is a protein called Carbon monoxide dehydrogenase/acetyl-CoA synthase beta subunit.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	629	Total	C	N	O	S	0	0
			4723	2967	814	902	40		
2	C	628	Total	C	N	O	S	0	0
			4714	2962	813	899	40		

- Molecule 3 is a protein called Ferredoxin.

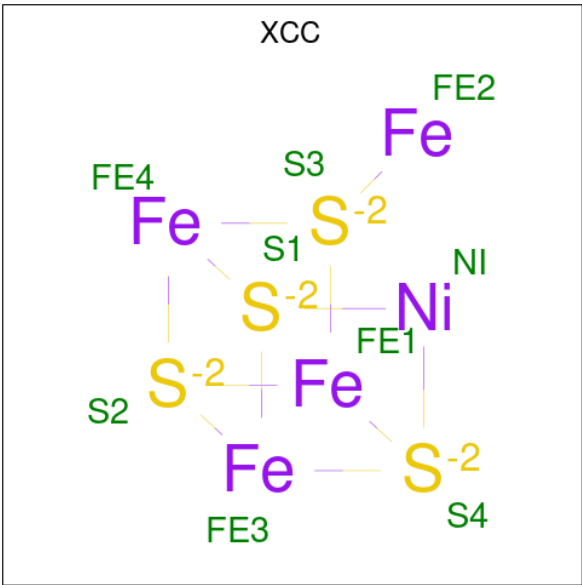
Mol	Chain	Residues	Atoms					AltConf	Trace
3	E	52	Total	C	N	O	S	0	0
			364	219	58	79	8		

- Molecule 4 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			AltConf
4	B	1	Total	Fe	S	0
			8	4	4	
4	C	1	Total	Fe	S	0
			8	4	4	
4	C	1	Total	Fe	S	0
			8	4	4	
4	E	1	Total	Fe	S	0
			8	4	4	
4	E	1	Total	Fe	S	0
			8	4	4	

- Molecule 5 is FE(4)-NI(1)-S(4) CLUSTER (CCD ID: XCC) (formula: Fe_4NiS_4) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
5	B	1	Total	Fe	Ni	S	0
			9	4	1	4	
5	C	1	Total	Fe	Ni	S	0
			9	4	1	4	

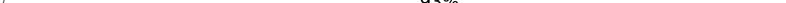
LEU	TYR	LEU	GLY	GLU	THR	GLU	THR	ALA
ASN	LEU	ASN	LEU	GLU	THR	VAL	GLN	CYS
ASP	ILE	ASP	ALA	SER	VAL	SER	PHE	GLN
ASN	ASP	SER	MET	SER	ILE	ILE	ALA	ASP
PHE	GLY	GLY	GLN	SER	GLN	GLN	PHE	GLY
ALA	GLY	GLY	GLY	GLY	GLY	GLY	HIS	ALA
MET	MET	GLY	GLY	GLY	GLY	VAL	VAL	CYS
ILE	ILE	VAL	VAL	THR	THR	THR	THR	VAL
CYS	CYS	GLN	THR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

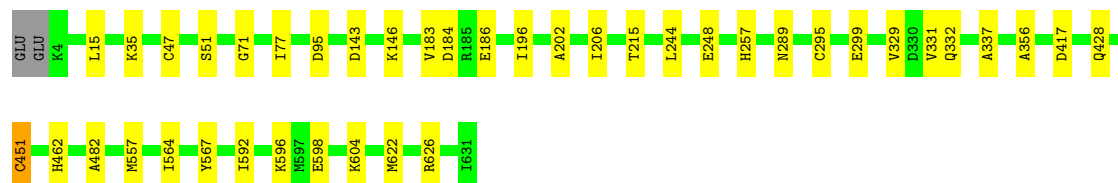
- Molecule 2: Carbon monoxide dehydrogenase/acetyl-CoA synthase beta subunit

Chain B:  95% 5%

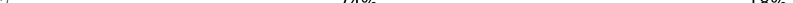


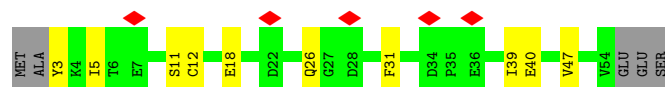
- Molecule 2: Carbon monoxide dehydrogenase/acetyl-CoA synthase beta subunit

Chain C:  93% 6%



- Molecule 3: Ferredoxin

Chain E:  9% 74% 18% 9%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	896132	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 ($e^-/\text{\AA}^2$)	70	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	165000	Depositor
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	4.367	Depositor
Minimum map value	-3.037	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.025	Depositor
Recommended contour level	0.213	Depositor
Map size (Å)	490.56, 490.56, 490.56	wwPDB
Map dimensions	560, 560, 560	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.876, 0.876, 0.876	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: SF4, XCC

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	0.19	0/2185	0.33	0/2970
1	D	0.21	0/2194	0.36	0/2981
2	B	0.25	0/4797	0.39	0/6488
2	C	0.25	0/4788	0.39	0/6476
3	E	0.21	0/369	0.47	0/502
All	All	0.23	0/14333	0.38	0/19417

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	2154	0	2264	8	0
1	D	2163	0	2277	11	0
2	B	4723	0	4803	22	0
2	C	4714	0	4797	25	0
3	E	364	0	329	10	0
4	B	8	0	0	1	0
4	C	16	0	0	0	0
4	E	16	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
5	B	9	0	0	1	0
5	C	9	0	0	1	0
All	All	14176	0	14470	67	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 2.

The worst 5 of 67 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:C:451:CYS:HB3	5:C:703:XCC:S2	2.33	0.69
2:B:451:CYS:HB3	5:B:702:XCC:S2	2.35	0.66
3:E:39:ILE:HG22	3:E:39:ILE:O	2.00	0.61
2:B:257:HIS:HB3	2:B:329:VAL:HG12	1.83	0.60
2:B:196:ILE:HD13	4:B:701:SF4:S1	2.42	0.59

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	288/708 (41%)	284 (99%)	4 (1%)	0	100	100
1	D	289/708 (41%)	285 (99%)	4 (1%)	0	100	100
2	B	627/630 (100%)	610 (97%)	16 (3%)	1 (0%)	43	44
2	C	626/630 (99%)	608 (97%)	17 (3%)	1 (0%)	43	44
3	E	50/57 (88%)	45 (90%)	5 (10%)	0	100	100
All	All	1880/2733 (69%)	1832 (97%)	46 (2%)	2 (0%)	49	50

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	C	332	GLN
2	B	332	GLN

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	232/583 (40%)	232 (100%)	0	100	100
1	D	233/583 (40%)	233 (100%)	0	100	100
2	B	518/519 (100%)	516 (100%)	2 (0%)	84	89
2	C	517/519 (100%)	514 (99%)	3 (1%)	78	86
3	E	43/47 (92%)	42 (98%)	1 (2%)	44	51
All	All	1543/2251 (68%)	1537 (100%)	6 (0%)	81	89

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

Mol	Chain	Res	Type
2	C	196	ILE
2	C	451	CYS
3	E	3	TYR
2	B	451	CYS
2	B	95	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 10 such sidechains are listed below:

Mol	Chain	Res	Type
2	C	357	HIS
2	C	388	ASN
1	D	5	GLN
2	B	10	GLN
2	B	199	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains [i](#)

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

7 ligands are 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	$\# Z > 2$
5	XCC	B	702	2	0,12,12	-	-	-		
4	SF4	C	701	2	0,12,12	-	-	-		
4	SF4	C	702	2	0,12,12	-	-	-		
4	SF4	E	102	3	0,12,12	-	-	-		
4	SF4	E	101	3	0,12,12	-	-	-		
4	SF4	B	701	2	0,12,12	-	-	-		
5	XCC	C	703	2	0,12,12	-	-	-		

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
5	XCC	B	702	2	-	-	0/4/4/4
4	SF4	C	701	2	-	-	0/6/5/5
4	SF4	C	702	2	-	-	0/6/5/5

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	SF4	E	102	3	-	-	0/6/5/5
4	SF4	E	101	3	-	-	0/6/5/5
4	SF4	B	701	2	-	-	0/6/5/5
5	XCC	C	703	2	-	-	0/4/4/4

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

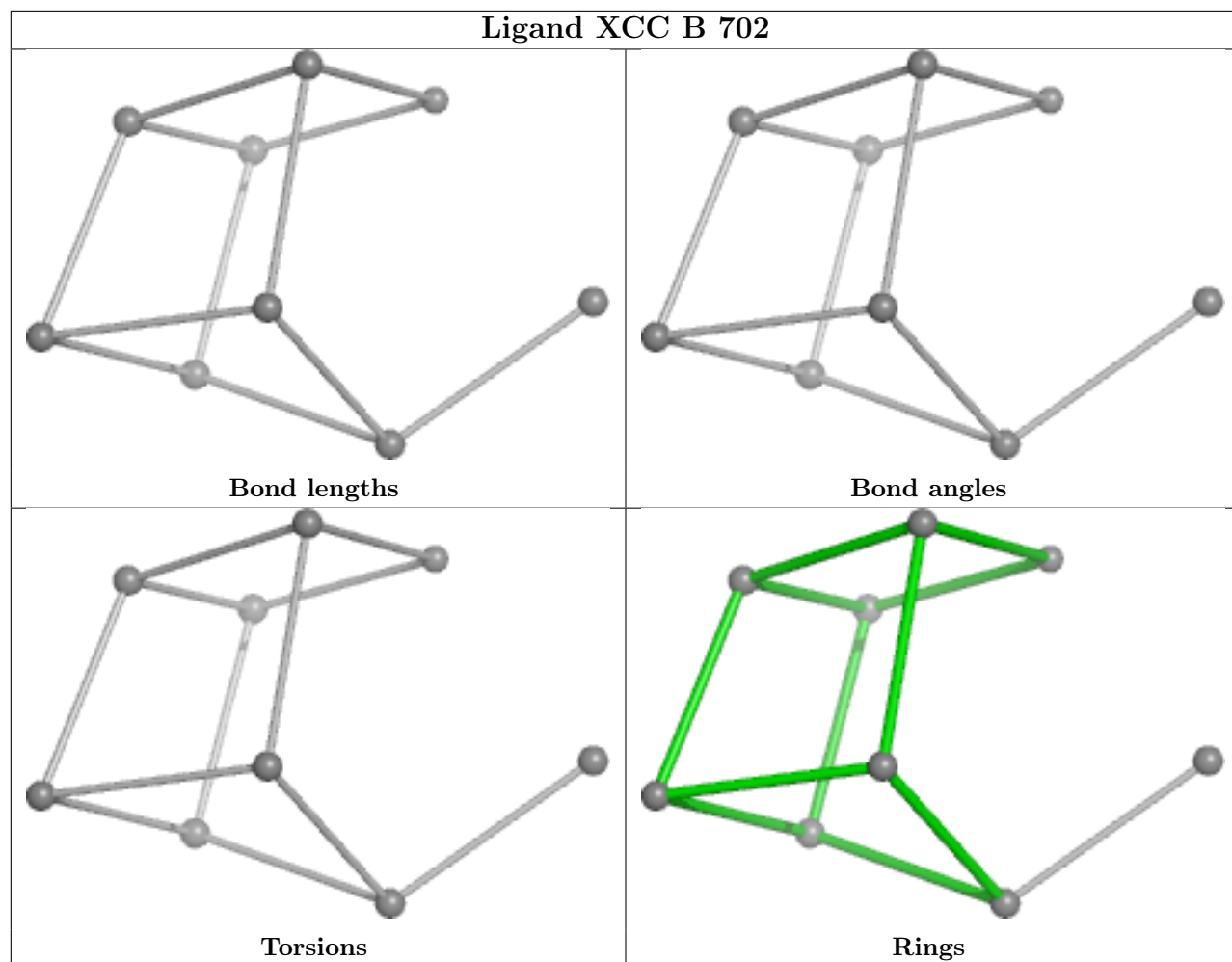
There are no torsion outliers.

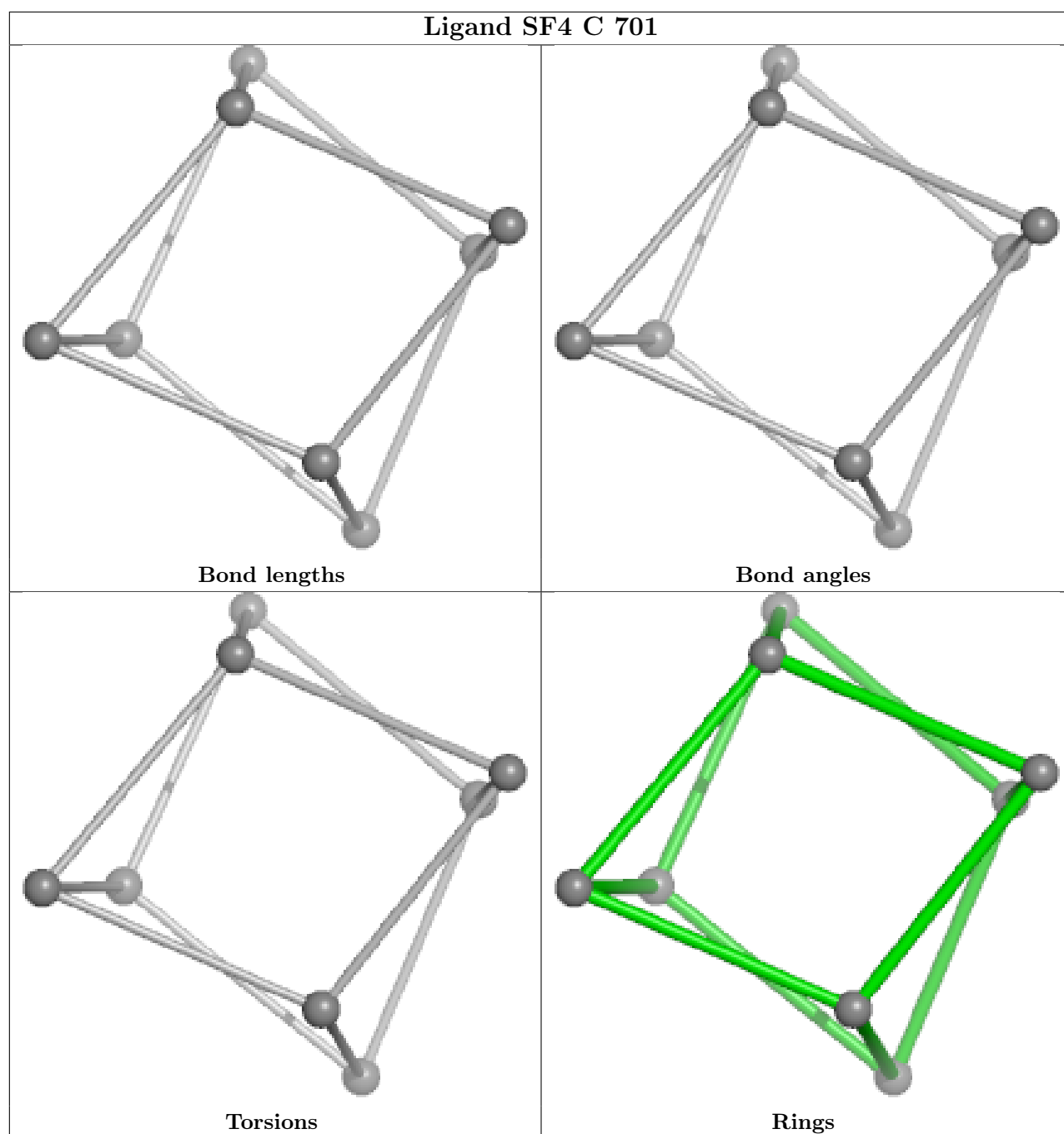
There are no ring outliers.

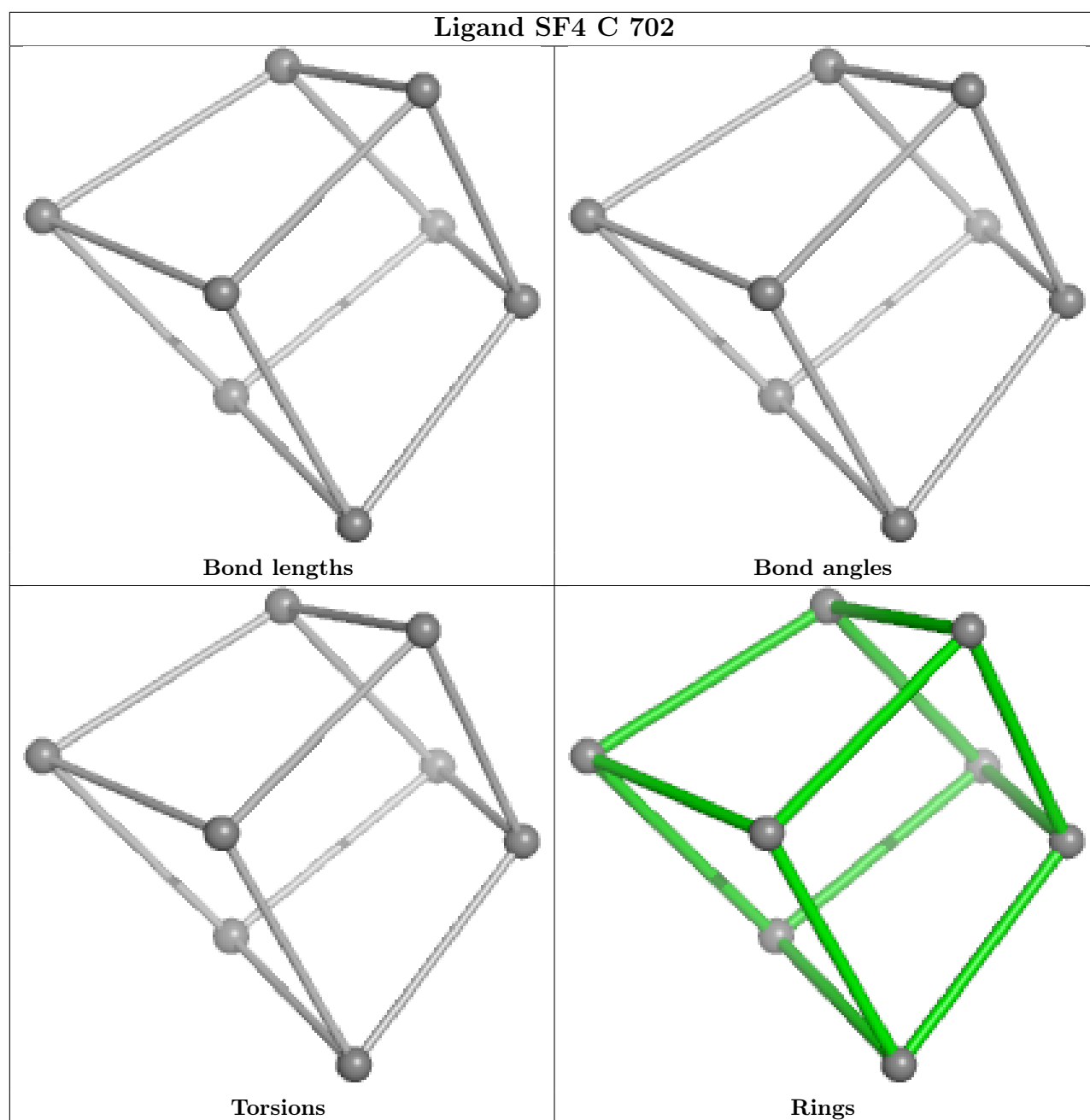
3 monomers are involved in 3 short contacts:

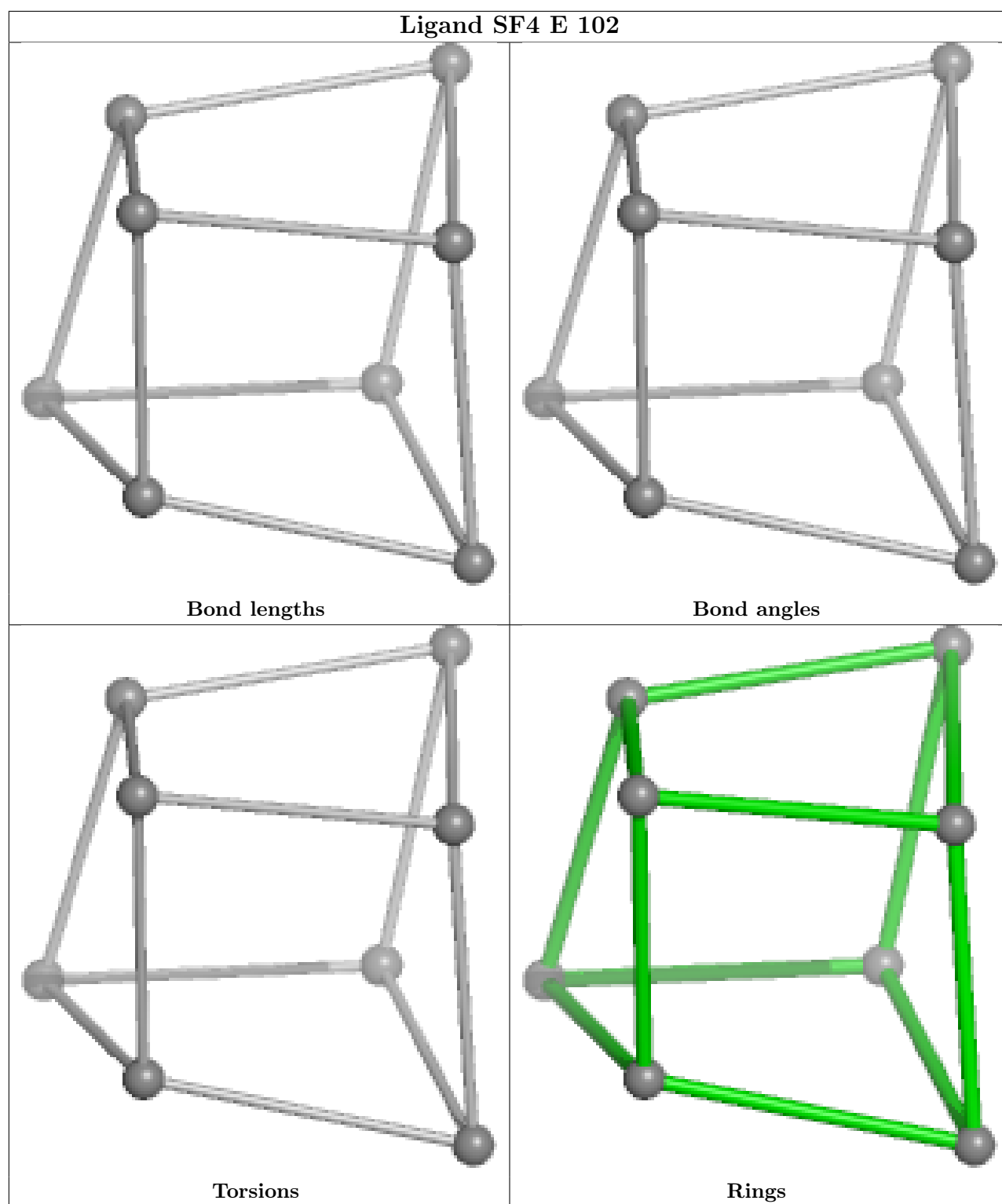
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	702	XCC	1	0
4	B	701	SF4	1	0
5	C	703	XCC	1	0

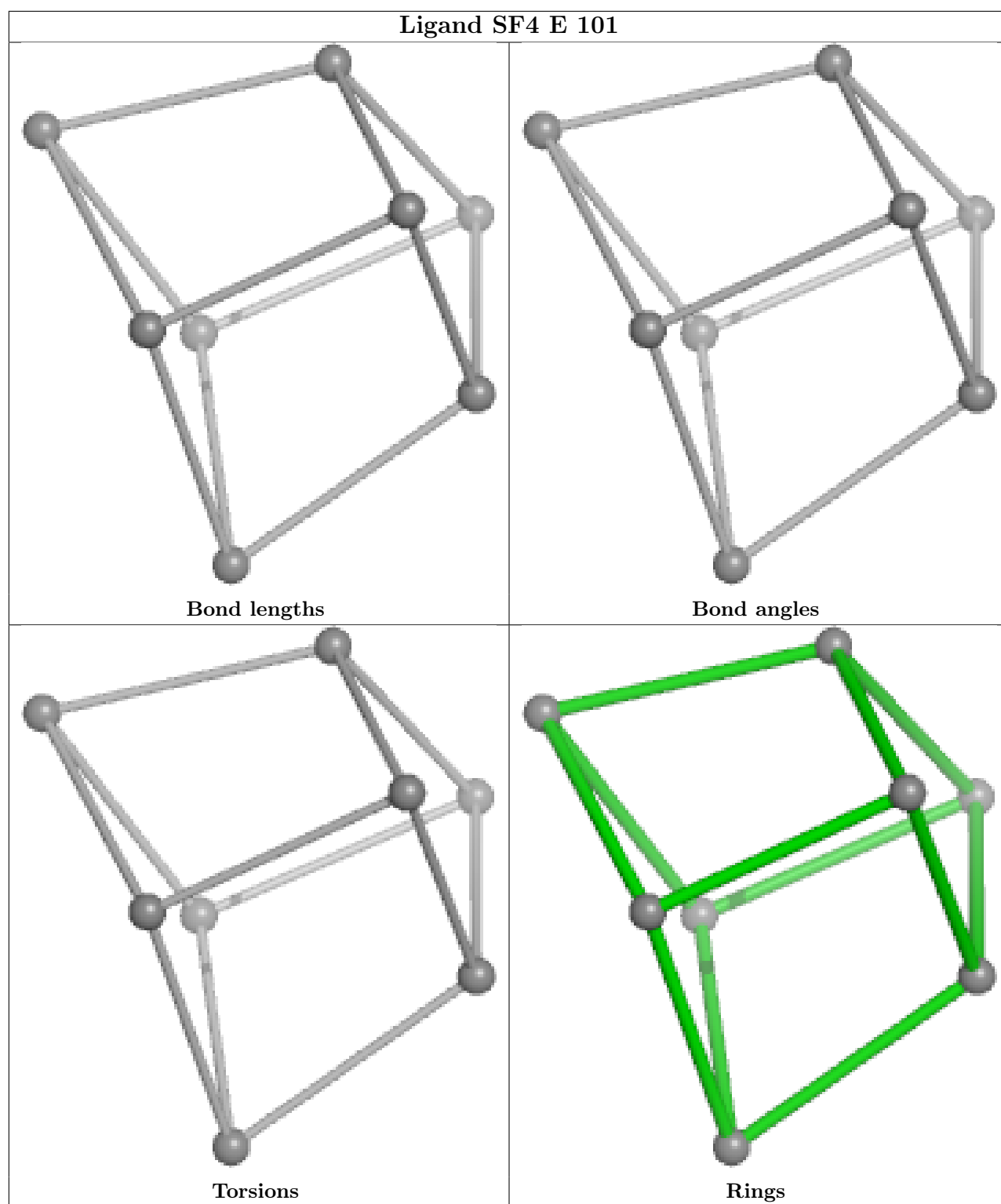
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. 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.

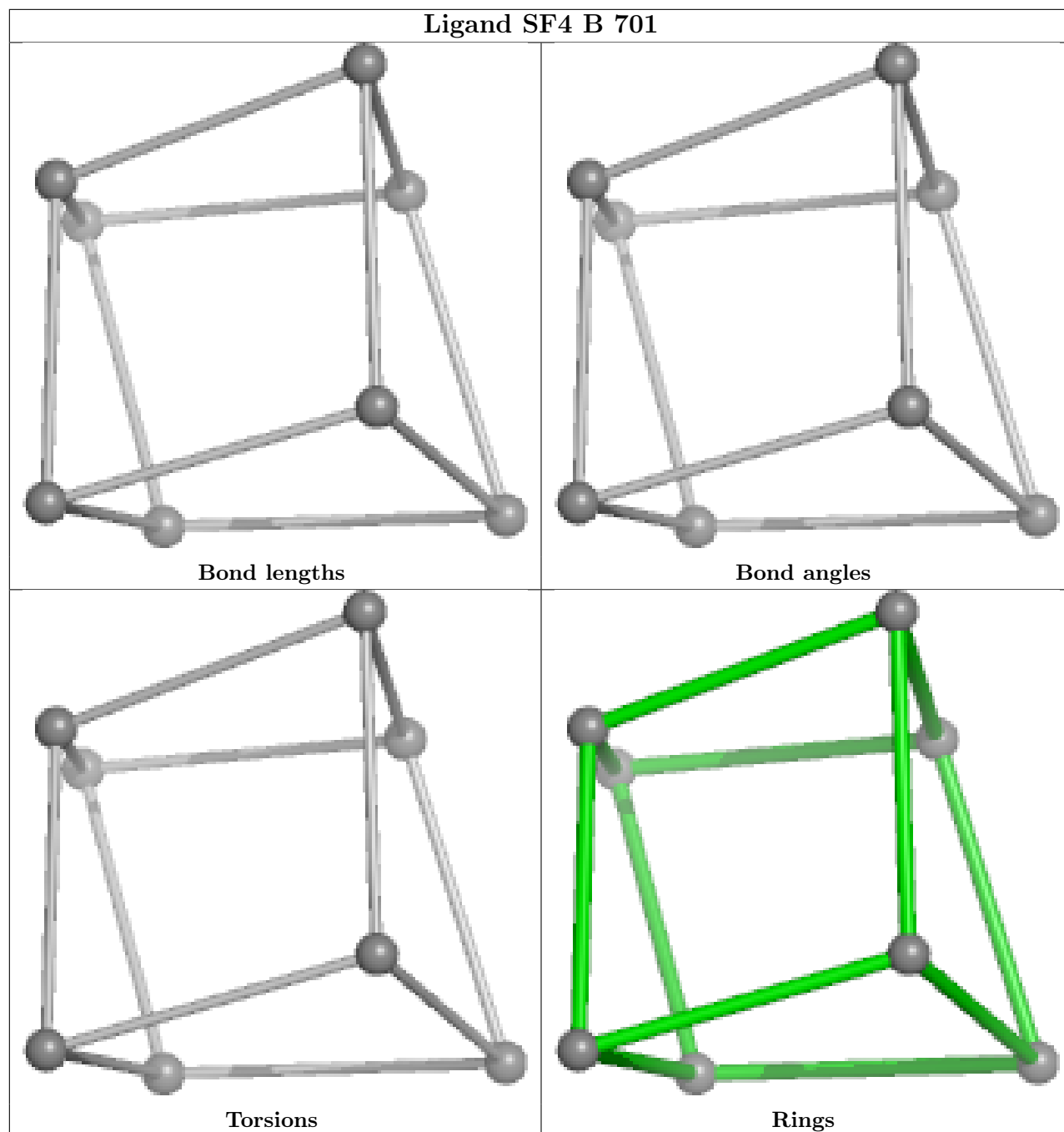


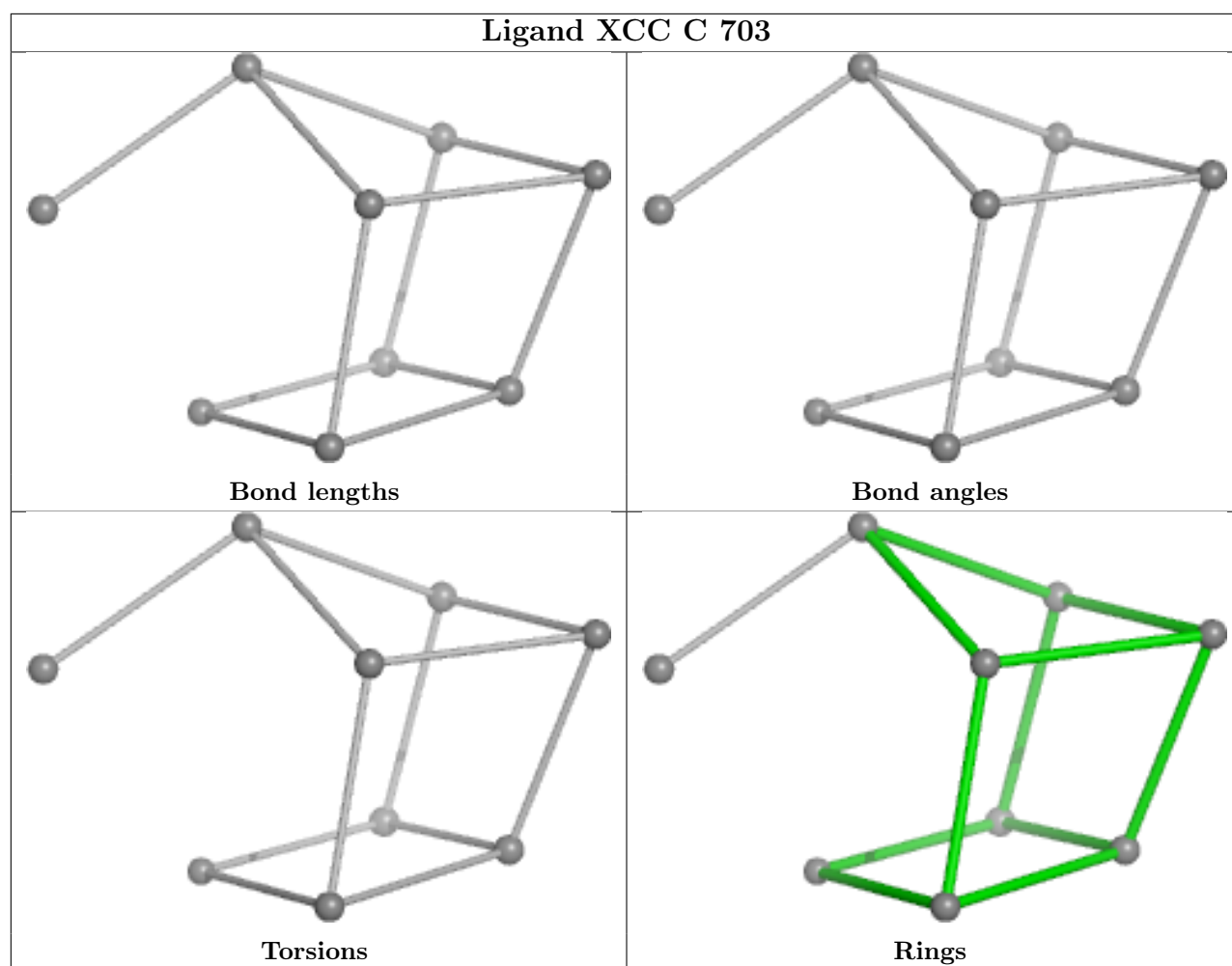












5.7 Other polymers [i](#)

There are no such residues in this entry.

5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

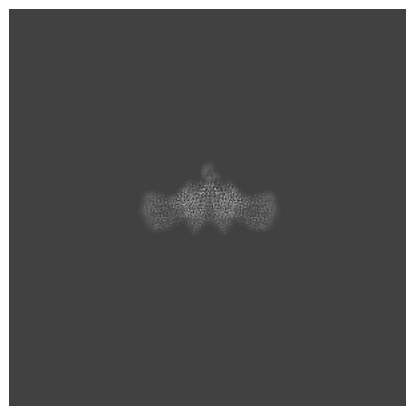
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-50909. These allow visual inspection of the internal detail of the map and identification of artifacts.

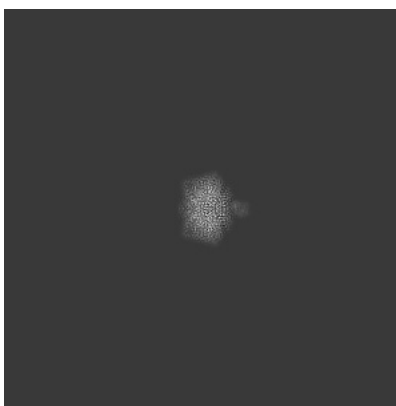
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

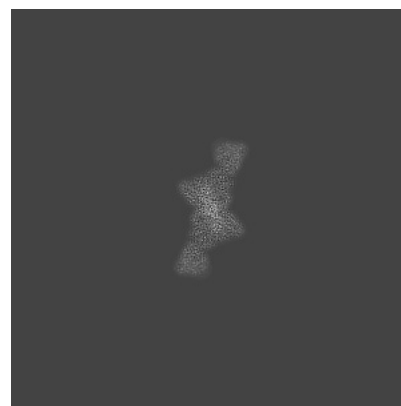
6.1.1 Primary map



X

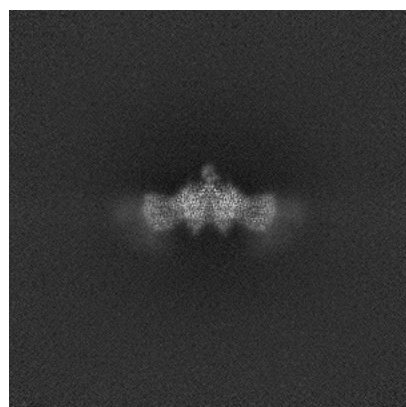


Y

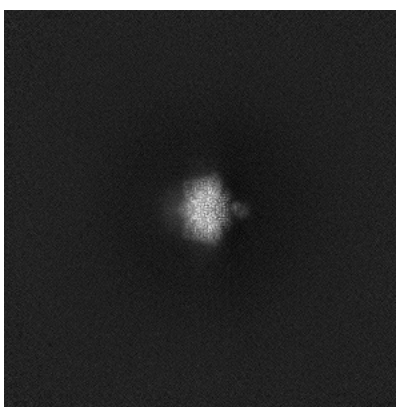


Z

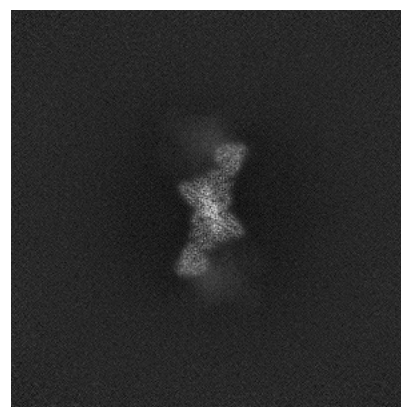
6.1.2 Raw map



X



Y

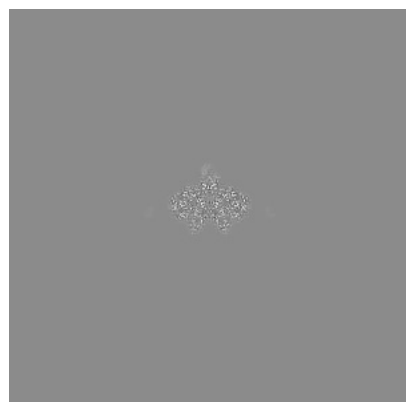


Z

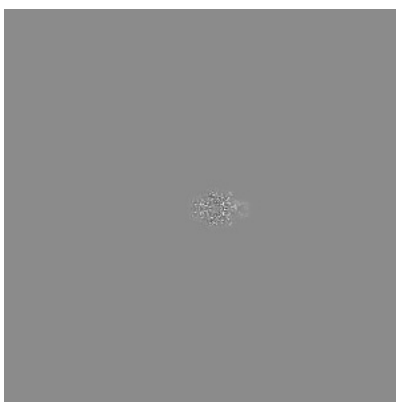
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

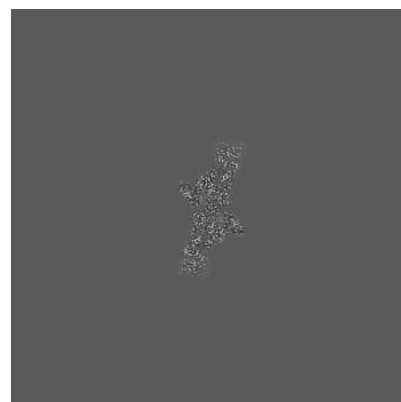
6.2.1 Primary map



X Index: 280

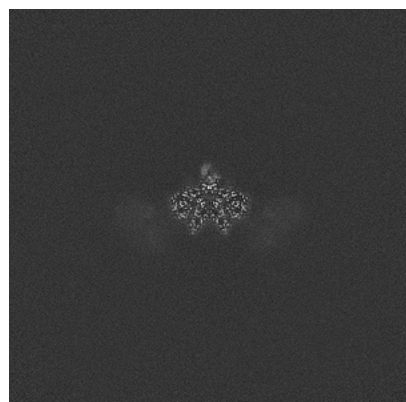


Y Index: 280

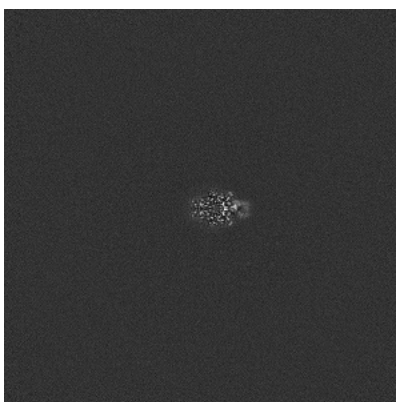


Z Index: 280

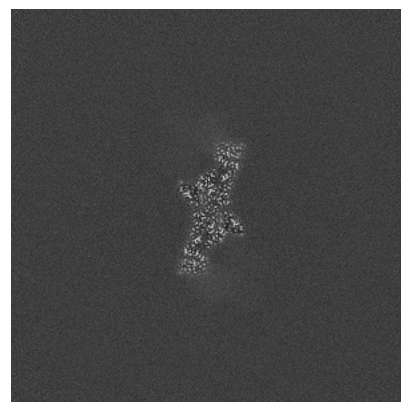
6.2.2 Raw map



X Index: 280



Y Index: 280

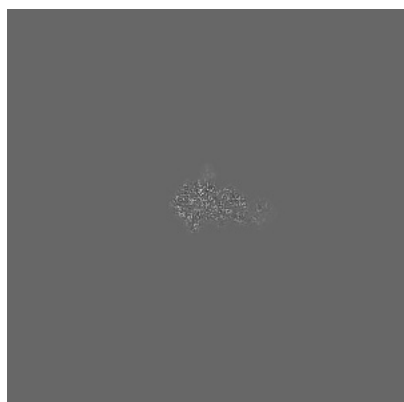


Z Index: 280

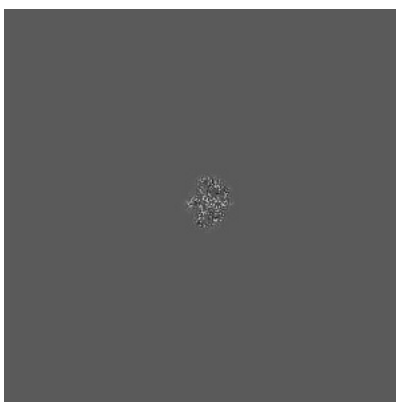
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

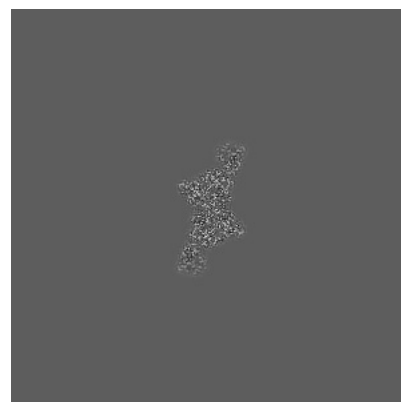
6.3.1 Primary map



X Index: 288

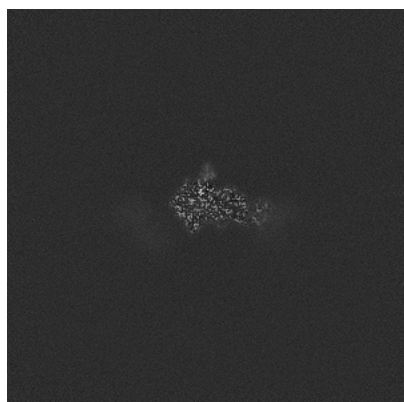


Y Index: 251

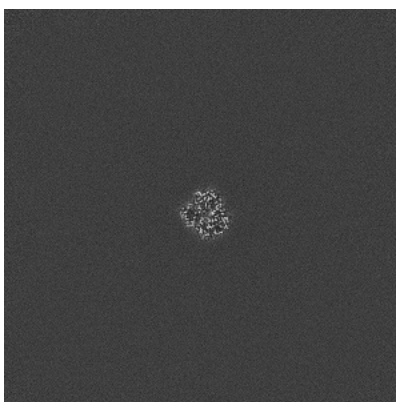


Z Index: 287

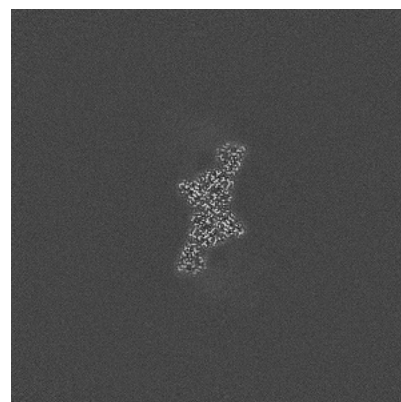
6.3.2 Raw map



X Index: 288



Y Index: 300

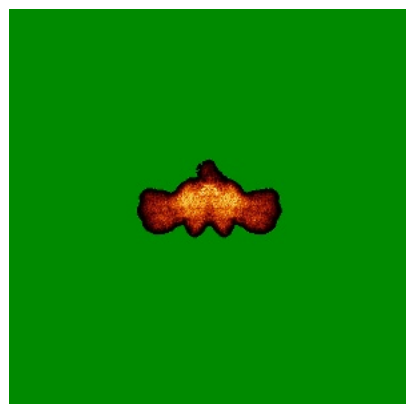


Z Index: 287

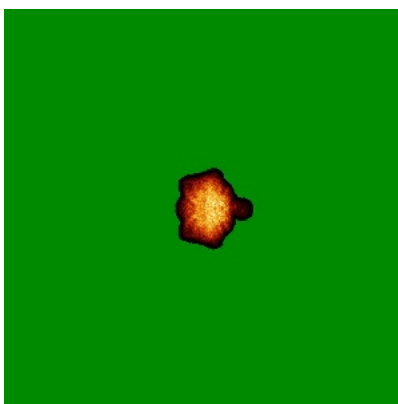
The images above show the largest variance slices of the map in three orthogonal directions.

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

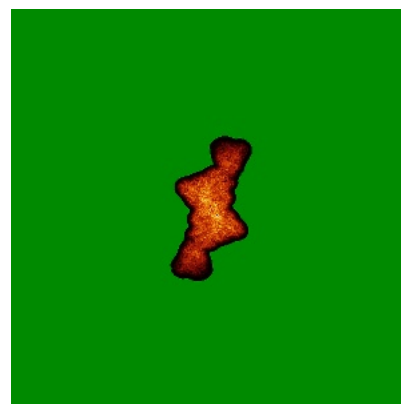
6.4.1 Primary map



X

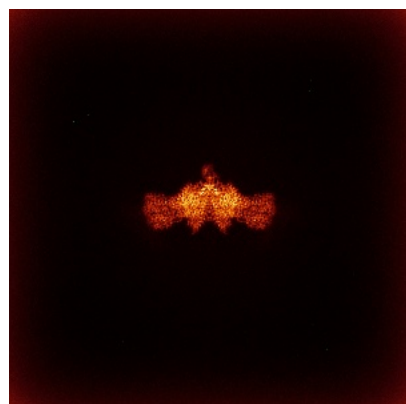


Y

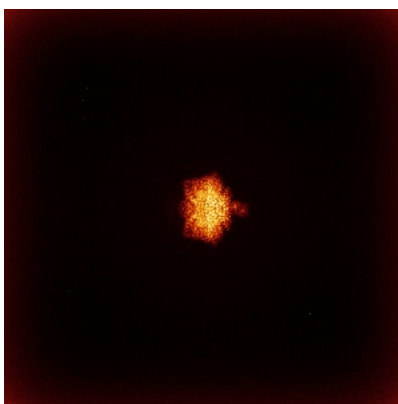


Z

6.4.2 Raw map



X



Y

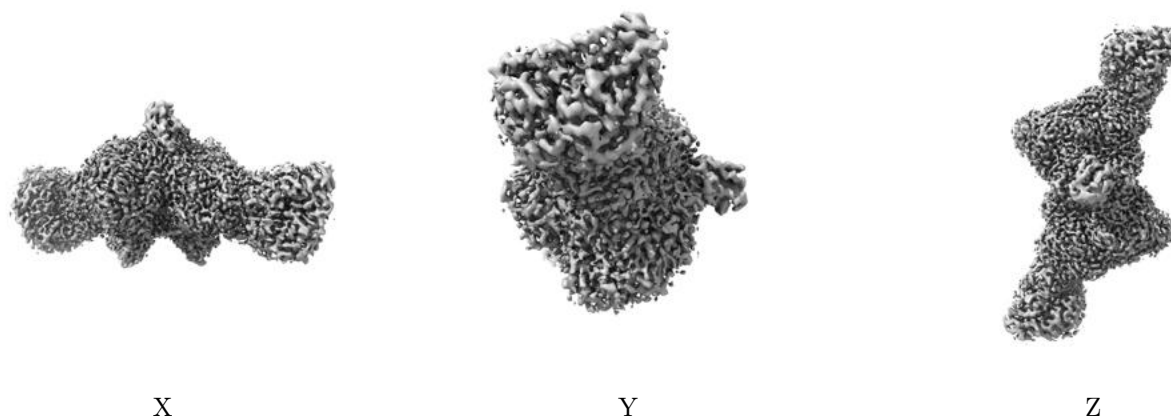


Z

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.

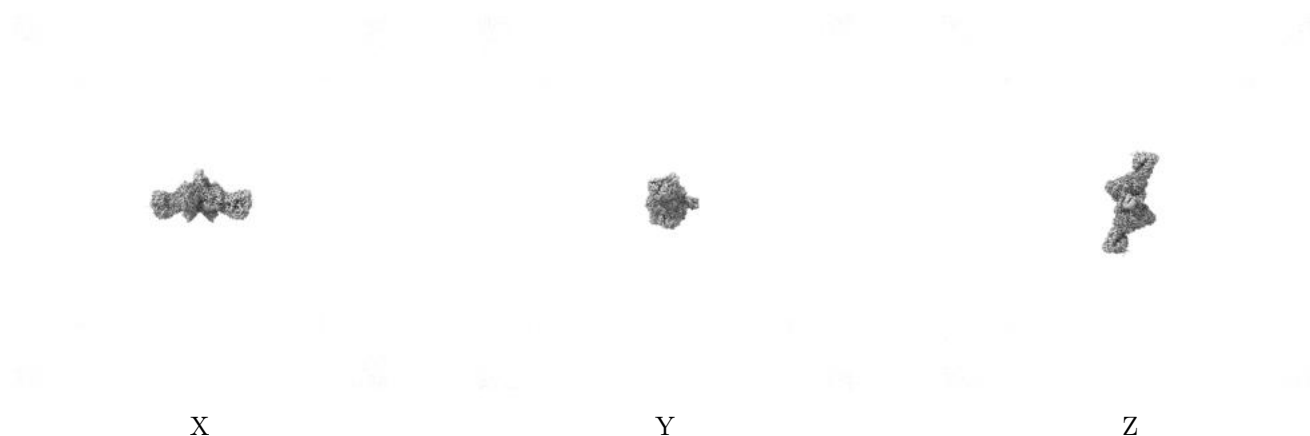
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.213. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



These images show the 3D surface of the raw map. The raw map's contour level was selected so that its surface encloses the same volume as the primary map does at its recommended contour level.

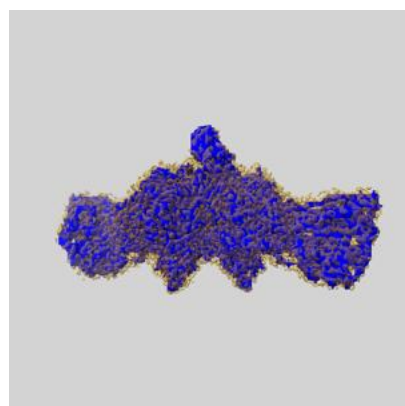
6.6 Mask visualisation [i](#)

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

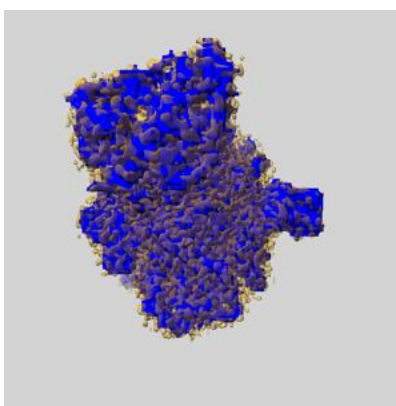
A mask typically either:

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

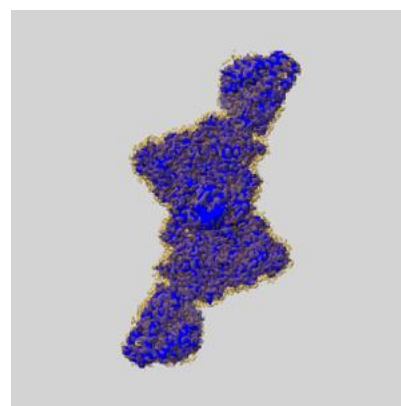
6.6.1 emd_50909_msk_1.map [i](#)



X



Y

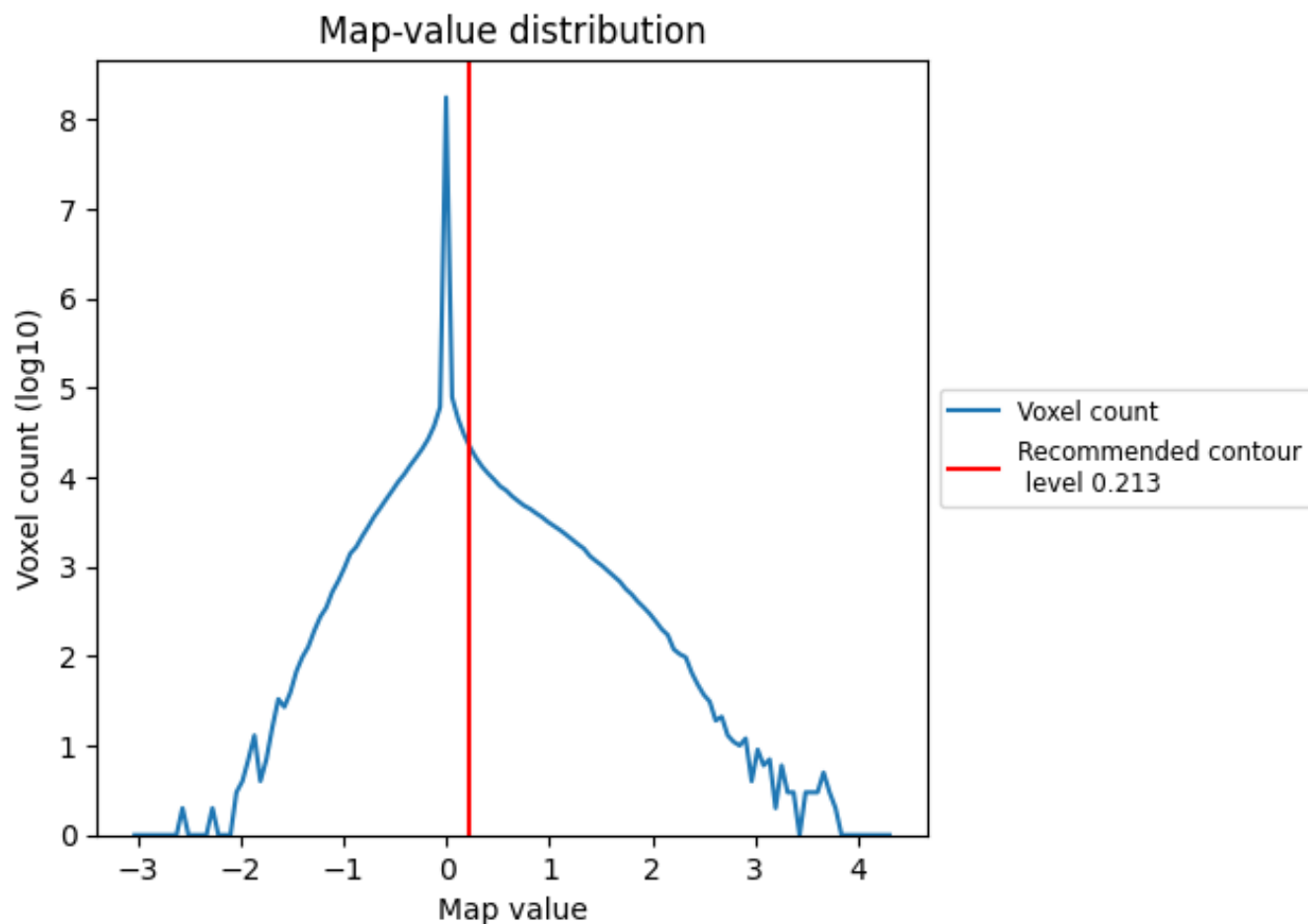


Z

7 Map analysis [i](#)

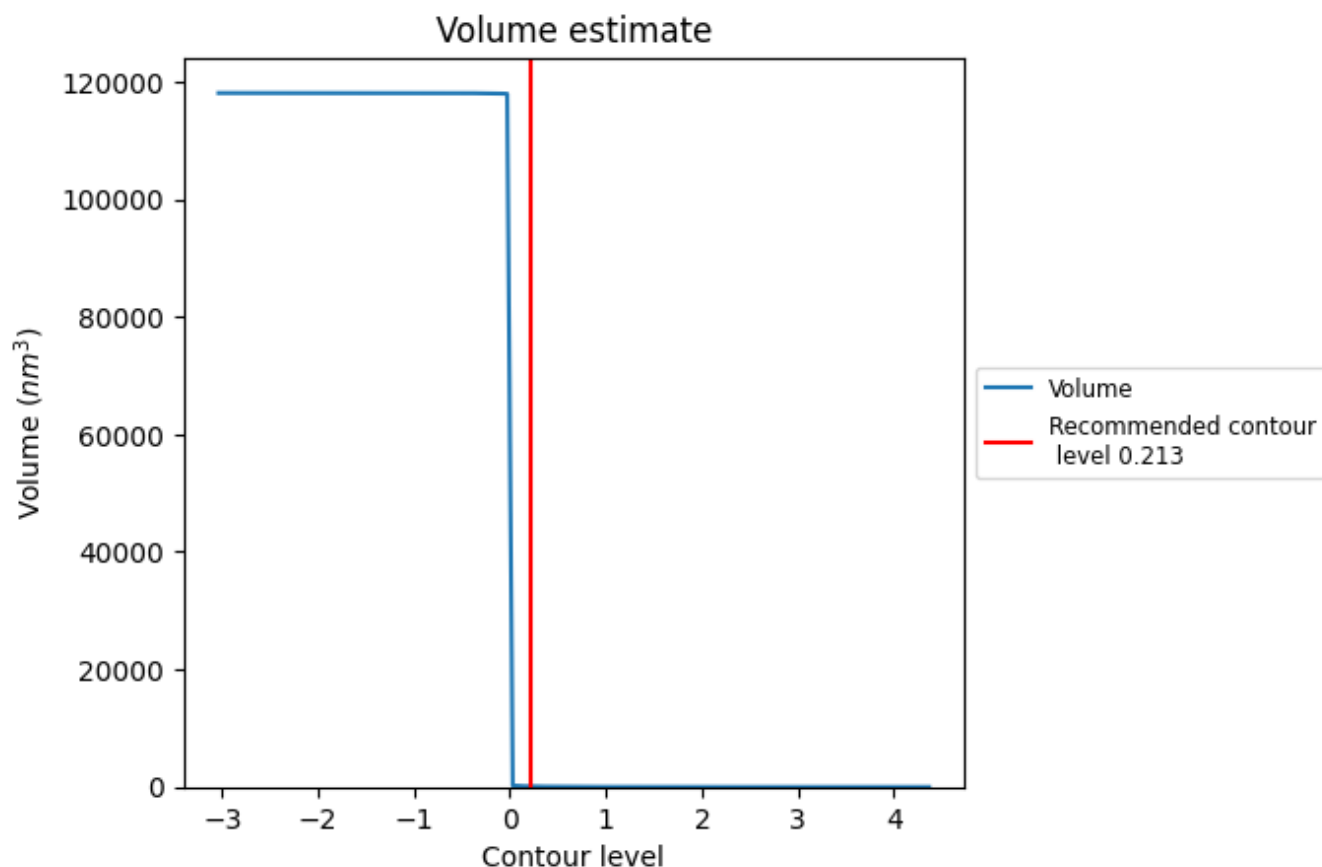
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



The map-value distribution is plotted in 128 intervals along the x-axis. The y-axis is logarithmic. A spike in this graph at zero usually indicates that the volume has been masked.

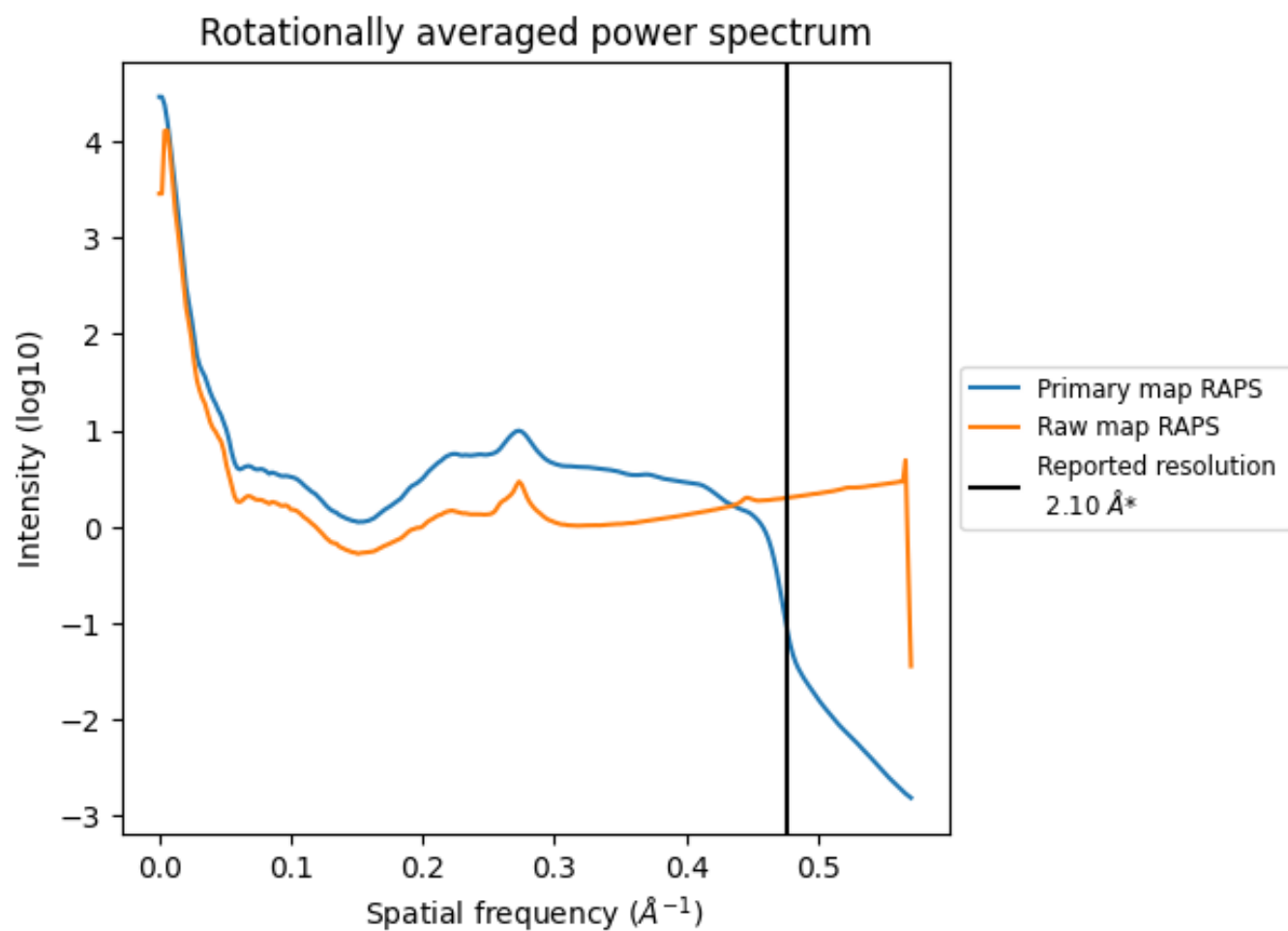
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 99 nm³; this corresponds to an approximate mass of 89 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 ⓘ

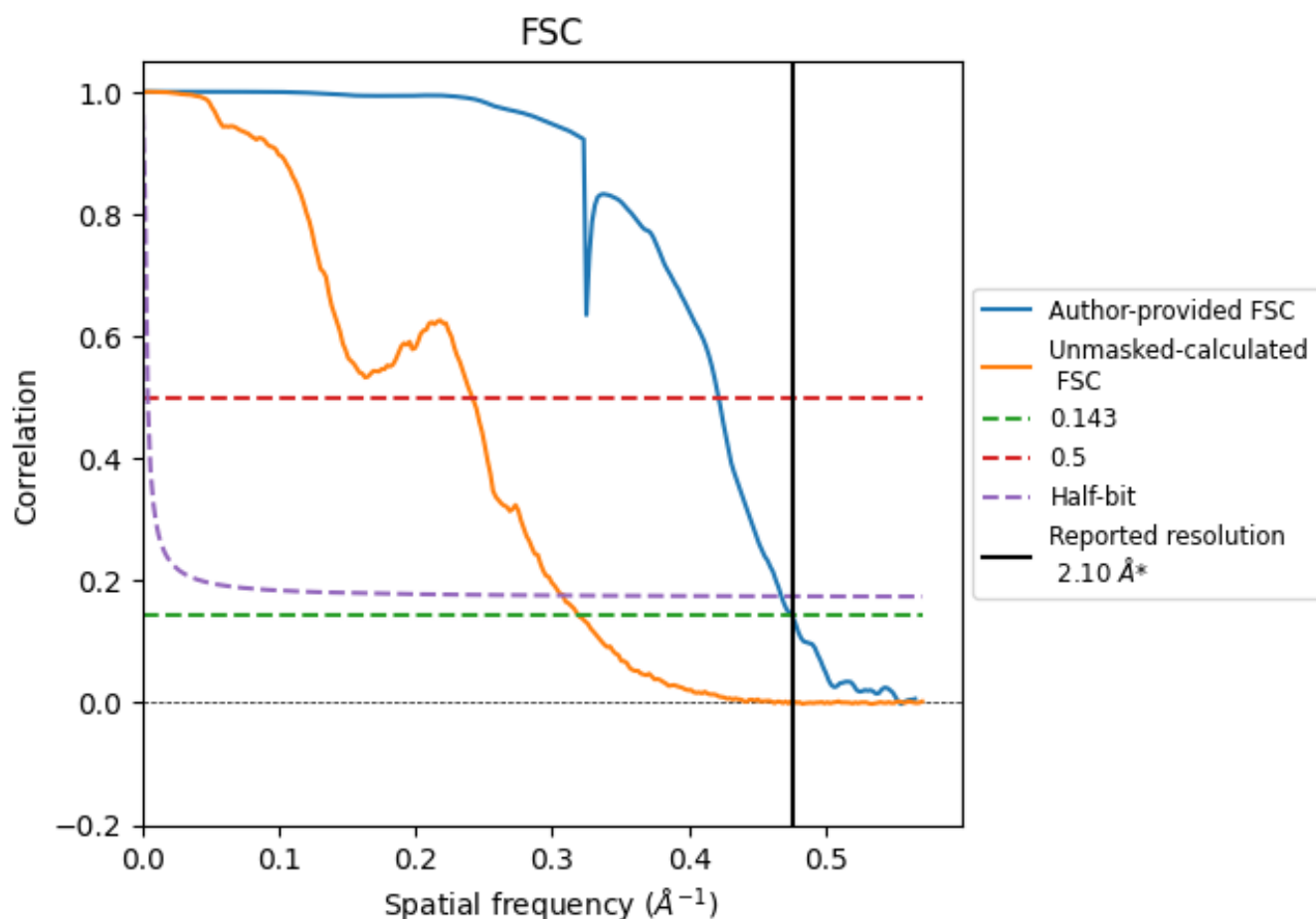


*Reported resolution corresponds to spatial frequency of 0.476 Å⁻¹

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 0.476 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

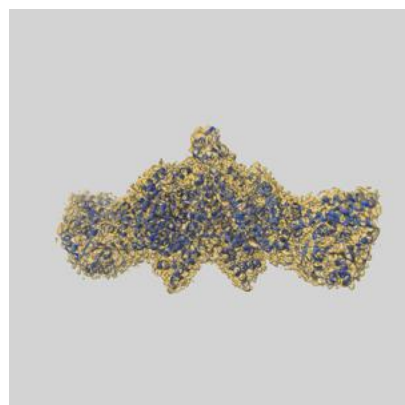
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.10	-	-
Author-provided FSC curve	2.10	2.37	2.14
Unmasked-calculated*	3.13	4.14	3.25

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.13 differs from the reported value 2.1 by more than 10 %

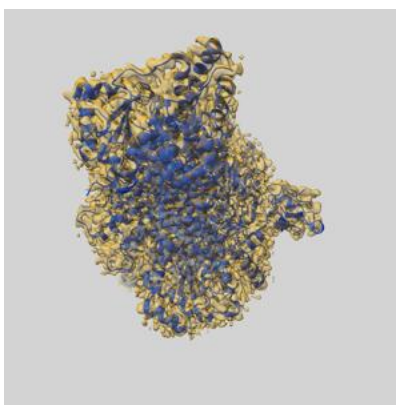
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-50909 and PDB model 9G03. Per-residue inclusion information can be found in section [3](#) on page [7](#).

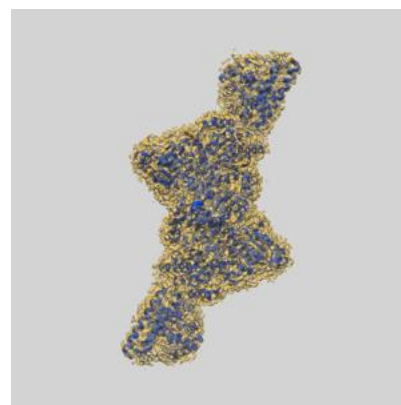
9.1 Map-model overlay [i](#)



X



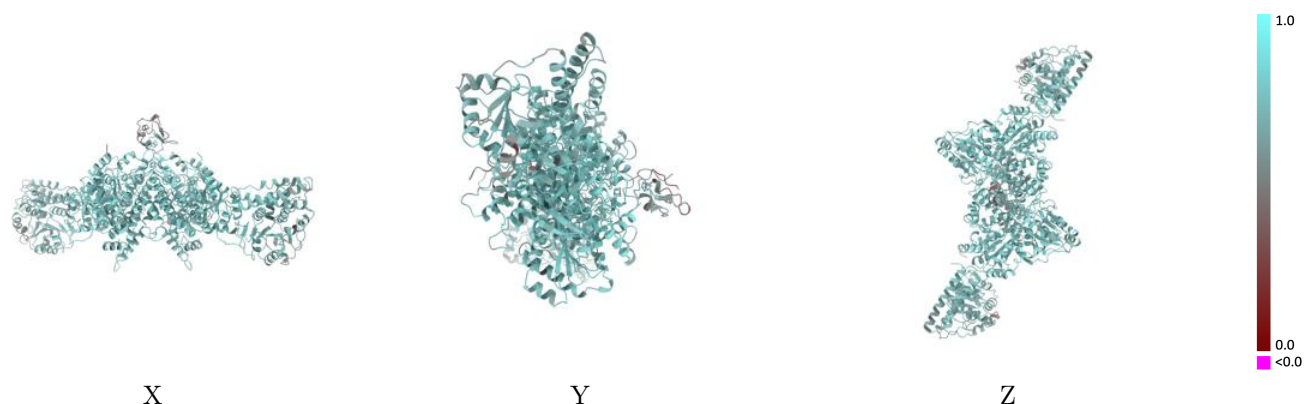
Y



Z

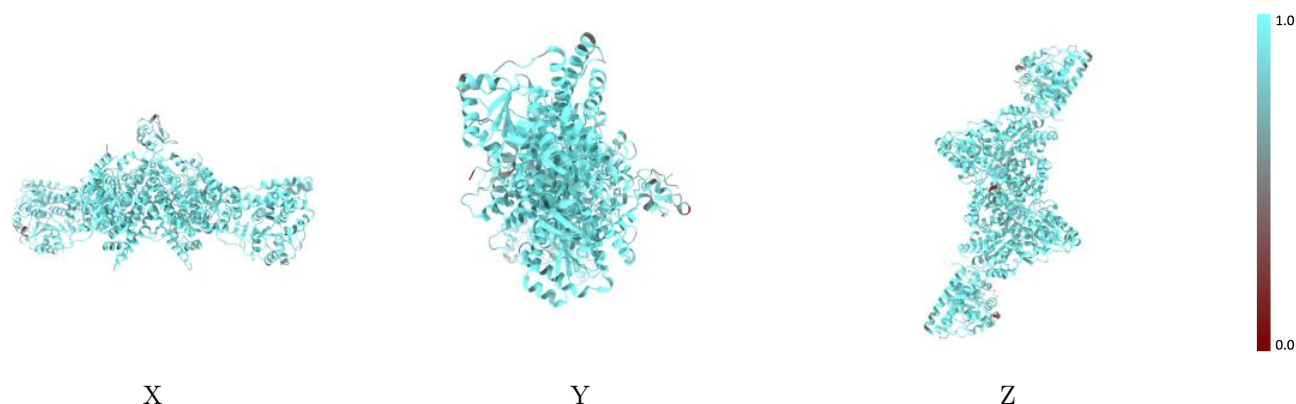
The images above show the 3D surface view of the map at the recommended contour level 0.213 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



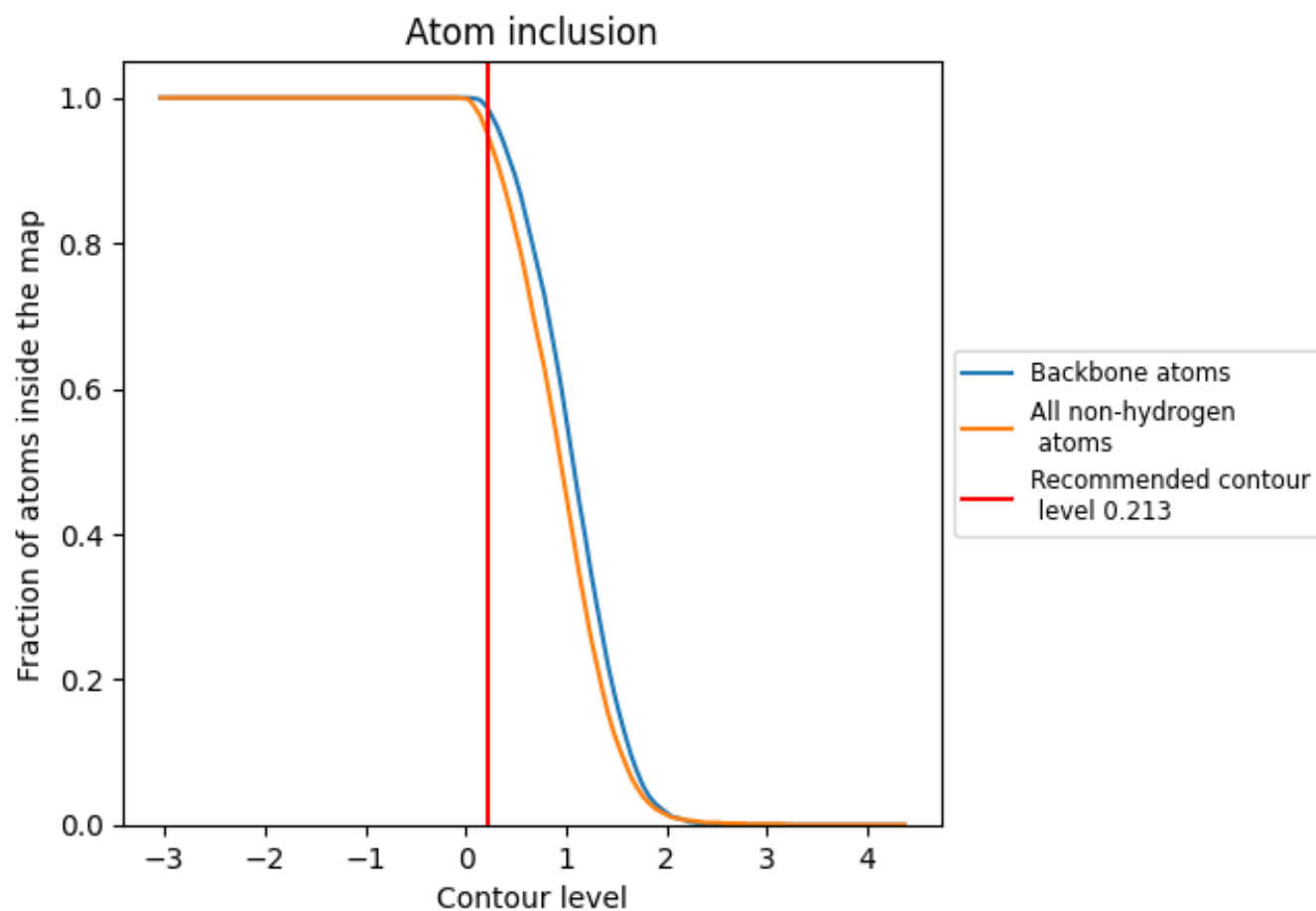
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.213).

9.4 Atom inclusion [i](#)



At the recommended contour level, 99% of all backbone atoms, 95% 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 (0.213) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9520	0.7120
A	0.9220	0.6610
B	0.9740	0.7440
C	0.9760	0.7450
D	0.9190	0.6610
E	0.7410	0.4970

