



wwPDB EM Validation Summary Report ⓘ

Dec 16, 2021 – 09:55 am GMT

PDB ID : 7PXG
EMDB ID : EMD-13702
Title : Verruculogen-bound Drosophila Slo channel
Authors : Raisch, T.; Brockmann, A.; Ebbinghaus-Kintscher, U.; Freigang, J.; Gutbrod, O.; Kubicek, J.; Maertens, B.; Hofnagel, O.; Raunser, S.
Deposited on : 2021-10-08
Resolution : 2.73 Å(reported)
Based on initial model : 5TJ6

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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

EMDB validation analysis : 0.0.0.dev97
Mogul : 1.8.4, CSD as541be (2020)
MolProbity : **FAILED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.24

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 2.73 Å.

There are no overall percentile quality scores available for this entry.

MolProbity failed to run properly - the sequence quality summary graphics cannot be shown.

2 Entry composition [i](#)

There are 7 unique types of molecules in this entry. The entry contains 28664 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 Isoform J of Calcium-activated potassium channel slowpoke.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	885	7023	4548	1148	1279	48	0	0
1	B	885	7023	4548	1148	1279	48	0	0
1	C	885	7023	4548	1148	1279	48	0	0
1	D	885	7023	4548	1148	1279	48	0	0

There are 52 discrepancies between the modelled and reference sequences:

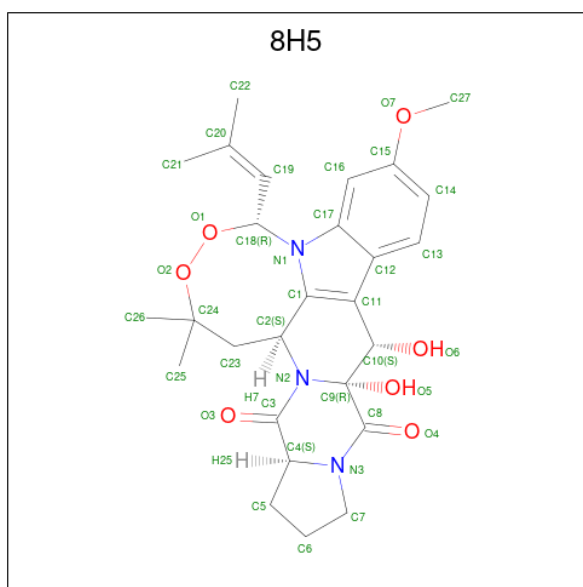
Chain	Residue	Modelled	Actual	Comment	Reference
A	281	ASP	ASN	conflict	UNP Q03720
A	328	ILE	MET	conflict	UNP Q03720
A	332	CYS	SER	conflict	UNP Q03720
A	338	ASP	GLU	conflict	UNP Q03720
A	340	ILE	VAL	conflict	UNP Q03720
A	342	THR	SER	conflict	UNP Q03720
A	343	ARG	GLY	conflict	UNP Q03720
A	344	ALA	ASN	conflict	UNP Q03720
A	349	THR	GLU	conflict	UNP Q03720
A	352	ASN	ARG	conflict	UNP Q03720
A	354	LYS	HIS	conflict	UNP Q03720
A	356	ARG	LYS	conflict	UNP Q03720
A	974	GLY	SER	conflict	UNP Q03720
B	281	ASP	ASN	conflict	UNP Q03720
B	328	ILE	MET	conflict	UNP Q03720
B	332	CYS	SER	conflict	UNP Q03720
B	338	ASP	GLU	conflict	UNP Q03720
B	340	ILE	VAL	conflict	UNP Q03720
B	342	THR	SER	conflict	UNP Q03720
B	343	ARG	GLY	conflict	UNP Q03720
B	344	ALA	ASN	conflict	UNP Q03720
B	349	THR	GLU	conflict	UNP Q03720

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Chain	Residue	Modelled	Actual	Comment	Reference
B	352	ASN	ARG	conflict	UNP Q03720
B	354	LYS	HIS	conflict	UNP Q03720
B	356	ARG	LYS	conflict	UNP Q03720
B	974	GLY	SER	conflict	UNP Q03720
C	281	ASP	ASN	conflict	UNP Q03720
C	328	ILE	MET	conflict	UNP Q03720
C	332	CYS	SER	conflict	UNP Q03720
C	338	ASP	GLU	conflict	UNP Q03720
C	340	ILE	VAL	conflict	UNP Q03720
C	342	THR	SER	conflict	UNP Q03720
C	343	ARG	GLY	conflict	UNP Q03720
C	344	ALA	ASN	conflict	UNP Q03720
C	349	THR	GLU	conflict	UNP Q03720
C	352	ASN	ARG	conflict	UNP Q03720
C	354	LYS	HIS	conflict	UNP Q03720
C	356	ARG	LYS	conflict	UNP Q03720
C	974	GLY	SER	conflict	UNP Q03720
D	281	ASP	ASN	conflict	UNP Q03720
D	328	ILE	MET	conflict	UNP Q03720
D	332	CYS	SER	conflict	UNP Q03720
D	338	ASP	GLU	conflict	UNP Q03720
D	340	ILE	VAL	conflict	UNP Q03720
D	342	THR	SER	conflict	UNP Q03720
D	343	ARG	GLY	conflict	UNP Q03720
D	344	ALA	ASN	conflict	UNP Q03720
D	349	THR	GLU	conflict	UNP Q03720
D	352	ASN	ARG	conflict	UNP Q03720
D	354	LYS	HIS	conflict	UNP Q03720
D	356	ARG	LYS	conflict	UNP Q03720
D	974	GLY	SER	conflict	UNP Q03720

- Molecule 2 is Verruculogen (three-letter code: 8H5) (formula: C₂₇H₃₃N₃O₇) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
			Total	C	N	O	
2	A	1	37	27	3	7	0
2	B	1	37	27	3	7	0
2	C	1	37	27	3	7	0
2	D	1	37	27	3	7	0

- Molecule 3 is CALCIUM ION (three-letter code: CA) (formula: Ca).

Mol	Chain	Residues	Atoms		AltConf
			Total	Ca	
3	A	2	2	2	0
3	B	2	2	2	0
3	C	2	2	2	0
3	D	2	2	2	0

- Molecule 4 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

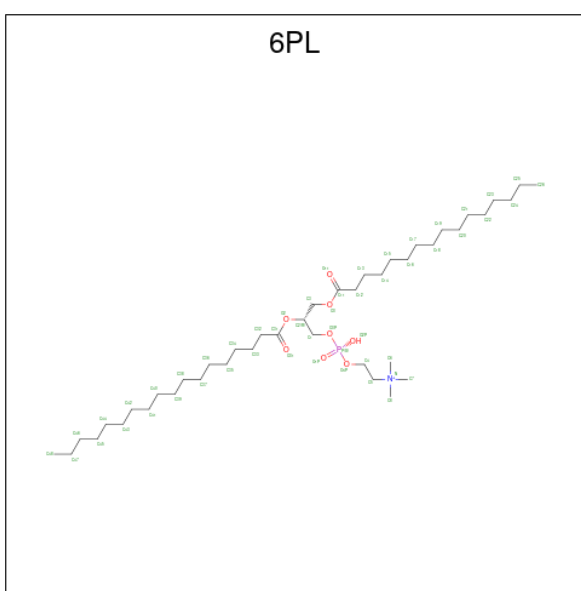
Mol	Chain	Residues	Atoms		AltConf
			Total	Mg	
4	A	1	1	1	0

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Mol	Chain	Residues	Atoms		AltConf
4	B	1	Total	Mg	0
			1	1	
4	C	1	Total	Mg	0
			1	1	
4	D	1	Total	Mg	0
			1	1	

- Molecule 5 is (4S,7R)-4-HYDROXY-N,N,N-TRIMETHYL-9-OXO-7-[(PALMITOYLOXY) METHYL]-3,5,8-TRIOXA-4-PHOSPHAHEXACOSAN-1-AMINIUM 4-OXIDE (three-letter code: 6PL) (formula: C₄₂H₈₅NO₈P).



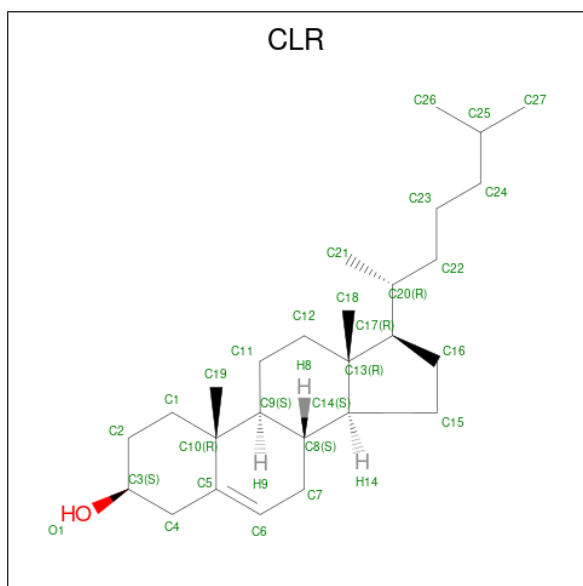
Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
5	A	1	Total	C	N	O	P	0
			35	25	1	8	1	
5	B	1	Total	C	N	O	P	0
			74	54	2	16	2	
5	B	1	Total	C	N	O	P	0
			74	54	2	16	2	
5	C	1	Total	C	N	O	P	0
			113	83	3	24	3	
5	C	1	Total	C	N	O	P	0
			113	83	3	24	3	
5	C	1	Total	C	N	O	P	0
			113	83	3	24	3	
5	D	1	Total	C	N	O	P	0
			74	54	2	16	2	

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Mol	Chain	Residues	Atoms				AltConf	
			Total	C	N	O		P
5	D	1	74	54	2	16	2	0

- Molecule 6 is CHOLESTEROL (three-letter code: CLR) (formula: $C_{27}H_{46}O$).



Mol	Chain	Residues	Atoms			AltConf
			Total	C	O	
6	A	1	28	27	1	0
6	B	1	28	27	1	0
6	C	1	28	27	1	0
6	D	1	28	27	1	0

- Molecule 7 is POTASSIUM ION (three-letter code: K) (formula: K).

Mol	Chain	Residues	Atoms		AltConf
			Total	K	
7	A	4	4	4	0

MolProbity failed to run properly - this section is therefore empty.

3 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C4	Depositor
Number of particles used	66077	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 ($e^-/\text{\AA}^2$)	80.4	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	39.447	Depositor
Minimum map value	-18.229	Depositor
Average map value	-0.000	Depositor
Map value standard deviation	1.000	Depositor
Recommended contour level	5.0	Depositor
Map size (Å)	313.6, 313.6, 313.6	wwPDB
Map dimensions	448, 448, 448	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.7, 0.7, 0.7	Depositor

4 Model quality [i](#)

4.1 Standard geometry [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.2 Too-close contacts [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

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4.3.2 Protein sidechains [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3.3 RNA [i](#)

MolProbity failed to run properly - this section is therefore empty.

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

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

4.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

4.6 Ligand geometry [i](#)

Of 32 ligands modelled in this entry, 16 are monoatomic - leaving 16 for Mogul analysis.

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
6	CLR	A	2006	-	31,31,31	0.52	0	48,48,48	1.26	4 (8%)
2	8H5	A	2001	-	30,42,42	0.99	2 (6%)	32,68,68	1.26	2 (6%)
5	6PL	A	2005	-	34,34,51	1.44	4 (11%)	40,42,59	1.09	2 (5%)
6	CLR	D	2007	-	31,31,31	0.51	0	48,48,48	1.27	4 (8%)
5	6PL	B	2006	-	34,34,51	1.44	4 (11%)	40,42,59	1.09	2 (5%)
5	6PL	D	2006	-	34,34,51	1.44	4 (11%)	40,42,59	1.09	2 (5%)
5	6PL	C	1207	-	34,34,51	1.44	4 (11%)	40,42,59	1.09	2 (5%)
6	CLR	B	2007	-	31,31,31	0.52	0	48,48,48	1.26	4 (8%)
5	6PL	C	1201	-	38,38,51	1.39	4 (10%)	44,46,59	1.08	2 (4%)
2	8H5	B	2001	-	30,42,42	1.00	2 (6%)	32,68,68	1.26	2 (6%)
5	6PL	D	2005	-	38,38,51	1.39	4 (10%)	44,46,59	1.08	2 (4%)
6	CLR	C	1208	-	31,31,31	0.52	0	48,48,48	1.26	4 (8%)
2	8H5	D	2001	-	30,42,42	1.01	2 (6%)	32,68,68	1.28	2 (6%)
2	8H5	C	1202	-	30,42,42	1.00	2 (6%)	32,68,68	1.27	2 (6%)
5	6PL	C	1206	-	38,38,51	1.39	4 (10%)	44,46,59	1.08	2 (4%)
5	6PL	B	2005	-	38,38,51	1.39	4 (10%)	44,46,59	1.08	2 (4%)

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
6	CLR	A	2006	-	-	7/10/68/68	0/4/4/4
2	8H5	A	2001	-	-	0/4/73/73	0/5/6/6
5	6PL	A	2005	-	-	15/38/38/55	-
6	CLR	D	2007	-	-	7/10/68/68	0/4/4/4
5	6PL	B	2006	-	-	15/38/38/55	-
5	6PL	D	2006	-	-	15/38/38/55	-
5	6PL	C	1207	-	-	15/38/38/55	-
6	CLR	B	2007	-	-	7/10/68/68	0/4/4/4
5	6PL	C	1201	-	-	15/42/42/55	-
2	8H5	B	2001	-	-	0/4/73/73	0/5/6/6

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	6PL	D	2005	-	-	15/42/42/55	-
6	CLR	C	1208	-	-	7/10/68/68	0/4/4/4
2	8H5	D	2001	-	-	0/4/73/73	0/5/6/6
2	8H5	C	1202	-	-	0/4/73/73	0/5/6/6
5	6PL	C	1206	-	-	15/42/42/55	-
5	6PL	B	2005	-	-	15/42/42/55	-

The worst 5 of 40 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
5	C	1207	6PL	O3-C11	3.91	1.44	1.33
5	A	2005	6PL	O3-C11	3.91	1.44	1.33
5	B	2005	6PL	O3-C11	3.89	1.44	1.33
5	D	2006	6PL	O3-C11	3.89	1.44	1.33
5	B	2006	6PL	O3-C11	3.89	1.44	1.33

The worst 5 of 40 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	2001	8H5	C11-C1-N1	5.18	109.95	106.66
2	C	1202	8H5	C11-C1-N1	5.16	109.94	106.66
2	A	2001	8H5	C11-C1-N1	5.13	109.92	106.66
2	B	2001	8H5	C11-C1-N1	5.10	109.90	106.66
5	C	1207	6PL	O2-C31-C32	4.01	120.15	111.50

There are no chirality outliers.

5 of 148 torsion outliers are listed below:

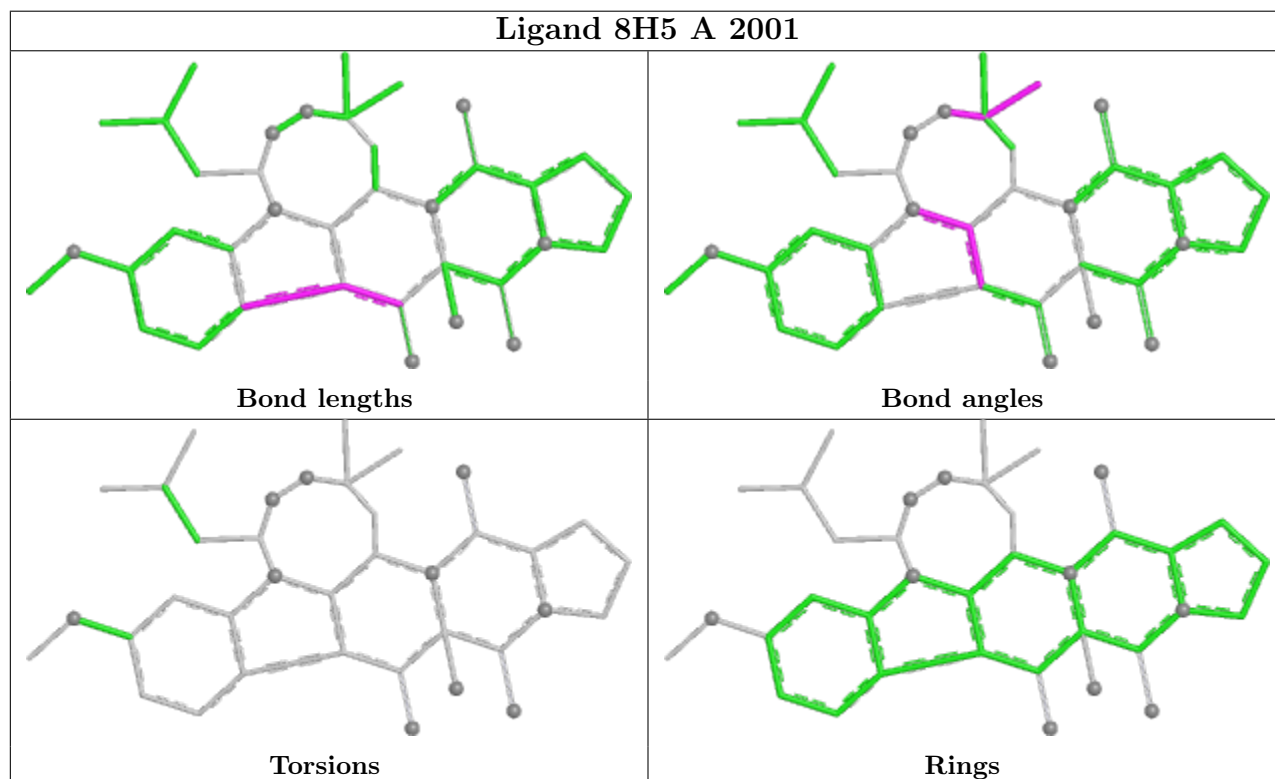
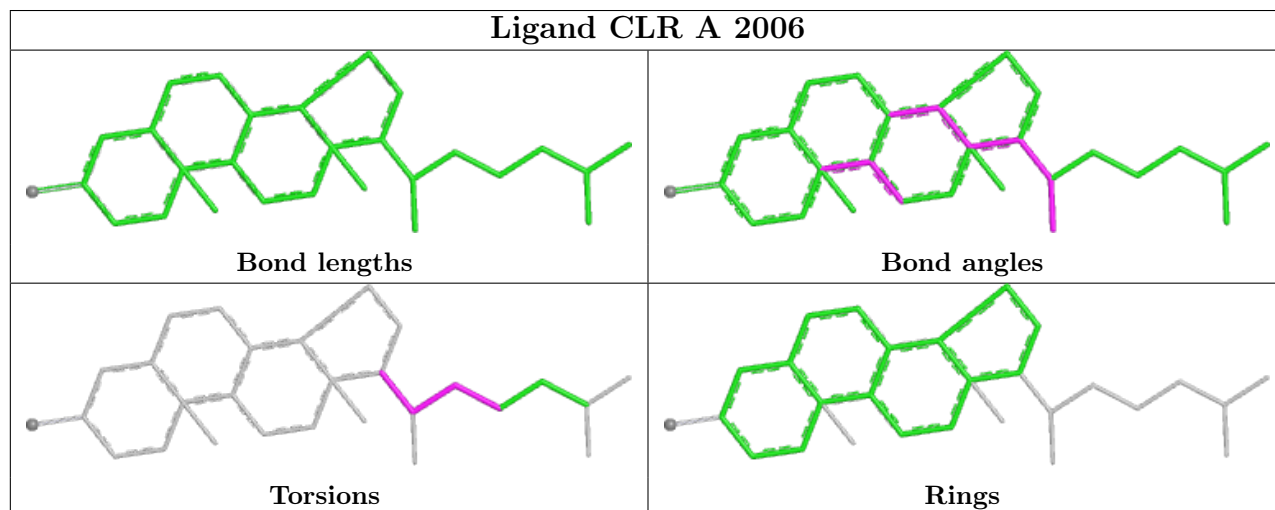
Mol	Chain	Res	Type	Atoms
5	A	2005	6PL	O2-C2-C3-O3
5	A	2005	6PL	C1-O3P-P-O1P
5	B	2005	6PL	C4-O4P-P-O3P
5	B	2005	6PL	C4-O4P-P-O1P
5	B	2005	6PL	C4-O4P-P-O2P

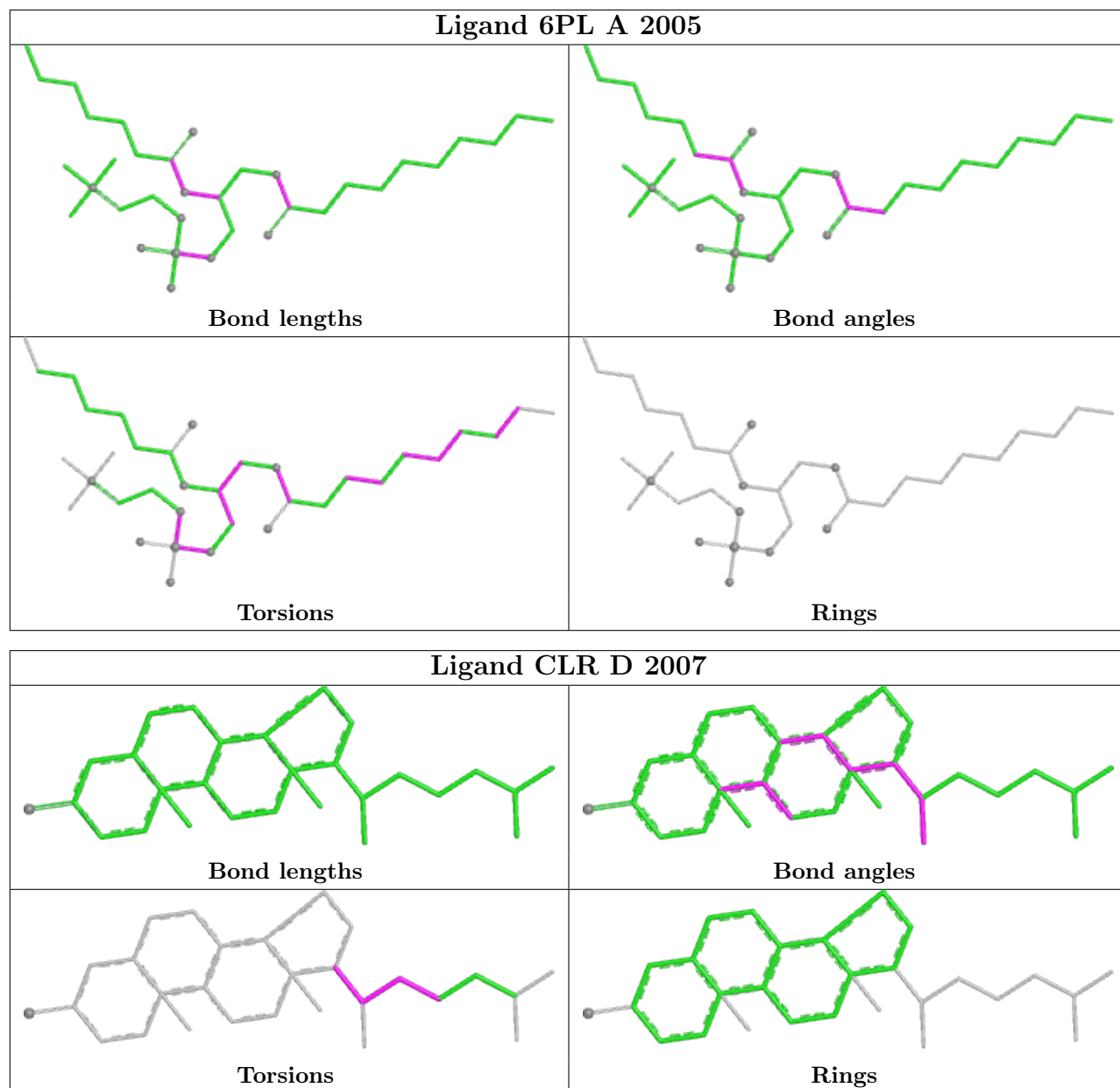
There are no ring outliers.

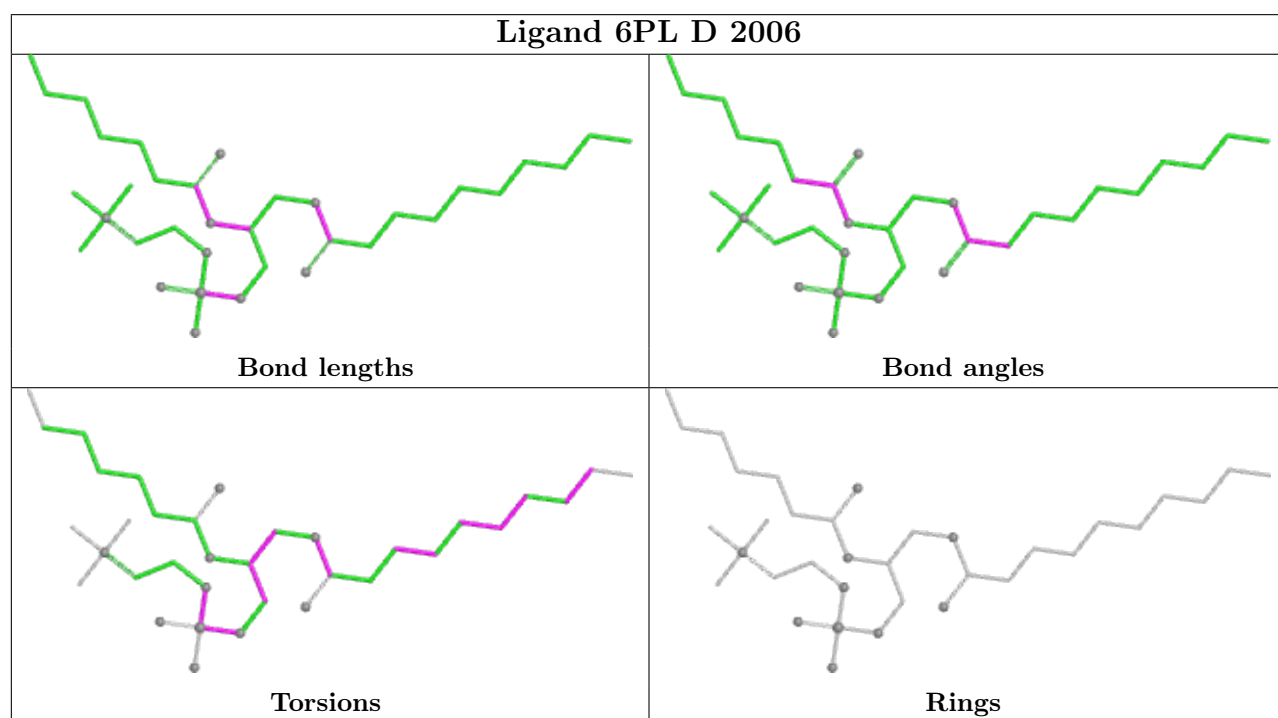
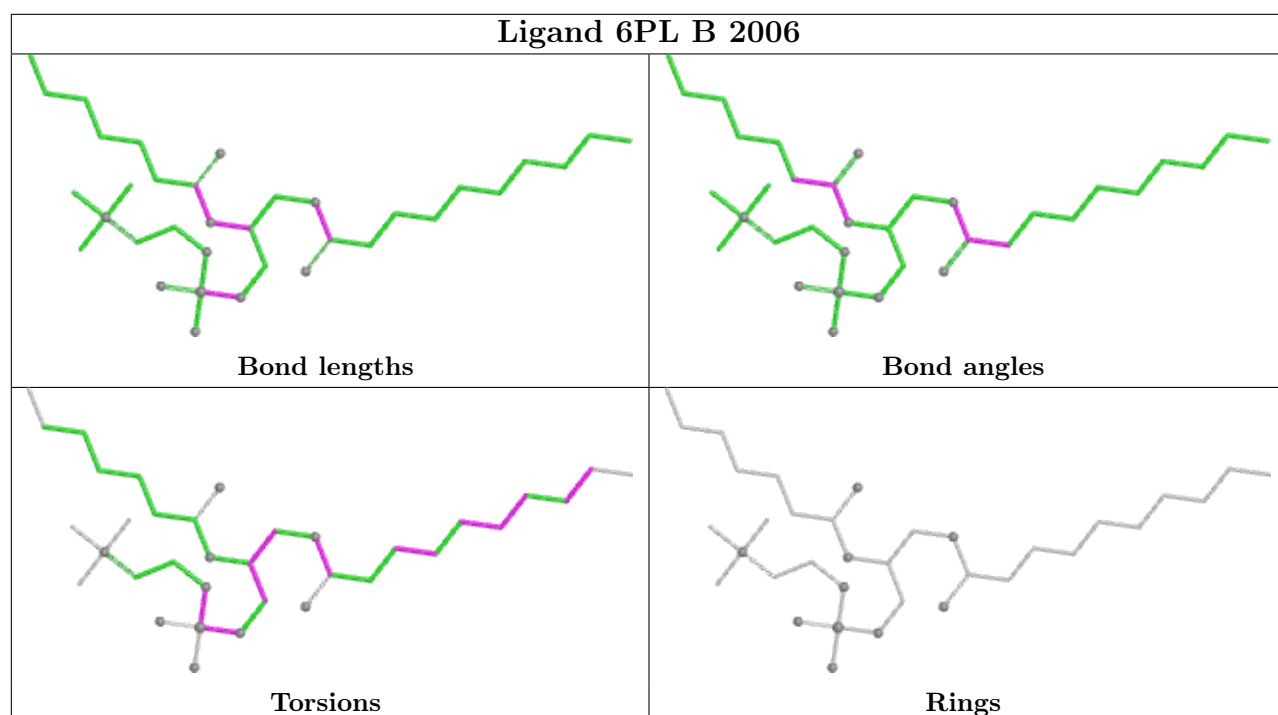
No monomer is involved in short contacts.

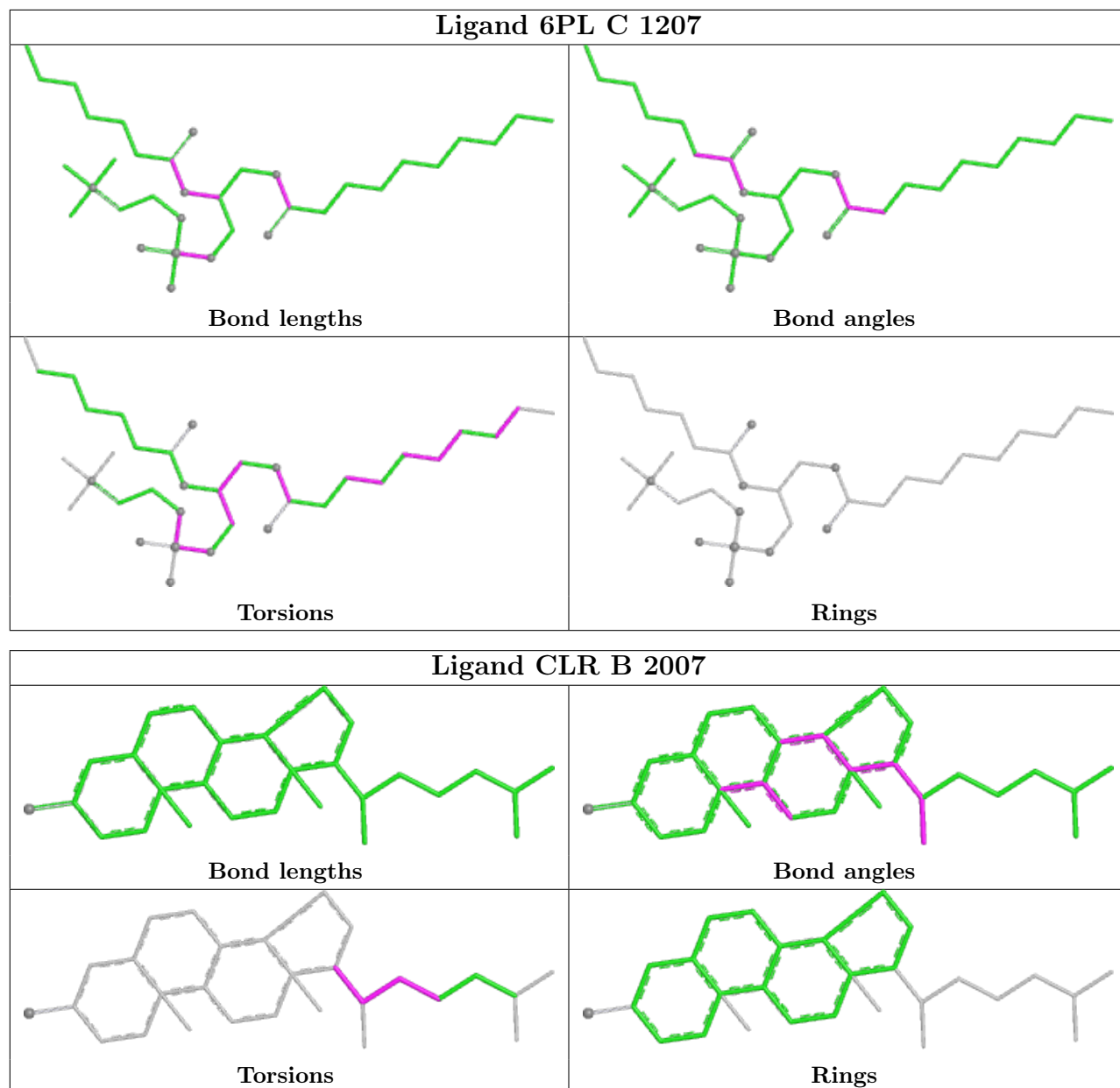
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will

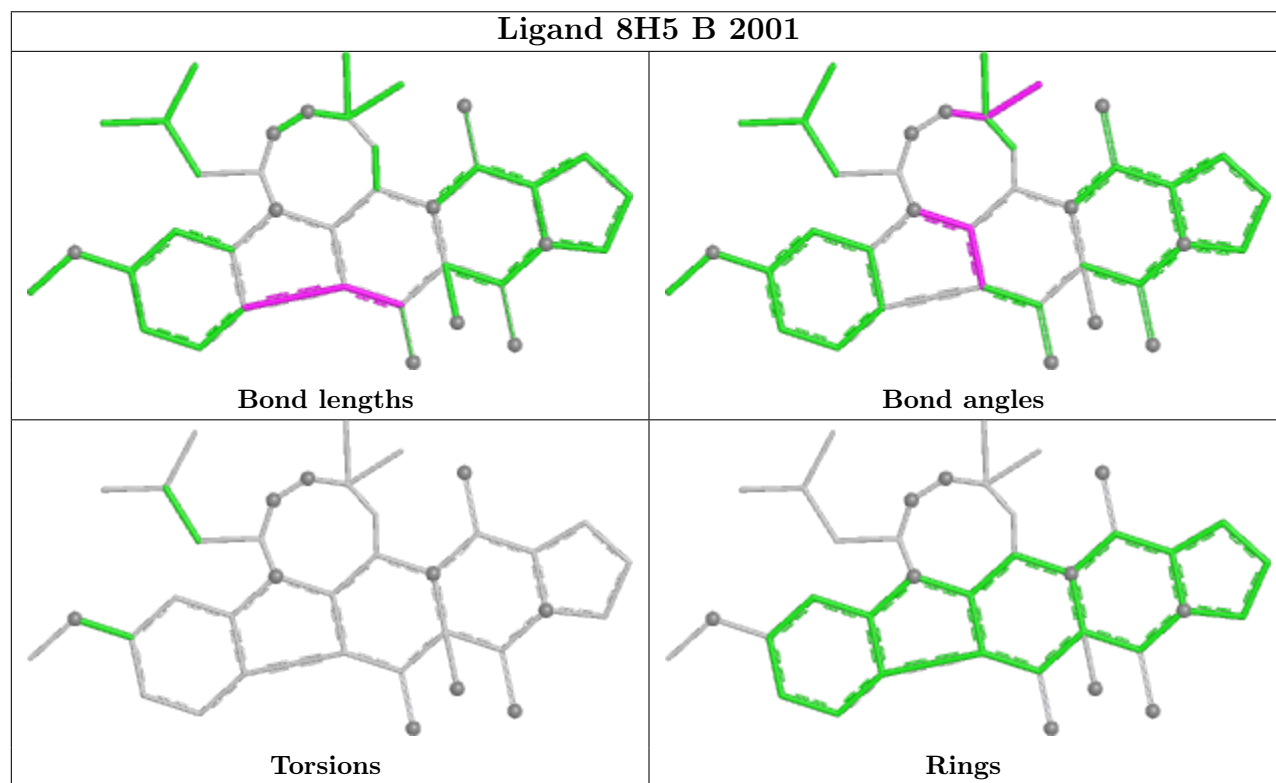
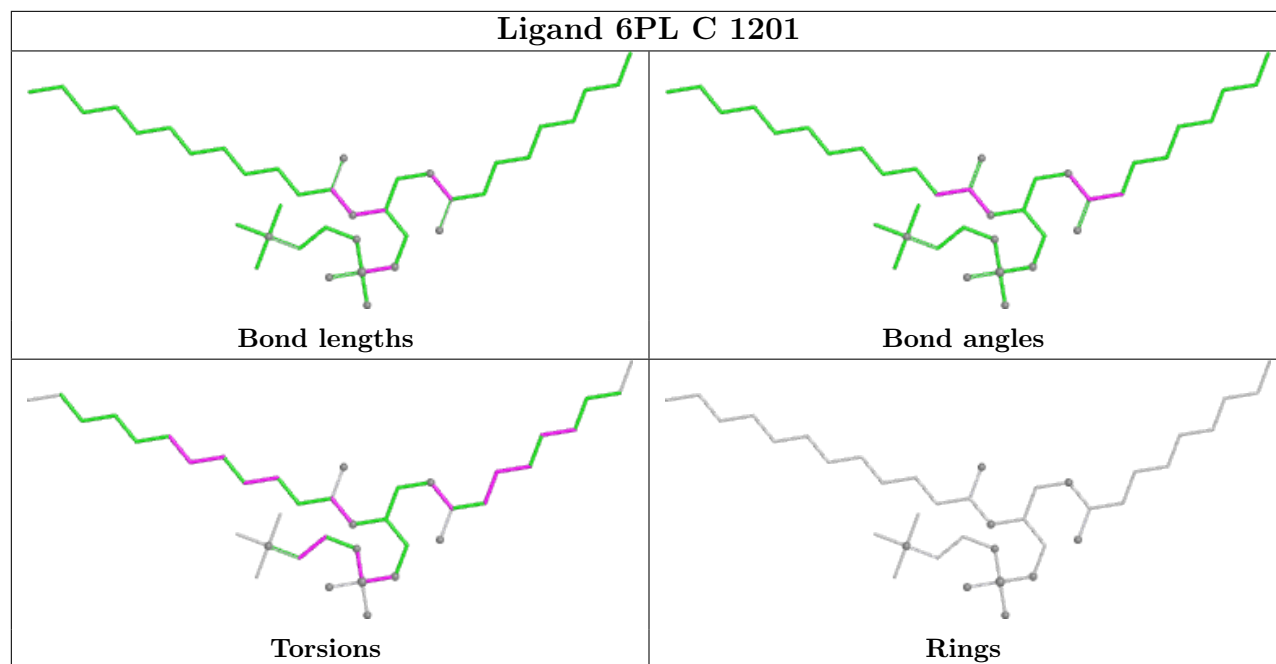
also be included. For torsion angles, if less than 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted 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.

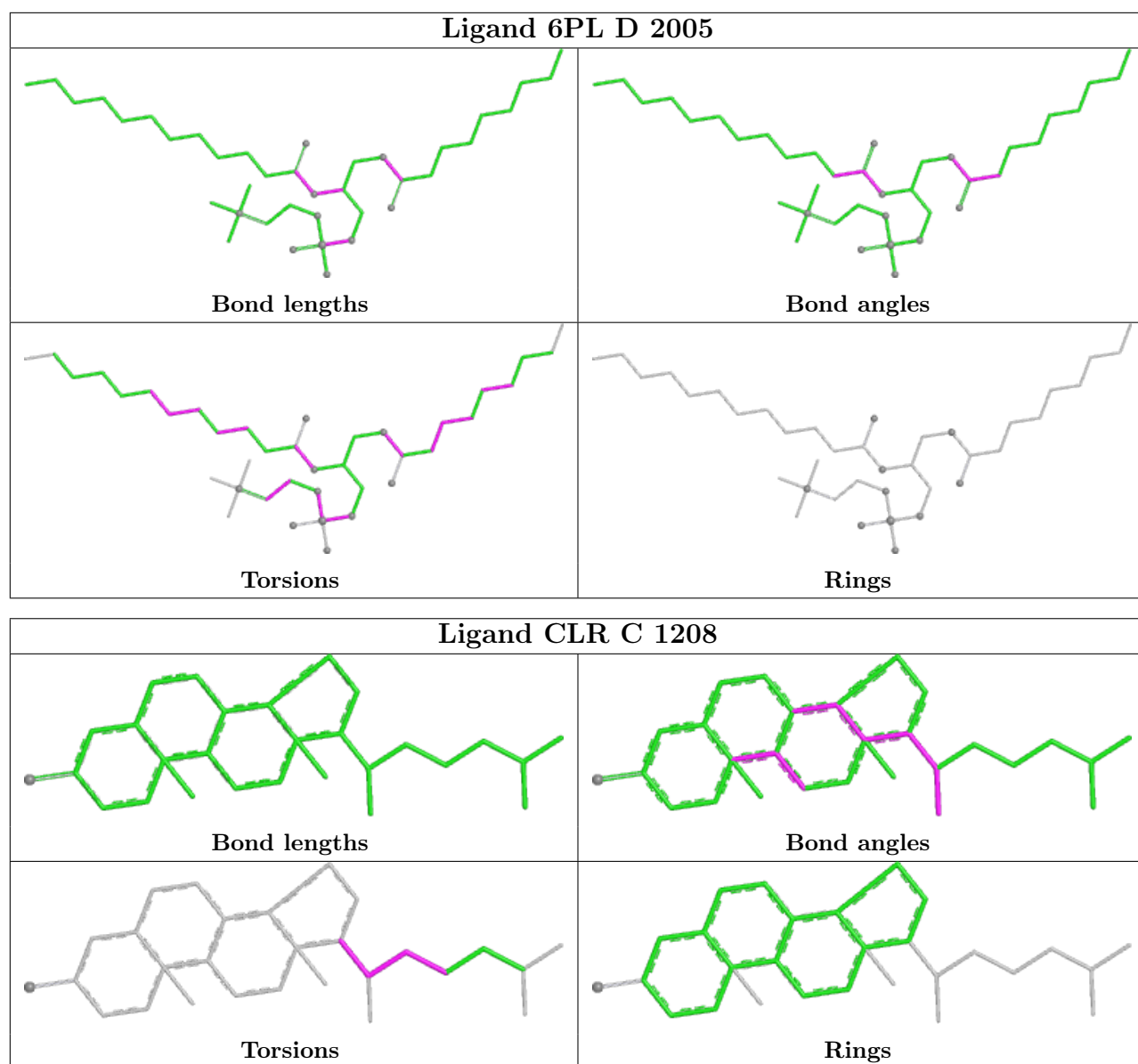


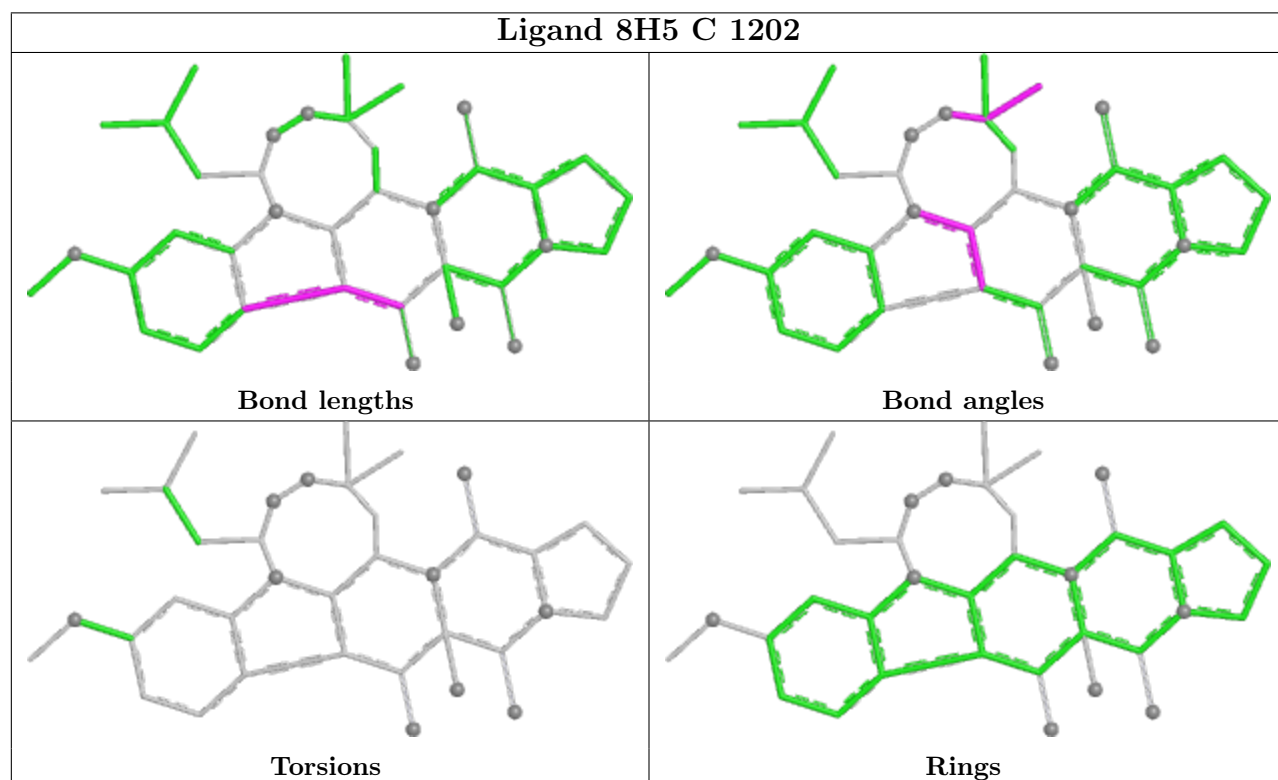
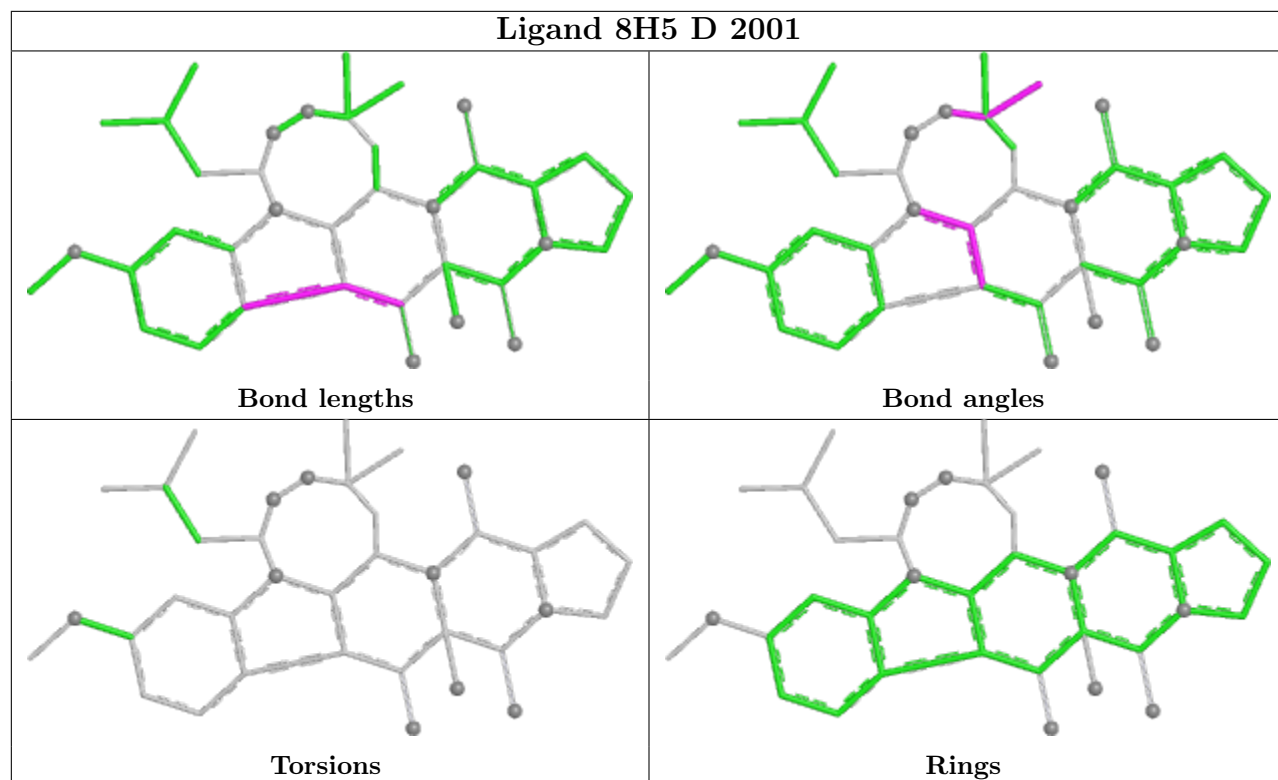


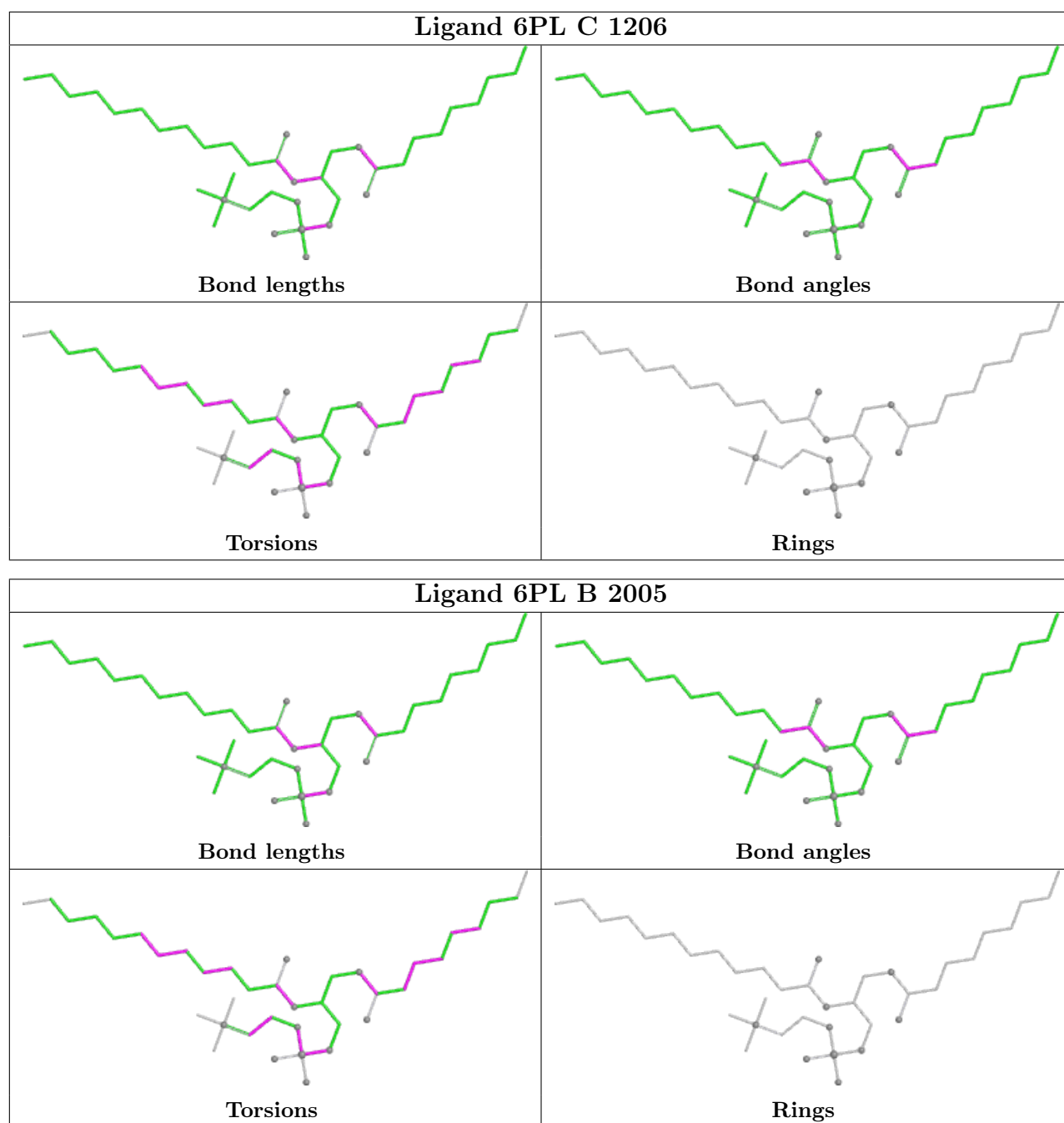












4.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

4.8 Polymer linkage issues [\(i\)](#)

There are no chain breaks in this entry.

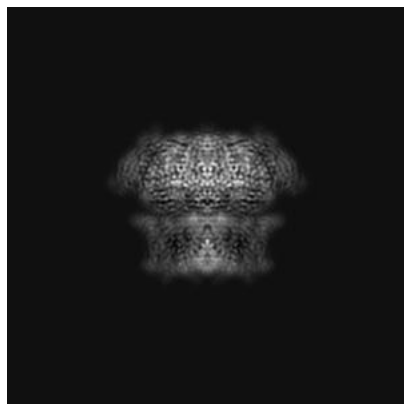
5 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-13702. These allow visual inspection of the internal detail of the map and identification of artifacts.

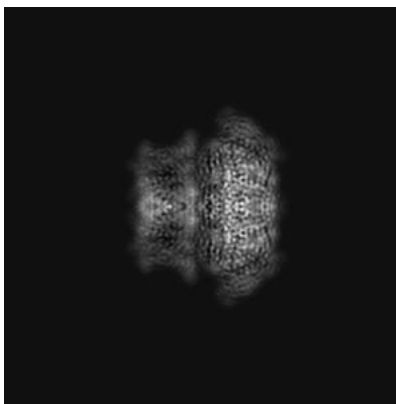
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

5.1 Orthogonal projections [i](#)

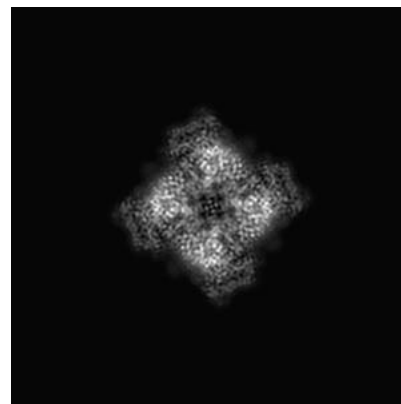
5.1.1 Primary map



X

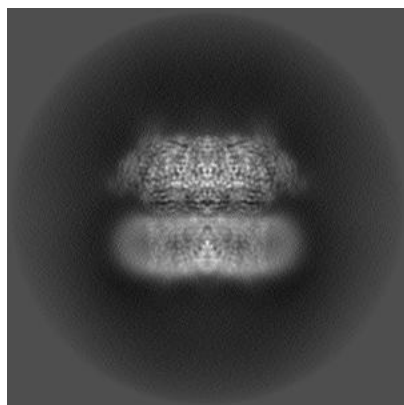


Y

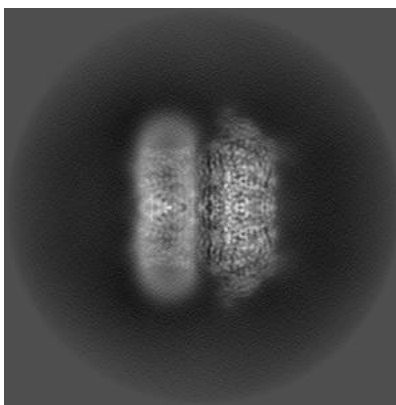


Z

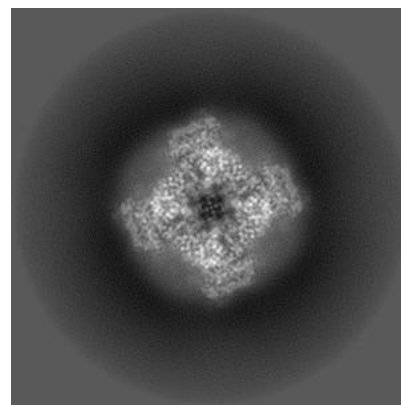
5.1.2 Raw map



X



Y

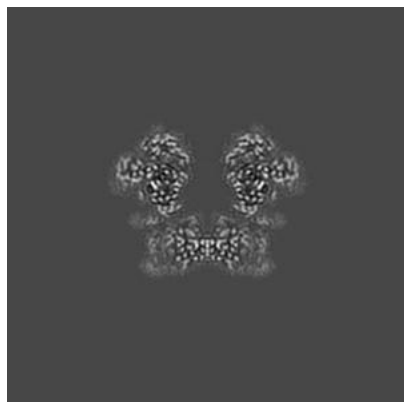


Z

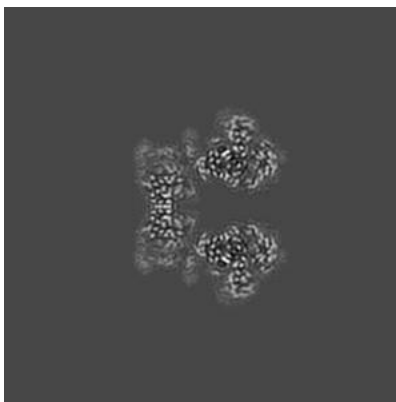
The images above show the map projected in three orthogonal directions.

5.2 Central slices [i](#)

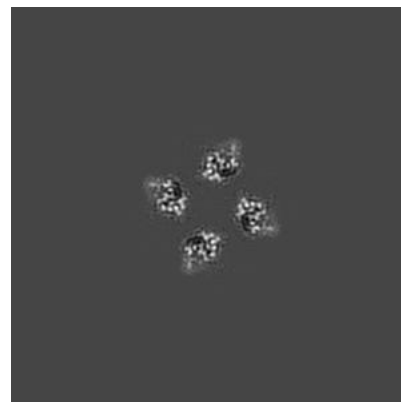
5.2.1 Primary map



X Index: 224

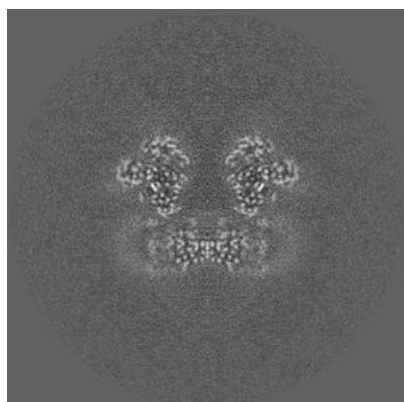


Y Index: 224

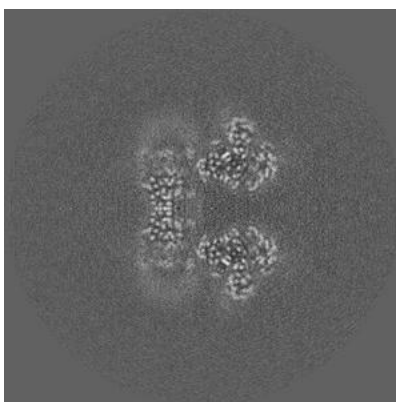


Z Index: 224

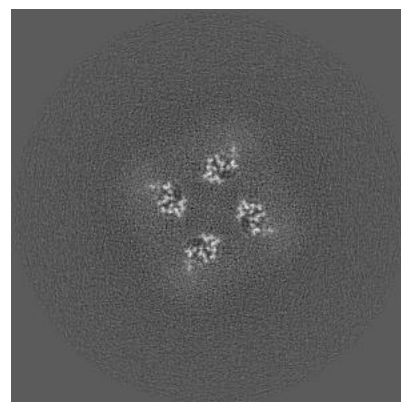
5.2.2 Raw map



X Index: 224



Y Index: 224

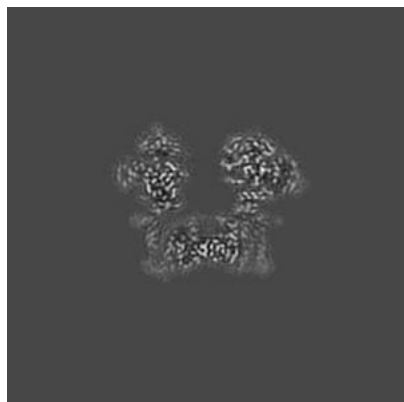


Z Index: 224

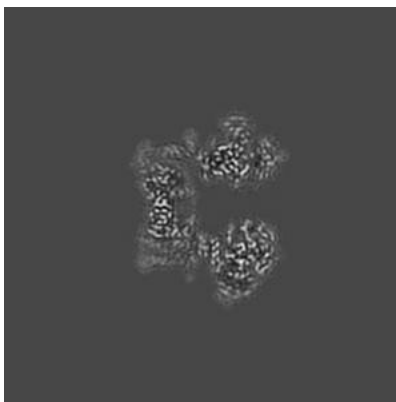
The images above show central slices of the map in three orthogonal directions.

5.3 Largest variance slices [i](#)

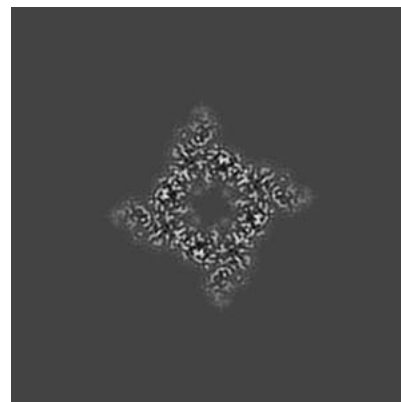
5.3.1 Primary map



X Index: 220

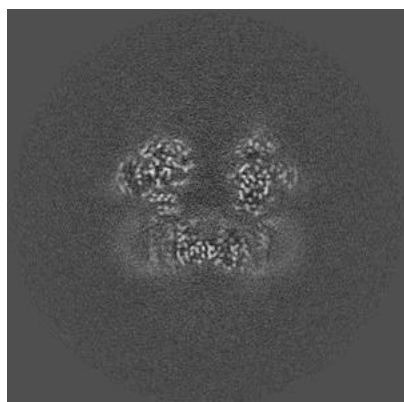


Y Index: 220

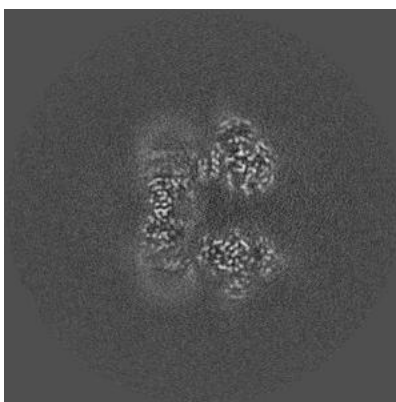


Z Index: 252

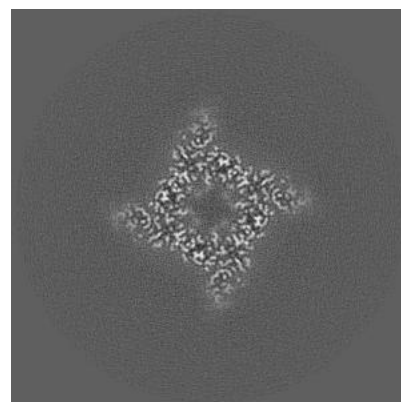
5.3.2 Raw map



X Index: 228



Y Index: 228

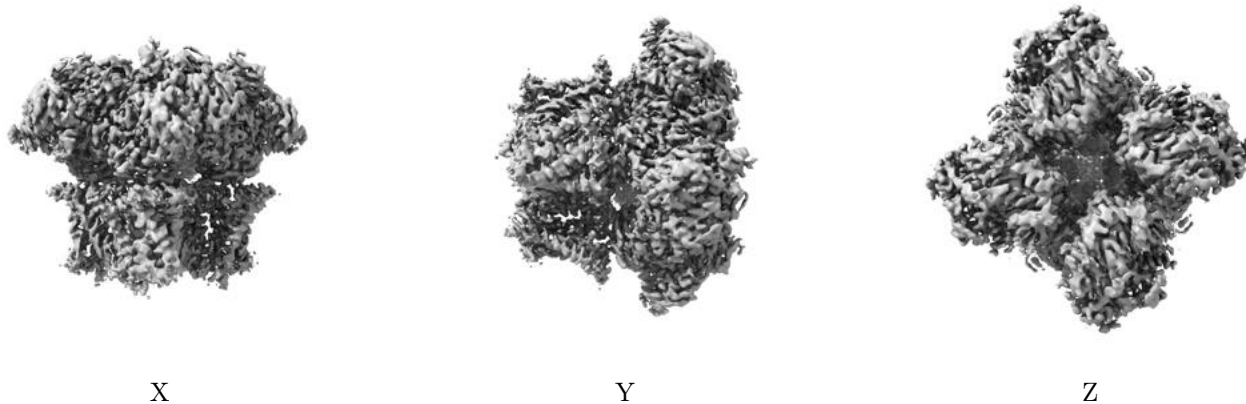


Z Index: 252

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

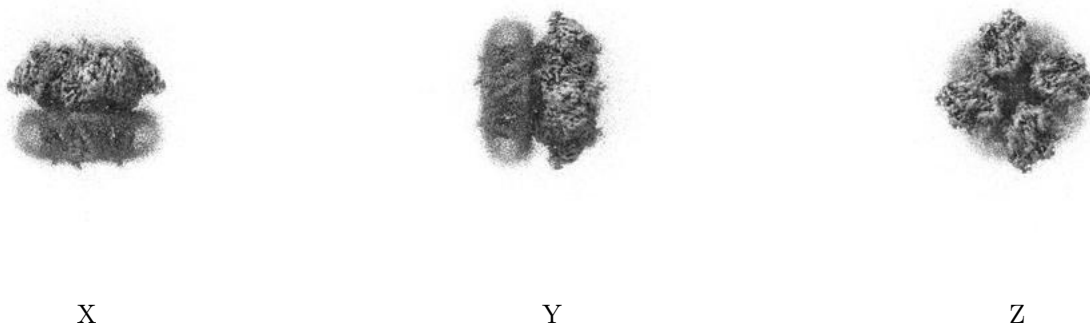
5.4 Orthogonal surface views [i](#)

5.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 5.0. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

5.4.2 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.

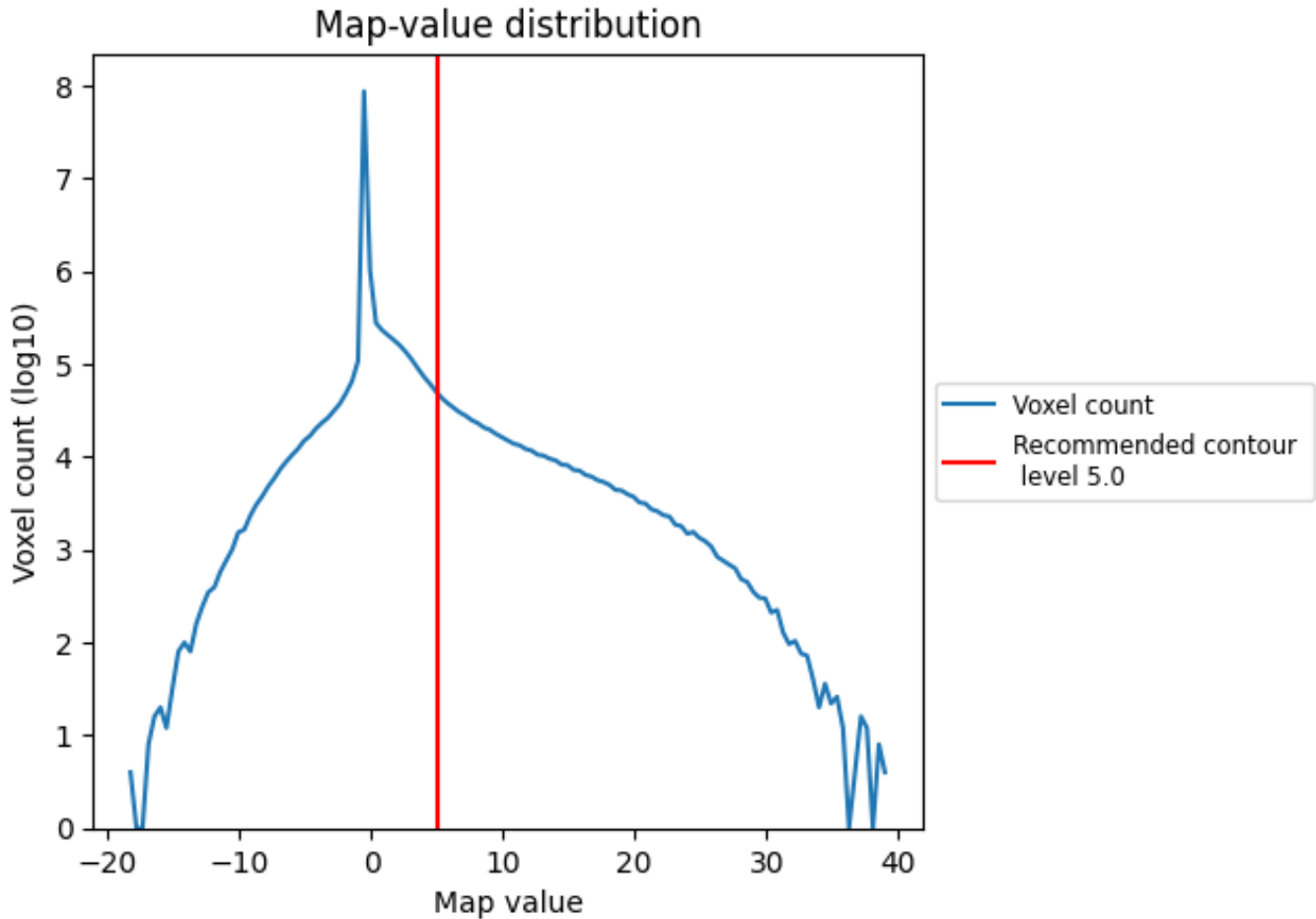
5.5 Mask visualisation [i](#)

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

6 Map analysis [i](#)

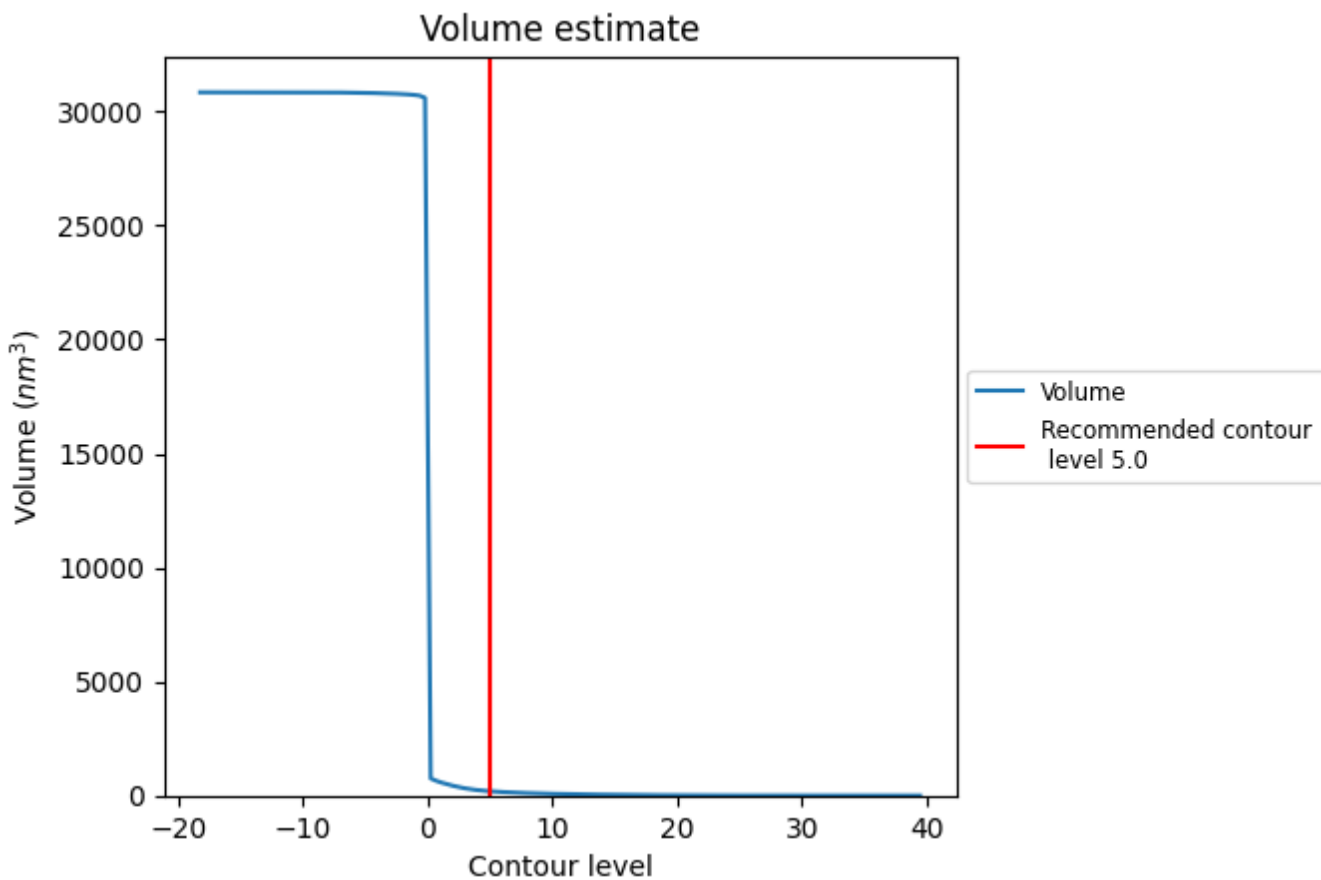
This section contains the results of statistical analysis of the map.

6.1 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.

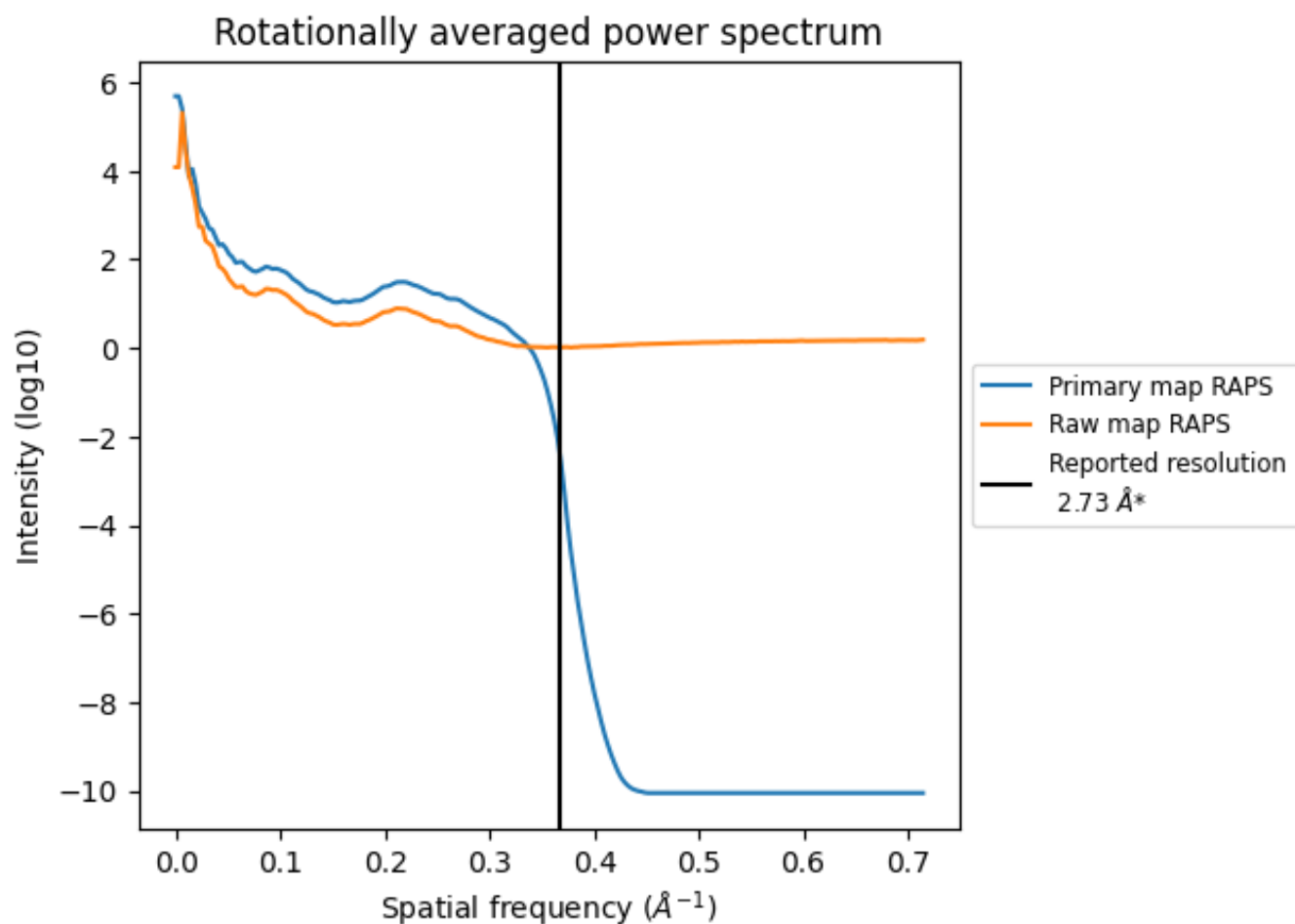
6.2 Volume estimate [\(i\)](#)



The volume at the recommended contour level is 189 nm³; this corresponds to an approximate mass of 171 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.

6.3 Rotationally averaged power spectrum [i](#)

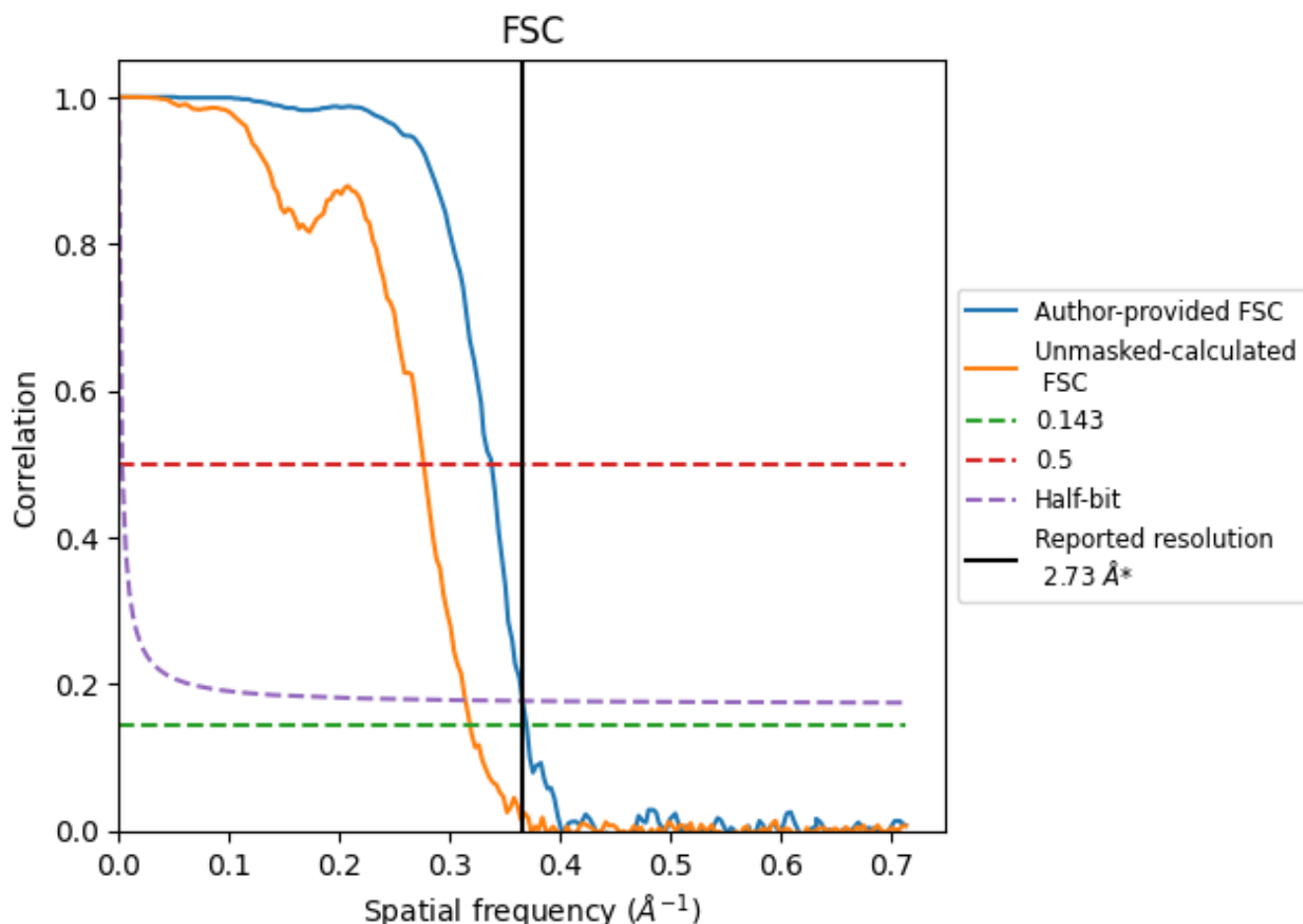


*Reported resolution corresponds to spatial frequency of 0.366 Å⁻¹

7 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.

7.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.366 Å⁻¹

7.2 Resolution estimates [i](#)

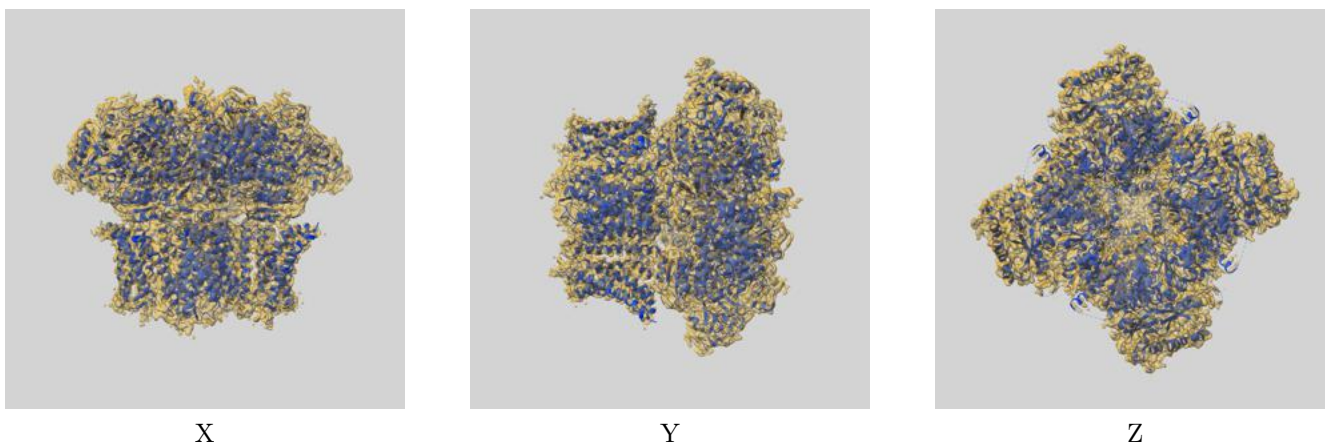
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.73	-	-
Author-provided FSC curve	2.70	2.96	2.73
Unmasked-calculated*	3.14	3.61	3.18

*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 3.14 differs from the reported value 2.73 by more than 10 %

8 Map-model fit [i](#)

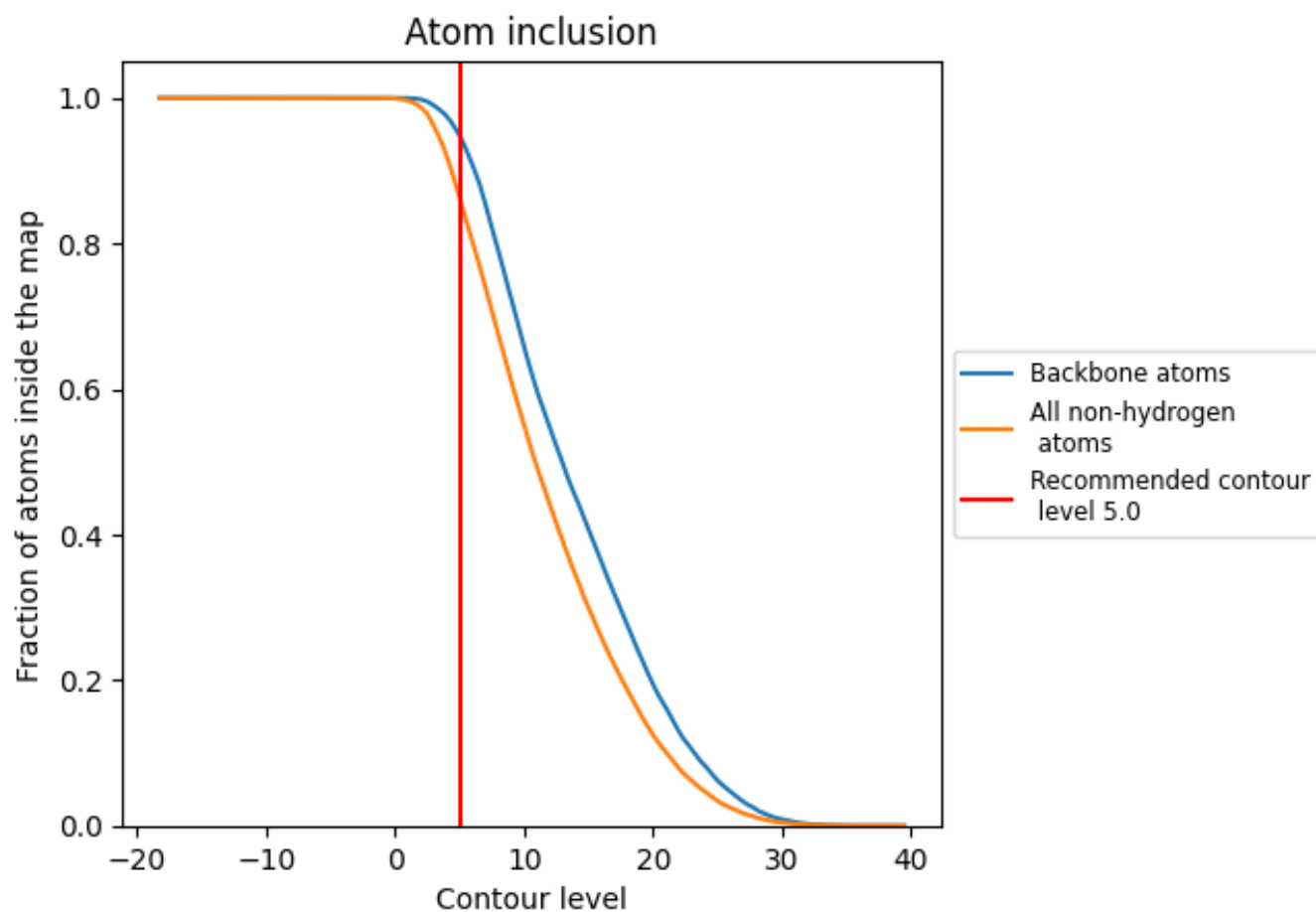
This section contains information regarding the fit between EMDB map EMD-13702 and PDB model 7PXG. Per-residue inclusion information can be found in section ?? on page ??.

8.1 Map-model overlay [i](#)



The images above show the 3D surface view of the map at the recommended contour level 5.0 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

8.2 Atom inclusion [i](#)



At the recommended contour level, 95% of all backbone atoms, 86% of all non-hydrogen atoms, are inside the map.