



## Full wwPDB EM Validation Report ⓘ

Dec 11, 2022 – 04:13 pm GMT

PDB ID : 6SP2  
EMDB ID : EMD-10279  
Title : CryoEM structure of SERINC from *Drosophila melanogaster*  
Authors : Pye, V.E.; Nans, A.; Cherepanov, P.  
Deposited on : 2019-08-30  
Resolution : 3.33 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](mailto: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>.

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

EMDB validation analysis : 0.0.1.dev43  
Mogul : 1.8.4, CSD as541be (2020)  
MolProbity : 4.02b-467  
buster-report : 1.1.7 (2018)  
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)  
MapQ : 1.9.9  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.31.3

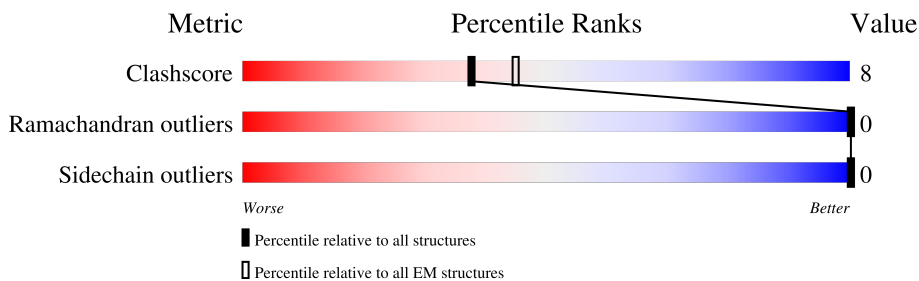
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*ELECTRON MICROSCOPY*

The reported resolution of this entry is 3.33 Å.

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



Metric	Whole archive (#Entries)	EM structures (#Entries)
Clashscore	158937	4297
Ramachandran outliers	154571	4023
Sidechain outliers	154315	3826

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 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	513	
1	B	513	
1	C	513	
1	D	513	
1	E	513	
1	F	513	

## 2 Entry composition i

There are 4 unique types of molecules in this entry. The entry contains 18498 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 Membrane protein TMS1d.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
1	A	365	2834	1872	439	497	26	0	0
1	B	365	2834	1872	439	497	26	0	0
1	C	365	2834	1872	439	497	26	0	0
1	D	365	2834	1872	439	497	26	0	0
1	E	365	2834	1872	439	497	26	0	0
1	F	365	2834	1872	439	497	26	0	0

There are 288 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	466	THR	-	expression tag	UNP Q9U6P4
A	467	GLY	-	expression tag	UNP Q9U6P4
A	468	GLY	-	expression tag	UNP Q9U6P4
A	469	SER	-	expression tag	UNP Q9U6P4
A	470	SER	-	expression tag	UNP Q9U6P4
A	471	GLY	-	expression tag	UNP Q9U6P4
A	472	LEU	-	expression tag	UNP Q9U6P4
A	473	GLU	-	expression tag	UNP Q9U6P4
A	474	VAL	-	expression tag	UNP Q9U6P4
A	475	LEU	-	expression tag	UNP Q9U6P4
A	476	PHE	-	expression tag	UNP Q9U6P4
A	477	GLN	-	expression tag	UNP Q9U6P4
A	478	GLY	-	expression tag	UNP Q9U6P4
A	479	PRO	-	expression tag	UNP Q9U6P4
A	480	GLY	-	expression tag	UNP Q9U6P4
A	481	SER	-	expression tag	UNP Q9U6P4
A	482	GLY	-	expression tag	UNP Q9U6P4
A	483	GLY	-	expression tag	UNP Q9U6P4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	484	SER	-	expression tag	UNP Q9U6P4
A	485	ALA	-	expression tag	UNP Q9U6P4
A	486	TRP	-	expression tag	UNP Q9U6P4
A	487	SER	-	expression tag	UNP Q9U6P4
A	488	HIS	-	expression tag	UNP Q9U6P4
A	489	PRO	-	expression tag	UNP Q9U6P4
A	490	GLN	-	expression tag	UNP Q9U6P4
A	491	PHE	-	expression tag	UNP Q9U6P4
A	492	GLU	-	expression tag	UNP Q9U6P4
A	493	LYS	-	expression tag	UNP Q9U6P4
A	494	GLY	-	expression tag	UNP Q9U6P4
A	495	GLY	-	expression tag	UNP Q9U6P4
A	496	GLY	-	expression tag	UNP Q9U6P4
A	497	SER	-	expression tag	UNP Q9U6P4
A	498	GLY	-	expression tag	UNP Q9U6P4
A	499	GLY	-	expression tag	UNP Q9U6P4
A	500	GLY	-	expression tag	UNP Q9U6P4
A	501	SER	-	expression tag	UNP Q9U6P4
A	502	GLY	-	expression tag	UNP Q9U6P4
A	503	GLY	-	expression tag	UNP Q9U6P4
A	504	SER	-	expression tag	UNP Q9U6P4
A	505	ALA	-	expression tag	UNP Q9U6P4
A	506	TRP	-	expression tag	UNP Q9U6P4
A	507	SER	-	expression tag	UNP Q9U6P4
A	508	HIS	-	expression tag	UNP Q9U6P4
A	509	PRO	-	expression tag	UNP Q9U6P4
A	510	GLN	-	expression tag	UNP Q9U6P4
A	511	PHE	-	expression tag	UNP Q9U6P4
A	512	GLU	-	expression tag	UNP Q9U6P4
A	513	LYS	-	expression tag	UNP Q9U6P4
B	466	THR	-	expression tag	UNP Q9U6P4
B	467	GLY	-	expression tag	UNP Q9U6P4
B	468	GLY	-	expression tag	UNP Q9U6P4
B	469	SER	-	expression tag	UNP Q9U6P4
B	470	SER	-	expression tag	UNP Q9U6P4
B	471	GLY	-	expression tag	UNP Q9U6P4
B	472	LEU	-	expression tag	UNP Q9U6P4
B	473	GLU	-	expression tag	UNP Q9U6P4
B	474	VAL	-	expression tag	UNP Q9U6P4
B	475	LEU	-	expression tag	UNP Q9U6P4
B	476	PHE	-	expression tag	UNP Q9U6P4
B	477	GLN	-	expression tag	UNP Q9U6P4

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Chain	Residue	Modelled	Actual	Comment	Reference
B	478	GLY	-	expression tag	UNP Q9U6P4
B	479	PRO	-	expression tag	UNP Q9U6P4
B	480	GLY	-	expression tag	UNP Q9U6P4
B	481	SER	-	expression tag	UNP Q9U6P4
B	482	GLY	-	expression tag	UNP Q9U6P4
B	483	GLY	-	expression tag	UNP Q9U6P4
B	484	SER	-	expression tag	UNP Q9U6P4
B	485	ALA	-	expression tag	UNP Q9U6P4
B	486	TRP	-	expression tag	UNP Q9U6P4
B	487	SER	-	expression tag	UNP Q9U6P4
B	488	HIS	-	expression tag	UNP Q9U6P4
B	489	PRO	-	expression tag	UNP Q9U6P4
B	490	GLN	-	expression tag	UNP Q9U6P4
B	491	PHE	-	expression tag	UNP Q9U6P4
B	492	GLU	-	expression tag	UNP Q9U6P4
B	493	LYS	-	expression tag	UNP Q9U6P4
B	494	GLY	-	expression tag	UNP Q9U6P4
B	495	GLY	-	expression tag	UNP Q9U6P4
B	496	GLY	-	expression tag	UNP Q9U6P4
B	497	SER	-	expression tag	UNP Q9U6P4
B	498	GLY	-	expression tag	UNP Q9U6P4
B	499	GLY	-	expression tag	UNP Q9U6P4
B	500	GLY	-	expression tag	UNP Q9U6P4
B	501	SER	-	expression tag	UNP Q9U6P4
B	502	GLY	-	expression tag	UNP Q9U6P4
B	503	GLY	-	expression tag	UNP Q9U6P4
B	504	SER	-	expression tag	UNP Q9U6P4
B	505	ALA	-	expression tag	UNP Q9U6P4
B	506	TRP	-	expression tag	UNP Q9U6P4
B	507	SER	-	expression tag	UNP Q9U6P4
B	508	HIS	-	expression tag	UNP Q9U6P4
B	509	PRO	-	expression tag	UNP Q9U6P4
B	510	GLN	-	expression tag	UNP Q9U6P4
B	511	PHE	-	expression tag	UNP Q9U6P4
B	512	GLU	-	expression tag	UNP Q9U6P4
B	513	LYS	-	expression tag	UNP Q9U6P4
C	466	THR	-	expression tag	UNP Q9U6P4
C	467	GLY	-	expression tag	UNP Q9U6P4
C	468	GLY	-	expression tag	UNP Q9U6P4
C	469	SER	-	expression tag	UNP Q9U6P4
C	470	SER	-	expression tag	UNP Q9U6P4
C	471	GLY	-	expression tag	UNP Q9U6P4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	472	LEU	-	expression tag	UNP Q9U6P4
C	473	GLU	-	expression tag	UNP Q9U6P4
C	474	VAL	-	expression tag	UNP Q9U6P4
C	475	LEU	-	expression tag	UNP Q9U6P4
C	476	PHE	-	expression tag	UNP Q9U6P4
C	477	GLN	-	expression tag	UNP Q9U6P4
C	478	GLY	-	expression tag	UNP Q9U6P4
C	479	PRO	-	expression tag	UNP Q9U6P4
C	480	GLY	-	expression tag	UNP Q9U6P4
C	481	SER	-	expression tag	UNP Q9U6P4
C	482	GLY	-	expression tag	UNP Q9U6P4
C	483	GLY	-	expression tag	UNP Q9U6P4
C	484	SER	-	expression tag	UNP Q9U6P4
C	485	ALA	-	expression tag	UNP Q9U6P4
C	486	TRP	-	expression tag	UNP Q9U6P4
C	487	SER	-	expression tag	UNP Q9U6P4
C	488	HIS	-	expression tag	UNP Q9U6P4
C	489	PRO	-	expression tag	UNP Q9U6P4
C	490	GLN	-	expression tag	UNP Q9U6P4
C	491	PHE	-	expression tag	UNP Q9U6P4
C	492	GLU	-	expression tag	UNP Q9U6P4
C	493	LYS	-	expression tag	UNP Q9U6P4
C	494	GLY	-	expression tag	UNP Q9U6P4
C	495	GLY	-	expression tag	UNP Q9U6P4
C	496	GLY	-	expression tag	UNP Q9U6P4
C	497	SER	-	expression tag	UNP Q9U6P4
C	498	GLY	-	expression tag	UNP Q9U6P4
C	499	GLY	-	expression tag	UNP Q9U6P4
C	500	GLY	-	expression tag	UNP Q9U6P4
C	501	SER	-	expression tag	UNP Q9U6P4
C	502	GLY	-	expression tag	UNP Q9U6P4
C	503	GLY	-	expression tag	UNP Q9U6P4
C	504	SER	-	expression tag	UNP Q9U6P4
C	505	ALA	-	expression tag	UNP Q9U6P4
C	506	TRP	-	expression tag	UNP Q9U6P4
C	507	SER	-	expression tag	UNP Q9U6P4
C	508	HIS	-	expression tag	UNP Q9U6P4
C	509	PRO	-	expression tag	UNP Q9U6P4
C	510	GLN	-	expression tag	UNP Q9U6P4
C	511	PHE	-	expression tag	UNP Q9U6P4
C	512	GLU	-	expression tag	UNP Q9U6P4
C	513	LYS	-	expression tag	UNP Q9U6P4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	466	THR	-	expression tag	UNP Q9U6P4
D	467	GLY	-	expression tag	UNP Q9U6P4
D	468	GLY	-	expression tag	UNP Q9U6P4
D	469	SER	-	expression tag	UNP Q9U6P4
D	470	SER	-	expression tag	UNP Q9U6P4
D	471	GLY	-	expression tag	UNP Q9U6P4
D	472	LEU	-	expression tag	UNP Q9U6P4
D	473	GLU	-	expression tag	UNP Q9U6P4
D	474	VAL	-	expression tag	UNP Q9U6P4
D	475	LEU	-	expression tag	UNP Q9U6P4
D	476	PHE	-	expression tag	UNP Q9U6P4
D	477	GLN	-	expression tag	UNP Q9U6P4
D	478	GLY	-	expression tag	UNP Q9U6P4
D	479	PRO	-	expression tag	UNP Q9U6P4
D	480	GLY	-	expression tag	UNP Q9U6P4
D	481	SER	-	expression tag	UNP Q9U6P4
D	482	GLY	-	expression tag	UNP Q9U6P4
D	483	GLY	-	expression tag	UNP Q9U6P4
D	484	SER	-	expression tag	UNP Q9U6P4
D	485	ALA	-	expression tag	UNP Q9U6P4
D	486	TRP	-	expression tag	UNP Q9U6P4
D	487	SER	-	expression tag	UNP Q9U6P4
D	488	HIS	-	expression tag	UNP Q9U6P4
D	489	PRO	-	expression tag	UNP Q9U6P4
D	490	GLN	-	expression tag	UNP Q9U6P4
D	491	PHE	-	expression tag	UNP Q9U6P4
D	492	GLU	-	expression tag	UNP Q9U6P4
D	493	LYS	-	expression tag	UNP Q9U6P4
D	494	GLY	-	expression tag	UNP Q9U6P4
D	495	GLY	-	expression tag	UNP Q9U6P4
D	496	GLY	-	expression tag	UNP Q9U6P4
D	497	SER	-	expression tag	UNP Q9U6P4
D	498	GLY	-	expression tag	UNP Q9U6P4
D	499	GLY	-	expression tag	UNP Q9U6P4
D	500	GLY	-	expression tag	UNP Q9U6P4
D	501	SER	-	expression tag	UNP Q9U6P4
D	502	GLY	-	expression tag	UNP Q9U6P4
D	503	GLY	-	expression tag	UNP Q9U6P4
D	504	SER	-	expression tag	UNP Q9U6P4
D	505	ALA	-	expression tag	UNP Q9U6P4
D	506	TRP	-	expression tag	UNP Q9U6P4
D	507	SER	-	expression tag	UNP Q9U6P4

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Chain	Residue	Modelled	Actual	Comment	Reference
D	508	HIS	-	expression tag	UNP Q9U6P4
D	509	PRO	-	expression tag	UNP Q9U6P4
D	510	GLN	-	expression tag	UNP Q9U6P4
D	511	PHE	-	expression tag	UNP Q9U6P4
D	512	GLU	-	expression tag	UNP Q9U6P4
D	513	LYS	-	expression tag	UNP Q9U6P4
E	466	THR	-	expression tag	UNP Q9U6P4
E	467	GLY	-	expression tag	UNP Q9U6P4
E	468	GLY	-	expression tag	UNP Q9U6P4
E	469	SER	-	expression tag	UNP Q9U6P4
E	470	SER	-	expression tag	UNP Q9U6P4
E	471	GLY	-	expression tag	UNP Q9U6P4
E	472	LEU	-	expression tag	UNP Q9U6P4
E	473	GLU	-	expression tag	UNP Q9U6P4
E	474	VAL	-	expression tag	UNP Q9U6P4
E	475	LEU	-	expression tag	UNP Q9U6P4
E	476	PHE	-	expression tag	UNP Q9U6P4
E	477	GLN	-	expression tag	UNP Q9U6P4
E	478	GLY	-	expression tag	UNP Q9U6P4
E	479	PRO	-	expression tag	UNP Q9U6P4
E	480	GLY	-	expression tag	UNP Q9U6P4
E	481	SER	-	expression tag	UNP Q9U6P4
E	482	GLY	-	expression tag	UNP Q9U6P4
E	483	GLY	-	expression tag	UNP Q9U6P4
E	484	SER	-	expression tag	UNP Q9U6P4
E	485	ALA	-	expression tag	UNP Q9U6P4
E	486	TRP	-	expression tag	UNP Q9U6P4
E	487	SER	-	expression tag	UNP Q9U6P4
E	488	HIS	-	expression tag	UNP Q9U6P4
E	489	PRO	-	expression tag	UNP Q9U6P4
E	490	GLN	-	expression tag	UNP Q9U6P4
E	491	PHE	-	expression tag	UNP Q9U6P4
E	492	GLU	-	expression tag	UNP Q9U6P4
E	493	LYS	-	expression tag	UNP Q9U6P4
E	494	GLY	-	expression tag	UNP Q9U6P4
E	495	GLY	-	expression tag	UNP Q9U6P4
E	496	GLY	-	expression tag	UNP Q9U6P4
E	497	SER	-	expression tag	UNP Q9U6P4
E	498	GLY	-	expression tag	UNP Q9U6P4
E	499	GLY	-	expression tag	UNP Q9U6P4
E	500	GLY	-	expression tag	UNP Q9U6P4
E	501	SER	-	expression tag	UNP Q9U6P4

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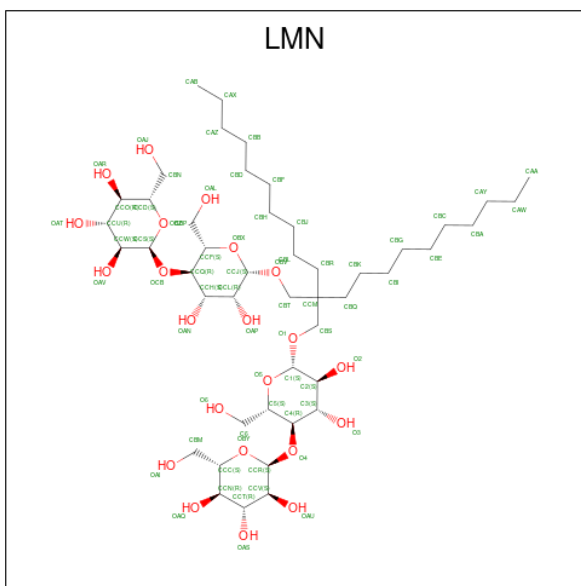
Chain	Residue	Modelled	Actual	Comment	Reference
E	502	GLY	-	expression tag	UNP Q9U6P4
E	503	GLY	-	expression tag	UNP Q9U6P4
E	504	SER	-	expression tag	UNP Q9U6P4
E	505	ALA	-	expression tag	UNP Q9U6P4
E	506	TRP	-	expression tag	UNP Q9U6P4
E	507	SER	-	expression tag	UNP Q9U6P4
E	508	HIS	-	expression tag	UNP Q9U6P4
E	509	PRO	-	expression tag	UNP Q9U6P4
E	510	GLN	-	expression tag	UNP Q9U6P4
E	511	PHE	-	expression tag	UNP Q9U6P4
E	512	GLU	-	expression tag	UNP Q9U6P4
E	513	LYS	-	expression tag	UNP Q9U6P4
F	466	THR	-	expression tag	UNP Q9U6P4
F	467	GLY	-	expression tag	UNP Q9U6P4
F	468	GLY	-	expression tag	UNP Q9U6P4
F	469	SER	-	expression tag	UNP Q9U6P4
F	470	SER	-	expression tag	UNP Q9U6P4
F	471	GLY	-	expression tag	UNP Q9U6P4
F	472	LEU	-	expression tag	UNP Q9U6P4
F	473	GLU	-	expression tag	UNP Q9U6P4
F	474	VAL	-	expression tag	UNP Q9U6P4
F	475	LEU	-	expression tag	UNP Q9U6P4
F	476	PHE	-	expression tag	UNP Q9U6P4
F	477	GLN	-	expression tag	UNP Q9U6P4
F	478	GLY	-	expression tag	UNP Q9U6P4
F	479	PRO	-	expression tag	UNP Q9U6P4
F	480	GLY	-	expression tag	UNP Q9U6P4
F	481	SER	-	expression tag	UNP Q9U6P4
F	482	GLY	-	expression tag	UNP Q9U6P4
F	483	GLY	-	expression tag	UNP Q9U6P4
F	484	SER	-	expression tag	UNP Q9U6P4
F	485	ALA	-	expression tag	UNP Q9U6P4
F	486	TRP	-	expression tag	UNP Q9U6P4
F	487	SER	-	expression tag	UNP Q9U6P4
F	488	HIS	-	expression tag	UNP Q9U6P4
F	489	PRO	-	expression tag	UNP Q9U6P4
F	490	GLN	-	expression tag	UNP Q9U6P4
F	491	PHE	-	expression tag	UNP Q9U6P4
F	492	GLU	-	expression tag	UNP Q9U6P4
F	493	LYS	-	expression tag	UNP Q9U6P4
F	494	GLY	-	expression tag	UNP Q9U6P4
F	495	GLY	-	expression tag	UNP Q9U6P4

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Chain	Residue	Modelled	Actual	Comment	Reference
F	496	GLY	-	expression tag	UNP Q9U6P4
F	497	SER	-	expression tag	UNP Q9U6P4
F	498	GLY	-	expression tag	UNP Q9U6P4
F	499	GLY	-	expression tag	UNP Q9U6P4
F	500	GLY	-	expression tag	UNP Q9U6P4
F	501	SER	-	expression tag	UNP Q9U6P4
F	502	GLY	-	expression tag	UNP Q9U6P4
F	503	GLY	-	expression tag	UNP Q9U6P4
F	504	SER	-	expression tag	UNP Q9U6P4
F	505	ALA	-	expression tag	UNP Q9U6P4
F	506	TRP	-	expression tag	UNP Q9U6P4
F	507	SER	-	expression tag	UNP Q9U6P4
F	508	HIS	-	expression tag	UNP Q9U6P4
F	509	PRO	-	expression tag	UNP Q9U6P4
F	510	GLN	-	expression tag	UNP Q9U6P4
F	511	PHE	-	expression tag	UNP Q9U6P4
F	512	GLU	-	expression tag	UNP Q9U6P4
F	513	LYS	-	expression tag	UNP Q9U6P4

- Molecule 2 is Lauryl Maltose Neopentyl Glycol (three-letter code: LMN) (formula:  $C_{47}H_{88}O_{22}$ ).



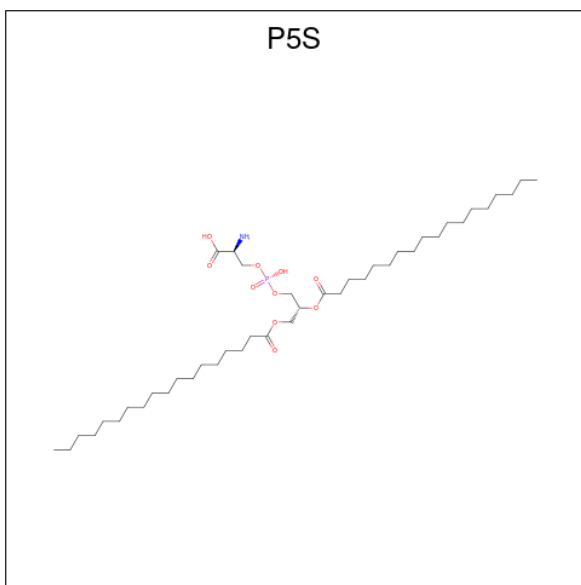
Mol	Chain	Residues	Atoms			AltConf
2	A	1	Total	C	O	0
			116	82	34	
2	A	1	Total	C	O	0
			116	82	34	

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Mol	Chain	Residues	Atoms			AltConf
			Total	C	O	
2	B	1	116	82	34	0
2	B	1	116	82	34	0
2	C	1	116	82	34	0
2	C	1	116	82	34	0
2	D	1	116	82	34	0
2	D	1	116	82	34	0
2	E	1	116	82	34	0
2	E	1	116	82	34	0
2	F	1	116	82	34	0
2	F	1	116	82	34	0

- Molecule 3 is O-[(R)-{[(2R)-2,3-bis(octadecanoyloxy)propyl]oxy}(hydroxy)phosphoryl]-L-serine (three-letter code: P5S) (formula:  $C_{42}H_{82}NO_{10}P$ ).



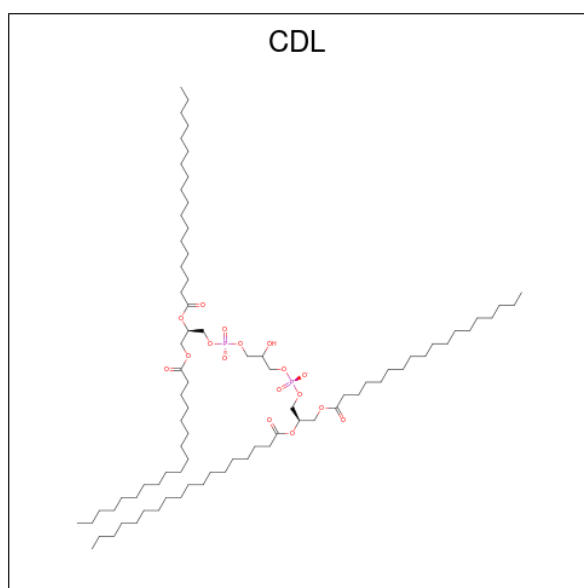
Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
3	A	1	41	29	1	10	1	0

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Mol	Chain	Residues	Atoms					AltConf
			Total	C	N	O	P	
3	B	1	Total 41	C 29	N 1	O 10	P 1	0
3	C	1	Total 41	C 29	N 1	O 10	P 1	0
3	D	1	Total 41	C 29	N 1	O 10	P 1	0
3	E	1	Total 41	C 29	N 1	O 10	P 1	0
3	F	1	Total 41	C 29	N 1	O 10	P 1	0

- Molecule 4 is CARDIOLIPIN (three-letter code: CDL) (formula:  $C_{81}H_{156}O_{17}P_2$ ).



Mol	Chain	Residues	Atoms				AltConf
			Total	C	O	P	
4	A	1	Total 92	C 73	O 17	P 2	0
4	B	1	Total 92	C 73	O 17	P 2	0
4	C	1	Total 92	C 73	O 17	P 2	0
4	D	1	Total 92	C 73	O 17	P 2	0
4	E	1	Total 92	C 73	O 17	P 2	0
4	F	1	Total 92	C 73	O 17	P 2	0









## 4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C6	Depositor
Number of particles used	159252	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	NONE	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ( $e^-/\text{\AA}^2$ )	50	Depositor
Minimum defocus (nm)	Not provided	
Maximum defocus (nm)	Not provided	
Magnification	Not provided	
Image detector	GATAN K2 QUANTUM (4k x 4k)	Depositor
Maximum map value	6.165	Depositor
Minimum map value	-3.928	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.136	Depositor
Recommended contour level	0.7	Depositor
Map size ( $\text{\AA}$ )	386.4, 386.4, 386.4	wwPDB
Map dimensions	280, 280, 280	wwPDB
Map angles ( $^\circ$ )	90.0, 90.0, 90.0	wwPDB
Pixel spacing ( $\text{\AA}$ )	1.38, 1.38, 1.38	Depositor

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: P5S, LMN, CDL

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	0.30	0/2909	0.46	1/3964 (0.0%)
1	B	0.30	0/2909	0.46	1/3964 (0.0%)
1	C	0.30	0/2909	0.46	1/3964 (0.0%)
1	D	0.30	0/2909	0.46	1/3964 (0.0%)
1	E	0.30	0/2909	0.46	1/3964 (0.0%)
1	F	0.30	0/2909	0.46	1/3964 (0.0%)
All	All	0.30	0/17454	0.46	6/23784 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	C	57	LEU	CA-CB-CG	5.22	127.31	115.30
1	B	57	LEU	CA-CB-CG	5.21	127.29	115.30
1	A	57	LEU	CA-CB-CG	5.21	127.28	115.30
1	D	57	LEU	CA-CB-CG	5.20	127.26	115.30
1	F	57	LEU	CA-CB-CG	5.20	127.25	115.30
1	E	57	LEU	CA-CB-CG	5.20	127.25	115.30

There are no chirality outliers.

There are no planarity outliers.

### 5.2 Too-close contacts

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	2834	0	2845	42	0
1	B	2834	0	2845	44	0
1	C	2834	0	2845	42	0
1	D	2834	0	2845	43	0
1	E	2834	0	2845	47	0
1	F	2834	0	2845	43	0
2	A	116	0	155	21	0
2	B	116	0	155	20	0
2	C	116	0	155	18	0
2	D	116	0	155	20	0
2	E	116	0	155	21	0
2	F	116	0	155	20	0
3	A	41	0	48	1	0
3	B	41	0	48	1	0
3	C	41	0	48	1	0
3	D	41	0	48	1	0
3	E	41	0	48	1	0
3	F	41	0	48	1	0
4	A	92	0	137	3	0
4	B	92	0	137	3	0
4	C	92	0	137	4	0
4	D	92	0	137	4	0
4	E	92	0	137	4	0
4	F	92	0	137	3	0
All	All	18498	0	19110	303	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 8.

All (303) 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:B:1002:LMN:O3	2:B:1002:LMN:OAU	1.59	1.19
2:A:604:LMN:OAU	2:A:604:LMN:O3	1.59	1.15
1:E:156:PHE:HZ	2:E:1002:LMN:HBJ	1.28	0.98
1:A:156:PHE:HZ	2:A:604:LMN:HBJ	1.28	0.97
1:F:156:PHE:HZ	2:F:1002:LMN:HBJ	1.28	0.97
1:B:144:PHE:O	1:B:148:PHE:HD2	1.48	0.96
1:B:156:PHE:HZ	2:B:1002:LMN:HBJ	1.30	0.95
1:C:144:PHE:O	1:C:148:PHE:HD2	1.48	0.95
1:E:144:PHE:O	1:E:148:PHE:HD2	1.48	0.95
1:A:144:PHE:O	1:A:148:PHE:HD2	1.48	0.95

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:144:PHE:O	1:F:148:PHE:HD2	1.48	0.95
1:D:144:PHE:O	1:D:148:PHE:HD2	1.48	0.94
1:D:156:PHE:HZ	2:D:1002:LMN:HBJ	1.30	0.93
1:C:156:PHE:HZ	2:C:1002:LMN:HBJ	1.34	0.90
1:D:156:PHE:CZ	2:D:1002:LMN:HBJ	2.09	0.88
1:F:156:PHE:CZ	2:F:1002:LMN:HBJ	2.09	0.88
1:E:156:PHE:CZ	2:E:1002:LMN:HBJ	2.09	0.87
1:A:156:PHE:CZ	2:A:604:LMN:HBJ	2.09	0.87
1:B:156:PHE:CZ	2:B:1002:LMN:HBJ	2.12	0.85
1:C:156:PHE:CZ	2:C:1002:LMN:HBJ	2.14	0.82
1:D:144:PHE:O	1:D:148:PHE:CD2	2.34	0.81
1:F:144:PHE:O	1:F:148:PHE:CD2	2.34	0.81
1:B:144:PHE:O	1:B:148:PHE:CD2	2.34	0.81
1:C:144:PHE:O	1:C:148:PHE:CD2	2.34	0.80
1:E:144:PHE:O	1:E:148:PHE:CD2	2.34	0.79
1:A:144:PHE:O	1:A:148:PHE:CD2	2.34	0.78
1:B:68:MET:CE	2:B:1002:LMN:HBK	2.14	0.77
1:A:68:MET:CE	2:A:604:LMN:HBK	2.16	0.76
1:D:68:MET:CE	2:D:1002:LMN:HBK	2.16	0.75
1:E:68:MET:CE	2:E:1002:LMN:HBK	2.16	0.75
1:C:68:MET:CE	2:C:1002:LMN:HBK	2.18	0.74
1:F:68:MET:CE	2:F:1002:LMN:HBK	2.19	0.73
1:E:72:ILE:CG2	2:E:1002:LMN:HBNA	2.19	0.73
1:D:68:MET:HE3	2:D:1002:LMN:HBK	1.72	0.71
1:A:72:ILE:CG2	2:A:604:LMN:HBNA	2.21	0.71
1:A:131:PRO:HB3	1:F:33:THR:HA	1.73	0.70
1:B:72:ILE:CG2	2:B:1002:LMN:HBNA	2.21	0.70
1:F:72:ILE:CG2	2:F:1002:LMN:HBNA	2.22	0.69
1:A:68:MET:HE1	2:A:604:LMN:HBK	1.74	0.69
1:A:61:LEU:HD23	1:B:72:ILE:HB	1.74	0.69
1:C:65:LEU:HD23	2:D:1002:LMN:H3	1.75	0.68
1:C:33:THR:HA	1:D:131:PRO:HB3	1.74	0.68
1:D:33:THR:HA	1:E:131:PRO:HB3	1.75	0.68
1:A:33:THR:HA	1:B:131:PRO:HB3	1.75	0.68
1:D:65:LEU:HD23	2:E:1002:LMN:H3	1.75	0.68
4:F:1001:CDL:H472	4:F:1001:CDL:H362	1.76	0.68
1:D:72:ILE:CG2	2:D:1002:LMN:HBNA	2.24	0.68
4:E:1001:CDL:H362	4:E:1001:CDL:H472	1.76	0.68
1:E:33:THR:HA	1:F:131:PRO:HB3	1.75	0.67
1:D:61:LEU:HD23	1:E:72:ILE:HB	1.77	0.67
4:A:603:CDL:H362	4:A:603:CDL:H472	1.76	0.67

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:61:LEU:HD23	1:D:72:ILE:HB	1.77	0.67
1:B:65:LEU:HD23	2:C:1002:LMN:H3	1.75	0.67
1:E:72:ILE:HG23	2:E:1002:LMN:HCO	1.77	0.67
1:A:65:LEU:HD23	2:B:1002:LMN:H3	1.75	0.67
1:C:68:MET:HE3	2:C:1002:LMN:HBK	1.77	0.67
1:B:33:THR:HA	1:C:131:PRO:HB3	1.75	0.66
1:B:68:MET:HE3	2:B:1002:LMN:HBK	1.76	0.66
1:B:61:LEU:HD23	1:C:72:ILE:HB	1.76	0.66
4:D:1001:CDL:H472	4:D:1001:CDL:H362	1.76	0.66
4:B:1001:CDL:H362	4:B:1001:CDL:H472	1.76	0.66
1:A:72:ILE:HB	1:F:61:LEU:HD23	1.76	0.66
2:A:604:LMN:H3	1:F:65:LEU:HD23	1.76	0.66
2:E:1003:LMN:HBHA	2:E:1003:LMN:HBKA	1.78	0.66
2:C:1003:LMN:HBHA	2:C:1003:LMN:HBKA	1.78	0.66
1:F:72:ILE:HG23	2:F:1002:LMN:HCO	1.76	0.66
4:C:1001:CDL:H472	4:C:1001:CDL:H362	1.76	0.66
1:E:65:LEU:HD23	2:F:1002:LMN:H3	1.75	0.66
1:D:72:ILE:HG23	2:D:1002:LMN:HCO	1.78	0.66
1:A:72:ILE:HG23	2:A:604:LMN:HCO	1.78	0.65
1:B:68:MET:HE1	2:B:1002:LMN:HBK	1.78	0.65
2:D:1003:LMN:HBHA	2:D:1003:LMN:HBKA	1.78	0.65
2:F:1002:LMN:H1	2:F:1002:LMN:OAP	1.97	0.65
2:B:1002:LMN:H1	2:B:1002:LMN:OAP	1.97	0.65
1:C:72:ILE:CG2	2:C:1002:LMN:HBNA	2.27	0.65
2:F:1003:LMN:HBHA	2:F:1003:LMN:HBKA	1.78	0.65
2:A:601:LMN:HBKA	2:A:601:LMN:HBHA	1.78	0.65
2:E:1002:LMN:H1	2:E:1002:LMN:OAP	1.97	0.64
2:B:1003:LMN:HBKA	2:B:1003:LMN:HBHA	1.78	0.64
2:D:1002:LMN:H1	2:D:1002:LMN:OAP	1.97	0.64
1:E:68:MET:HE3	2:E:1002:LMN:HBK	1.78	0.64
2:A:604:LMN:H1	2:A:604:LMN:OAP	1.97	0.64
2:C:1002:LMN:H1	2:C:1002:LMN:OAP	1.97	0.64
1:E:61:LEU:HD23	1:F:72:ILE:HB	1.79	0.64
1:F:68:MET:HE1	2:F:1002:LMN:HBK	1.79	0.64
1:B:72:ILE:HG23	2:B:1002:LMN:HCO	1.80	0.63
1:E:68:MET:HE1	2:E:1002:LMN:HBK	1.79	0.63
1:C:72:ILE:HG23	2:C:1002:LMN:HCO	1.82	0.62
1:C:57:LEU:HB2	1:C:326:VAL:HG21	1.82	0.61
1:D:57:LEU:HB2	1:D:326:VAL:HG21	1.82	0.61
1:A:57:LEU:HB2	1:A:326:VAL:HG21	1.82	0.61
1:B:57:LEU:HB2	1:B:326:VAL:HG21	1.82	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:57:LEU:HB2	1:F:326:VAL:HG21	1.82	0.60
2:E:1002:LMN:H6A	2:E:1002:LMN:HBP	1.84	0.60
1:F:152:ILE:HD11	2:F:1002:LMN:HAWA	1.84	0.60
1:E:57:LEU:HB2	1:E:326:VAL:HG21	1.82	0.60
1:A:68:MET:HE3	2:A:604:LMN:HBK	1.84	0.60
2:A:604:LMN:H6A	2:A:604:LMN:HBP	1.84	0.60
1:C:68:MET:HE1	2:C:1002:LMN:HBK	1.83	0.60
2:B:1002:LMN:H6A	2:B:1002:LMN:HBP	1.84	0.59
2:D:1002:LMN:HBP	2:D:1002:LMN:H6A	1.84	0.59
2:F:1002:LMN:H6A	2:F:1002:LMN:HBP	1.84	0.59
1:E:152:ILE:HD11	2:E:1002:LMN:HAWA	1.86	0.58
2:C:1002:LMN:HBP	2:C:1002:LMN:H6A	1.84	0.58
1:D:152:ILE:HD11	2:D:1002:LMN:HAWA	1.86	0.58
1:F:68:MET:HE3	2:F:1002:LMN:HBK	1.84	0.58
1:A:152:ILE:HD11	2:A:604:LMN:HAWA	1.85	0.57
1:C:152:ILE:HD11	2:C:1002:LMN:HAWA	1.86	0.57
1:D:68:MET:HE1	2:D:1002:LMN:HBK	1.86	0.56
1:F:152:ILE:CD1	2:F:1002:LMN:HAWA	2.36	0.56
1:D:152:ILE:CD1	2:D:1002:LMN:HAWA	2.36	0.55
1:B:255:ILE:HD11	1:B:449:TYR:CD2	2.42	0.55
1:C:255:ILE:HD11	1:C:449:TYR:CD2	2.42	0.55
1:E:72:ILE:HG21	2:E:1002:LMN:HBNA	1.88	0.55
1:A:255:ILE:HD11	1:A:449:TYR:CD2	2.42	0.55
1:F:36:SER:HA	1:F:39:ARG:HE	1.72	0.55
1:A:36:SER:HA	1:A:39:ARG:HE	1.72	0.55
1:E:36:SER:HA	1:E:39:ARG:HE	1.72	0.55
1:B:36:SER:HA	1:B:39:ARG:HE	1.72	0.55
1:D:255:ILE:HD11	1:D:449:TYR:CD2	2.42	0.55
1:F:255:ILE:HD11	1:F:449:TYR:CD2	2.42	0.54
1:E:255:ILE:HD11	1:E:449:TYR:CD2	2.42	0.54
1:B:152:ILE:HD11	2:B:1002:LMN:HAWA	1.88	0.54
1:A:152:ILE:CD1	2:A:604:LMN:HAWA	2.37	0.54
1:C:152:ILE:CD1	2:C:1002:LMN:HAWA	2.38	0.54
1:B:72:ILE:HG21	2:B:1002:LMN:HBNA	1.89	0.54
1:D:36:SER:HA	1:D:39:ARG:HE	1.72	0.54
1:C:36:SER:HA	1:C:39:ARG:HE	1.72	0.54
1:E:152:ILE:CD1	2:E:1002:LMN:HAWA	2.38	0.54
1:F:72:ILE:HG21	2:F:1002:LMN:HBNA	1.90	0.53
1:B:260:ILE:HD12	1:B:273:LEU:HD21	1.92	0.52
1:E:260:ILE:HD12	1:E:273:LEU:HD21	1.92	0.52
1:A:260:ILE:HD12	1:A:273:LEU:HD21	1.92	0.52

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:260:ILE:HD12	1:C:273:LEU:HD21	1.92	0.52
1:D:260:ILE:HD12	1:D:273:LEU:HD21	1.92	0.52
1:F:260:ILE:HD12	1:F:273:LEU:HD21	1.92	0.52
1:B:141:PRO:HG3	2:B:1003:LMN:HCJ	1.92	0.52
1:A:72:ILE:HG21	2:A:604:LMN:HBNA	1.90	0.51
1:C:141:PRO:HG3	2:C:1003:LMN:HCJ	1.92	0.51
1:A:141:PRO:HG3	2:A:601:LMN:HCJ	1.92	0.51
1:C:200:ALA:HB1	1:C:206:TYR:HB2	1.93	0.51
1:A:200:ALA:HB1	1:A:206:TYR:HB2	1.93	0.50
1:E:141:PRO:HG3	2:E:1003:LMN:HCJ	1.92	0.50
1:F:141:PRO:HG3	2:F:1003:LMN:HCJ	1.92	0.50
1:D:200:ALA:HB1	1:D:206:TYR:HB2	1.93	0.50
1:B:152:ILE:CD1	2:B:1002:LMN:HAWA	2.41	0.50
1:F:200:ALA:HB1	1:F:206:TYR:HB2	1.93	0.50
1:C:72:ILE:HG21	2:C:1002:LMN:HBNA	1.94	0.50
1:D:141:PRO:HG3	2:D:1003:LMN:HCJ	1.92	0.50
1:B:200:ALA:HB1	1:B:206:TYR:HB2	1.93	0.49
1:D:152:ILE:O	1:D:155:ILE:HG22	2.13	0.49
1:D:72:ILE:HG21	2:D:1002:LMN:HBNA	1.95	0.49
1:E:152:ILE:O	1:E:155:ILE:HG22	2.13	0.49
1:A:152:ILE:O	1:A:155:ILE:HG22	2.13	0.49
1:C:152:ILE:O	1:C:155:ILE:HG22	2.13	0.49
1:E:200:ALA:HB1	1:E:206:TYR:HB2	1.93	0.49
1:F:152:ILE:O	1:F:155:ILE:HG22	2.13	0.48
1:A:222:LEU:O	1:A:225:ILE:HG22	2.13	0.48
1:F:222:LEU:O	1:F:225:ILE:HG22	2.13	0.48
1:E:72:ILE:HD13	2:E:1002:LMN:HCH	1.96	0.48
1:B:152:ILE:O	1:B:155:ILE:HG22	2.13	0.47
1:C:61:LEU:CD2	1:D:72:ILE:HB	2.44	0.47
1:A:61:LEU:CD2	1:B:72:ILE:HB	2.43	0.47
1:B:222:LEU:O	1:B:225:ILE:HG22	2.13	0.47
2:B:1003:LMN:HCH	2:B:1003:LMN:HCS	1.71	0.47
1:D:222:LEU:O	1:D:225:ILE:HG22	2.13	0.47
1:E:222:LEU:O	1:E:225:ILE:HG22	2.13	0.47
1:C:222:LEU:O	1:C:225:ILE:HG22	2.13	0.47
1:A:72:ILE:HD13	2:A:604:LMN:HCH	1.97	0.47
1:E:204:ARG:NH2	1:E:208:TYR:OH	2.42	0.47
1:B:61:LEU:CD2	1:C:72:ILE:HB	2.44	0.46
1:A:97:TYR:CD2	1:A:425:ILE:HD12	2.51	0.46
1:D:97:TYR:CD2	1:D:425:ILE:HD12	2.51	0.46
1:F:97:TYR:CD2	1:F:425:ILE:HD12	2.51	0.46

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:C:97:TYR:CD2	1:C:425:ILE:HD12	2.51	0.46
1:C:328:PHE:HB3	4:D:1001:CDL:OA7	2.16	0.46
1:A:72:ILE:HB	1:F:61:LEU:CD2	2.45	0.46
1:D:328:PHE:HB3	4:E:1001:CDL:OA7	2.16	0.46
1:A:328:PHE:HB3	4:B:1001:CDL:OA7	2.16	0.46
1:D:72:ILE:CD1	2:D:1002:LMN:HCJ	2.46	0.46
1:E:97:TYR:CD2	1:E:425:ILE:HD12	2.51	0.46
1:F:72:ILE:HD13	2:F:1002:LMN:HCH	1.98	0.46
1:B:97:TYR:CD2	1:B:425:ILE:HD12	2.51	0.46
1:D:72:ILE:HD13	2:D:1002:LMN:HCH	1.98	0.45
1:D:204:ARG:NH2	1:D:208:TYR:OH	2.42	0.45
4:A:603:CDL:OA7	1:F:328:PHE:HB3	2.16	0.45
1:B:259:SER:O	1:B:271:SER:OG	2.31	0.45
1:B:328:PHE:HB3	4:C:1001:CDL:OA7	2.16	0.45
1:E:328:PHE:HB3	4:F:1001:CDL:OA7	2.16	0.45
1:D:49:GLY:HA3	1:D:111:MET:SD	2.57	0.45
1:E:49:GLY:HA3	1:E:111:MET:SD	2.57	0.45
1:B:36:SER:HB2	1:B:39:ARG:HH21	1.82	0.45
1:B:72:ILE:HD13	2:B:1002:LMN:HCH	1.98	0.45
1:C:49:GLY:HA3	1:C:111:MET:SD	2.57	0.45
1:D:36:SER:HB2	1:D:39:ARG:HH21	1.82	0.45
1:E:35:ALA:O	1:E:38:SER:OG	2.27	0.45
1:C:36:SER:HB2	1:C:39:ARG:HH21	1.82	0.45
1:A:49:GLY:HA3	1:A:111:MET:SD	2.57	0.44
1:C:190:HIS:HA	1:C:275:GLN:HE21	1.82	0.44
1:D:190:HIS:HA	1:D:275:GLN:HE21	1.82	0.44
1:E:36:SER:HB2	1:E:39:ARG:HH21	1.82	0.44
1:B:190:HIS:HA	1:B:275:GLN:HE21	1.83	0.44
1:B:35:ALA:O	1:B:38:SER:OG	2.27	0.44
1:B:49:GLY:HA3	1:B:111:MET:SD	2.57	0.44
1:A:36:SER:HB2	1:A:39:ARG:HH21	1.82	0.44
1:F:49:GLY:HA3	1:F:111:MET:SD	2.57	0.44
1:C:261:LEU:HD11	1:C:264:VAL:HG23	2.00	0.44
1:E:61:LEU:CD2	1:F:72:ILE:HB	2.48	0.44
2:F:1002:LMN:HCL	2:F:1002:LMN:HBT	1.82	0.44
1:A:190:HIS:HA	1:A:275:GLN:HE21	1.82	0.44
1:E:190:HIS:HA	1:E:275:GLN:HE21	1.82	0.44
1:B:261:LEU:HD11	1:B:264:VAL:HG23	2.00	0.43
1:C:265:GLN:HG2	1:C:271:SER:HB3	2.00	0.43
1:F:261:LEU:HD11	1:F:264:VAL:HG23	2.00	0.43
1:D:265:GLN:HG2	1:D:271:SER:HB3	2.00	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:190:HIS:HA	1:F:275:GLN:HE21	1.82	0.43
1:F:265:GLN:HG2	1:F:271:SER:HB3	2.00	0.43
1:B:265:GLN:HG2	1:B:271:SER:HB3	2.00	0.43
1:D:35:ALA:O	1:D:38:SER:OG	2.27	0.43
1:F:36:SER:HB2	1:F:39:ARG:HH21	1.82	0.43
1:F:204:ARG:NH2	1:F:208:TYR:OH	2.42	0.43
1:A:72:ILE:CD1	2:A:604:LMN:HCJ	2.49	0.43
2:E:1002:LMN:HBI	2:E:1002:LMN:HBR	2.01	0.43
2:F:1002:LMN:HBI	2:F:1002:LMN:HBR	2.01	0.43
2:D:1002:LMN:HBR	2:D:1002:LMN:HBI	2.01	0.43
2:D:1003:LMN:HCS	2:D:1003:LMN:HCH	1.71	0.43
4:C:1001:CDL:H251	4:C:1001:CDL:H222	1.91	0.43
2:C:1002:LMN:HBI	2:C:1002:LMN:HBR	2.01	0.43
1:D:261:LEU:HD11	1:D:264:VAL:HG23	2.00	0.43
1:A:261:LEU:HD11	1:A:264:VAL:HG23	2.00	0.43
1:D:91:LEU:HG	1:D:422:HIS:CE1	2.54	0.43
1:B:216:LEU:HA	1:B:219:ILE:HG12	2.02	0.42
2:D:1002:LMN:HCL	2:D:1002:LMN:HBT	1.82	0.42
2:A:604:LMN:HBR	2:A:604:LMN:HBI	2.01	0.42
2:B:1002:LMN:HBR	2:B:1002:LMN:HBI	2.01	0.42
1:D:216:LEU:HA	1:D:219:ILE:HG12	2.01	0.42
1:E:72:ILE:CD1	2:E:1002:LMN:HCJ	2.49	0.42
2:E:1003:LMN:HCH	2:E:1003:LMN:HCS	1.70	0.42
2:A:601:LMN:HBL	2:A:601:LMN:HBSA	1.84	0.42
1:F:411:VAL:HA	1:F:414:THR:HG22	2.01	0.42
1:A:216:LEU:HA	1:A:219:ILE:HG12	2.01	0.42
1:B:411:VAL:HA	1:B:414:THR:HG22	2.01	0.42
1:C:411:VAL:HA	1:C:414:THR:HG22	2.01	0.42
1:E:91:LEU:HG	1:E:422:HIS:CE1	2.54	0.42
4:E:1001:CDL:H251	4:E:1001:CDL:H222	1.91	0.42
1:F:72:ILE:CD1	2:F:1002:LMN:HCJ	2.50	0.42
1:A:91:LEU:HG	1:A:422:HIS:CE1	2.54	0.42
1:A:265:GLN:HG2	1:A:271:SER:HB3	2.00	0.42
1:E:216:LEU:HA	1:E:219:ILE:HG12	2.01	0.42
1:C:91:LEU:HG	1:C:422:HIS:CE1	2.54	0.42
1:E:261:LEU:HD11	1:E:264:VAL:HG23	2.00	0.42
1:E:265:GLN:HG2	1:E:271:SER:HB3	2.00	0.42
1:F:91:LEU:HG	1:F:422:HIS:CE1	2.54	0.42
1:A:411:VAL:HA	1:A:414:THR:HG22	2.01	0.42
1:A:204:ARG:NH2	1:A:208:TYR:OH	2.42	0.42
2:A:601:LMN:HCH	2:A:601:LMN:HCS	1.71	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:259:SER:O	1:E:271:SER:OG	2.31	0.42
2:F:1003:LMN:HCH	2:F:1003:LMN:HCS	1.71	0.42
4:A:603:CDL:OA9	4:A:603:CDL:O1	2.38	0.41
4:B:1001:CDL:O1	4:B:1001:CDL:OA9	2.38	0.41
1:D:411:VAL:HA	1:D:414:THR:HG22	2.01	0.41
2:E:1002:LMN:HCL	2:E:1002:LMN:HBT	1.82	0.41
3:E:1004:P5S:H49A	3:E:1004:P5S:H27A	2.03	0.41
1:B:91:LEU:HG	1:B:422:HIS:CE1	2.54	0.41
4:E:1001:CDL:OA9	4:E:1001:CDL:O1	2.38	0.41
1:B:204:ARG:NH2	1:B:208:TYR:OH	2.42	0.41
2:B:1003:LMN:HBG	2:B:1003:LMN:HBAA	1.88	0.41
3:B:1004:P5S:H49A	3:B:1004:P5S:H27A	2.02	0.41
1:C:204:ARG:NH2	1:C:208:TYR:OH	2.42	0.41
1:C:259:SER:O	1:C:271:SER:OG	2.31	0.41
1:F:255:ILE:HD11	1:F:449:TYR:HD2	1.86	0.41
1:A:255:ILE:HD11	1:A:449:TYR:HD2	1.85	0.41
2:C:1003:LMN:HBSA	2:C:1003:LMN:HBL	1.84	0.41
1:F:216:LEU:HA	1:F:219:ILE:HG12	2.01	0.41
4:D:1001:CDL:H251	4:D:1001:CDL:H222	1.91	0.41
1:E:404:PHE:HD1	1:E:404:PHE:HA	1.71	0.41
1:E:411:VAL:HA	1:E:414:THR:HG22	2.01	0.41
1:C:56:ALA:O	1:C:57:LEU:HG	2.21	0.41
1:C:216:LEU:HA	1:C:219:ILE:HG12	2.01	0.41
4:C:1001:CDL:OA9	4:C:1001:CDL:O1	2.38	0.41
1:D:61:LEU:CD2	1:E:72:ILE:HB	2.46	0.41
4:D:1001:CDL:O1	4:D:1001:CDL:OA9	2.38	0.41
1:A:268:LEU:HA	1:A:269:PRO:HD3	1.93	0.41
3:D:1004:P5S:H27A	3:D:1004:P5S:H49A	2.03	0.41
3:F:1004:P5S:H49A	3:F:1004:P5S:H27A	2.03	0.41
1:D:125:VAL:HG21	1:D:397:TRP:CZ3	2.56	0.41
1:E:56:ALA:O	1:E:57:LEU:HG	2.21	0.41
2:E:1003:LMN:HBSA	2:E:1003:LMN:HBL	1.84	0.41
2:C:1003:LMN:HBG	2:C:1003:LMN:HBAA	1.88	0.41
1:E:125:VAL:HG21	1:E:397:TRP:CZ3	2.56	0.41
3:C:1004:P5S:H49A	3:C:1004:P5S:H27A	2.03	0.40
1:E:255:ILE:HD11	1:E:449:TYR:HD2	1.86	0.40
3:A:602:P5S:H49A	3:A:602:P5S:H27A	2.03	0.40
1:B:56:ALA:O	1:B:57:LEU:HG	2.21	0.40
1:C:45:ILE:HG12	1:C:144:PHE:CE1	2.57	0.40
1:D:56:ALA:O	1:D:57:LEU:HG	2.21	0.40
1:F:125:VAL:HG21	1:F:397:TRP:CZ3	2.56	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
4:F:1001:CDL:OA9	4:F:1001:CDL:O1	2.38	0.40
1:B:255:ILE:HD11	1:B:449:TYR:HD2	1.86	0.40
1:B:125:VAL:HG21	1:B:397:TRP:CZ3	2.56	0.40
1:C:125:VAL:HG21	1:C:397:TRP:CZ3	2.56	0.40
1:E:45:ILE:HG12	1:E:144:PHE:CE1	2.57	0.40
1:F:45:ILE:HG12	1:F:144:PHE:CE1	2.57	0.40

There are no symmetry-related clashes.

## 5.3 Torsion angles [i](#)

### 5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	357/513 (70%)	340 (95%)	17 (5%)	0	100	100
1	B	357/513 (70%)	340 (95%)	17 (5%)	0	100	100
1	C	357/513 (70%)	341 (96%)	16 (4%)	0	100	100
1	D	357/513 (70%)	340 (95%)	17 (5%)	0	100	100
1	E	357/513 (70%)	340 (95%)	17 (5%)	0	100	100
1	F	357/513 (70%)	340 (95%)	17 (5%)	0	100	100
All	All	2142/3078 (70%)	2041 (95%)	101 (5%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	312/416 (75%)	312 (100%)	0	100	100
1	B	312/416 (75%)	312 (100%)	0	100	100
1	C	312/416 (75%)	312 (100%)	0	100	100
1	D	312/416 (75%)	312 (100%)	0	100	100
1	E	312/416 (75%)	312 (100%)	0	100	100
1	F	312/416 (75%)	312 (100%)	0	100	100
All	All	1872/2496 (75%)	1872 (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. 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 monosaccharides in this entry.

## 5.6 Ligand geometry [i](#)

24 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	$\# Z  > 2$
4	CDL	A	603	-	91,91,99	0.30	0	97,103,111	0.37	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	LMN	B	1002	-	72,72,72	0.17	0	96,98,98	0.48	0
3	P5S	A	602	-	39,40,53	0.29	0	43,47,60	0.30	0
3	P5S	D	1004	-	39,40,53	0.29	0	43,47,60	0.30	0
2	LMN	E	1003	-	48,48,72	0.14	0	61,62,98	0.33	0
2	LMN	A	604	-	72,72,72	0.17	0	96,98,98	0.48	0
4	CDL	B	1001	-	91,91,99	0.30	0	97,103,111	0.37	0
2	LMN	D	1003	-	48,48,72	0.13	0	61,62,98	0.33	0
3	P5S	B	1004	-	39,40,53	0.29	0	43,47,60	0.30	0
4	CDL	F	1001	-	91,91,99	0.30	0	97,103,111	0.37	0
4	CDL	E	1001	-	91,91,99	0.30	0	97,103,111	0.37	0
2	LMN	F	1003	-	48,48,72	0.13	0	61,62,98	0.33	0
2	LMN	A	601	-	48,48,72	0.14	0	61,62,98	0.33	0
2	LMN	C	1003	-	48,48,72	0.14	0	61,62,98	0.33	0
2	LMN	F	1002	-	72,72,72	0.17	0	96,98,98	0.48	0
3	P5S	E	1004	-	39,40,53	0.29	0	43,47,60	0.30	0
3	P5S	F	1004	-	39,40,53	0.29	0	43,47,60	0.30	0
2	LMN	C	1002	-	72,72,72	0.17	0	96,98,98	0.48	0
2	LMN	E	1002	-	72,72,72	0.17	0	96,98,98	0.48	0
3	P5S	C	1004	-	39,40,53	0.29	0	43,47,60	0.30	0
2	LMN	B	1003	-	48,48,72	0.14	0	61,62,98	0.33	0
4	CDL	D	1001	-	91,91,99	0.30	0	97,103,111	0.37	0
2	LMN	D	1002	-	72,72,72	0.17	0	96,98,98	0.48	0
4	CDL	C	1001	-	91,91,99	0.30	0	97,103,111	0.37	0

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. '2' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	CDL	A	603	-	-	29/102/102/110	-
2	LMN	B	1002	-	-	12/50/130/130	0/4/4/4
3	P5S	A	602	-	-	9/46/46/59	-
3	P5S	D	1004	-	-	9/46/46/59	-
2	LMN	E	1003	-	-	13/39/79/130	0/2/2/4
2	LMN	A	604	-	-	12/50/130/130	0/4/4/4
4	CDL	B	1001	-	-	29/102/102/110	-
2	LMN	D	1003	-	-	13/39/79/130	0/2/2/4
3	P5S	B	1004	-	-	9/46/46/59	-
4	CDL	F	1001	-	-	29/102/102/110	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	CDL	E	1001	-	-	29/102/102/110	-
2	LMN	F	1003	-	-	13/39/79/130	0/2/2/4
2	LMN	A	601	-	-	13/39/79/130	0/2/2/4
2	LMN	C	1003	-	-	13/39/79/130	0/2/2/4
2	LMN	F	1002	-	-	12/50/130/130	0/4/4/4
3	P5S	E	1004	-	-	9/46/46/59	-
3	P5S	F	1004	-	-	9/46/46/59	-
2	LMN	C	1002	-	-	12/50/130/130	0/4/4/4
2	LMN	E	1002	-	-	12/50/130/130	0/4/4/4
3	P5S	C	1004	-	-	9/46/46/59	-
2	LMN	B	1003	-	-	13/39/79/130	0/2/2/4
4	CDL	D	1001	-	-	29/102/102/110	-
2	LMN	D	1002	-	-	12/50/130/130	0/4/4/4
4	CDL	C	1001	-	-	29/102/102/110	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (378) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	601	LMN	CBL-CBR-CCM-CBQ
2	A	601	LMN	CBL-CBR-CCM-CBS
2	A	601	LMN	CBL-CBR-CCM-CBT
2	A	604	LMN	O5-C1-O1-CBS
2	A	604	LMN	OBX-CCJ-OBV-CBT
2	B	1002	LMN	O5-C1-O1-CBS
2	B	1002	LMN	OBX-CCJ-OBV-CBT
2	B	1003	LMN	CBL-CBR-CCM-CBQ
2	B	1003	LMN	CBL-CBR-CCM-CBS
2	B	1003	LMN	CBL-CBR-CCM-CBT
2	C	1002	LMN	O5-C1-O1-CBS
2	C	1002	LMN	OBX-CCJ-OBV-CBT
2	C	1003	LMN	CBL-CBR-CCM-CBQ
2	C	1003	LMN	CBL-CBR-CCM-CBS
2	C	1003	LMN	CBL-CBR-CCM-CBT
2	D	1002	LMN	O5-C1-O1-CBS
2	D	1002	LMN	OBX-CCJ-OBV-CBT

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Mol	Chain	Res	Type	Atoms
2	D	1003	LMN	CBL-CBR-CCM-CBQ
2	D	1003	LMN	CBL-CBR-CCM-CBS
2	D	1003	LMN	CBL-CBR-CCM-CBT
2	E	1002	LMN	O5-C1-O1-CBS
2	E	1002	LMN	OBX-CCJ-OBV-CBT
2	E	1003	LMN	CBL-CBR-CCM-CBQ
2	E	1003	LMN	CBL-CBR-CCM-CBS
2	E	1003	LMN	CBL-CBR-CCM-CBT
2	F	1002	LMN	O5-C1-O1-CBS
2	F	1002	LMN	OBX-CCJ-OBV-CBT
2	F	1003	LMN	CBL-CBR-CCM-CBQ
2	F	1003	LMN	CBL-CBR-CCM-CBS
2	F	1003	LMN	CBL-CBR-CCM-CBT
3	A	602	P5S	O-C-CA-CB
3	A	602	P5S	OXT-C-CA-CB
3	A	602	P5S	N-CA-CB-OG
3	A	602	P5S	CB-OG-P12-O15
3	B	1004	P5S	O-C-CA-CB
3	B	1004	P5S	OXT-C-CA-CB
3	B	1004	P5S	N-CA-CB-OG
3	B	1004	P5S	CB-OG-P12-O15
3	C	1004	P5S	O-C-CA-CB
3	C	1004	P5S	OXT-C-CA-CB
3	C	1004	P5S	N-CA-CB-OG
3	C	1004	P5S	CB-OG-P12-O15
3	D	1004	P5S	O-C-CA-CB
3	D	1004	P5S	OXT-C-CA-CB
3	D	1004	P5S	N-CA-CB-OG
3	D	1004	P5S	CB-OG-P12-O15
3	E	1004	P5S	O-C-CA-CB
3	E	1004	P5S	OXT-C-CA-CB
3	E	1004	P5S	N-CA-CB-OG
3	E	1004	P5S	CB-OG-P12-O15
3	F	1004	P5S	O-C-CA-CB
3	F	1004	P5S	OXT-C-CA-CB
3	F	1004	P5S	N-CA-CB-OG
3	F	1004	P5S	CB-OG-P12-O15
4	A	603	CDL	CA3-OA5-PA1-OA3
4	A	603	CDL	CA4-CA3-OA5-PA1
4	B	1001	CDL	CA3-OA5-PA1-OA3
4	B	1001	CDL	CA4-CA3-OA5-PA1
4	C	1001	CDL	CA3-OA5-PA1-OA3

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Mol	Chain	Res	Type	Atoms
4	C	1001	CDL	CA4-CA3-OA5-PA1
4	D	1001	CDL	CA3-OA5-PA1-OA3
4	D	1001	CDL	CA4-CA3-OA5-PA1
4	E	1001	CDL	CA3-OA5-PA1-OA3
4	E	1001	CDL	CA4-CA3-OA5-PA1
4	F	1001	CDL	CA3-OA5-PA1-OA3
4	F	1001	CDL	CA4-CA3-OA5-PA1
2	A	601	LMN	CCH-CCQ-OCB-CCS
2	B	1003	LMN	CCH-CCQ-OCB-CCS
2	C	1003	LMN	CCH-CCQ-OCB-CCS
2	D	1003	LMN	CCH-CCQ-OCB-CCS
2	E	1003	LMN	CCH-CCQ-OCB-CCS
2	F	1003	LMN	CCH-CCQ-OCB-CCS
2	A	604	LMN	CCH-CCQ-OCB-CCS
2	B	1002	LMN	CCH-CCQ-OCB-CCS
2	C	1002	LMN	CCH-CCQ-OCB-CCS
2	D	1002	LMN	CCH-CCQ-OCB-CCS
2	E	1002	LMN	CCH-CCQ-OCB-CCS
2	F	1002	LMN	CCH-CCQ-OCB-CCS
2	A	604	LMN	OBY-CCR-O4-C4
2	B	1002	LMN	OBY-CCR-O4-C4
2	C	1002	LMN	OBY-CCR-O4-C4
2	D	1002	LMN	OBY-CCR-O4-C4
2	E	1002	LMN	OBY-CCR-O4-C4
2	F	1002	LMN	OBY-CCR-O4-C4
2	A	604	LMN	CCV-CCR-O4-C4
2	B	1002	LMN	CCV-CCR-O4-C4
2	C	1002	LMN	CCV-CCR-O4-C4
2	D	1002	LMN	CCV-CCR-O4-C4
2	E	1002	LMN	CCV-CCR-O4-C4
2	F	1002	LMN	CCV-CCR-O4-C4
3	A	602	P5S	OXT-C-CA-N
3	B	1004	P5S	OXT-C-CA-N
3	C	1004	P5S	OXT-C-CA-N
3	D	1004	P5S	OXT-C-CA-N
3	E	1004	P5S	OXT-C-CA-N
3	F	1004	P5S	OXT-C-CA-N
2	A	604	LMN	C2-C1-O1-CBS
2	B	1002	LMN	C2-C1-O1-CBS
2	C	1002	LMN	C2-C1-O1-CBS
2	D	1002	LMN	C2-C1-O1-CBS
2	E	1002	LMN	C2-C1-O1-CBS

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Mol	Chain	Res	Type	Atoms
2	F	1002	LMN	C2-C1-O1-CBS
4	A	603	CDL	CB7-C71-C72-C73
4	B	1001	CDL	CB7-C71-C72-C73
4	C	1001	CDL	CB7-C71-C72-C73
4	D	1001	CDL	CB7-C71-C72-C73
4	E	1001	CDL	CB7-C71-C72-C73
4	F	1001	CDL	CB7-C71-C72-C73
4	A	603	CDL	CA3-OA5-PA1-OA2
4	B	1001	CDL	CA3-OA5-PA1-OA2
4	C	1001	CDL	CA3-OA5-PA1-OA2
4	D	1001	CDL	CA3-OA5-PA1-OA2
4	E	1001	CDL	CA3-OA5-PA1-OA2
4	F	1001	CDL	CA3-OA5-PA1-OA2
4	B	1001	CDL	C73-C74-C75-C76
4	C	1001	CDL	C73-C74-C75-C76
4	D	1001	CDL	C73-C74-C75-C76
4	F	1001	CDL	C73-C74-C75-C76
4	A	603	CDL	C73-C74-C75-C76
4	E	1001	CDL	C73-C74-C75-C76
2	A	601	LMN	CCL-CCJ-OBV-CBT
2	A	604	LMN	CCL-CCJ-OBV-CBT
2	B	1002	LMN	CCL-CCJ-OBV-CBT
2	B	1003	LMN	CCL-CCJ-OBV-CBT
2	C	1002	LMN	CCL-CCJ-OBV-CBT
2	C	1003	LMN	CCL-CCJ-OBV-CBT
2	D	1002	LMN	CCL-CCJ-OBV-CBT
2	D	1003	LMN	CCL-CCJ-OBV-CBT
2	E	1002	LMN	CCL-CCJ-OBV-CBT
2	E	1003	LMN	CCL-CCJ-OBV-CBT
2	F	1002	LMN	CCL-CCJ-OBV-CBT
2	F	1003	LMN	CCL-CCJ-OBV-CBT
4	A	603	CDL	C41-C42-C43-C44
4	B	1001	CDL	C41-C42-C43-C44
4	C	1001	CDL	C41-C42-C43-C44
4	D	1001	CDL	C41-C42-C43-C44
4	E	1001	CDL	C41-C42-C43-C44
4	F	1001	CDL	C41-C42-C43-C44
3	A	602	P5S	C-CA-CB-OG
3	B	1004	P5S	C-CA-CB-OG
3	C	1004	P5S	C-CA-CB-OG
3	D	1004	P5S	C-CA-CB-OG
3	E	1004	P5S	C-CA-CB-OG

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Mol	Chain	Res	Type	Atoms
3	F	1004	P5S	C-CA-CB-OG
2	A	601	LMN	OBX-CCJ-OBV-CBT
2	B	1003	LMN	OBX-CCJ-OBV-CBT
2	C	1003	LMN	OBX-CCJ-OBV-CBT
2	D	1003	LMN	OBX-CCJ-OBV-CBT
2	E	1003	LMN	OBX-CCJ-OBV-CBT
2	F	1003	LMN	OBX-CCJ-OBV-CBT
2	A	604	LMN	O5-C5-C6-O6
2	B	1002	LMN	O5-C5-C6-O6
2	C	1002	LMN	O5-C5-C6-O6
2	D	1002	LMN	O5-C5-C6-O6
2	E	1002	LMN	O5-C5-C6-O6
2	F	1002	LMN	O5-C5-C6-O6
2	A	604	LMN	OAI-CBM-CCC-OBY
2	B	1002	LMN	OAI-CBM-CCC-OBY
2	C	1002	LMN	OAI-CBM-CCC-OBY
2	D	1002	LMN	OAI-CBM-CCC-OBY
2	E	1002	LMN	OAI-CBM-CCC-OBY
2	F	1002	LMN	OAI-CBM-CCC-OBY
2	B	1002	LMN	OAJ-CBN-CCD-OBZ
2	C	1002	LMN	OAJ-CBN-CCD-OBZ
2	E	1002	LMN	OAJ-CBN-CCD-OBZ
2	F	1002	LMN	OAJ-CBN-CCD-OBZ
2	A	604	LMN	CBI-CBK-CBQ-CCM
2	B	1002	LMN	CBI-CBK-CBQ-CCM
2	C	1002	LMN	CBI-CBK-CBQ-CCM
2	D	1002	LMN	CBI-CBK-CBQ-CCM
2	E	1002	LMN	CBI-CBK-CBQ-CCM
2	F	1002	LMN	CBI-CBK-CBQ-CCM
2	A	604	LMN	OAJ-CBN-CCD-OBZ
2	D	1002	LMN	OAJ-CBN-CCD-OBZ
2	A	601	LMN	CBB-CBD-CBF-CBH
2	C	1003	LMN	CBB-CBD-CBF-CBH
2	E	1003	LMN	CBB-CBD-CBF-CBH
2	F	1003	LMN	CBB-CBD-CBF-CBH
2	B	1003	LMN	CBB-CBD-CBF-CBH
2	D	1003	LMN	CBB-CBD-CBF-CBH
4	A	603	CDL	C78-C79-C80-C81
4	B	1001	CDL	C78-C79-C80-C81
4	F	1001	CDL	C78-C79-C80-C81
4	C	1001	CDL	C78-C79-C80-C81
4	D	1001	CDL	C78-C79-C80-C81

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Mol	Chain	Res	Type	Atoms
4	E	1001	CDL	C78-C79-C80-C81
2	E	1003	LMN	OAJ-CBN-CCD-OBZ
2	A	601	LMN	OAJ-CBN-CCD-OBZ
2	B	1003	LMN	OAJ-CBN-CCD-OBZ
2	C	1003	LMN	OAJ-CBN-CCD-OBZ
2	D	1003	LMN	OAJ-CBN-CCD-OBZ
2	F	1003	LMN	OAJ-CBN-CCD-OBZ
4	A	603	CDL	C71-C72-C73-C74
4	B	1001	CDL	C71-C72-C73-C74
4	C	1001	CDL	C71-C72-C73-C74
4	D	1001	CDL	C71-C72-C73-C74
4	E	1001	CDL	C71-C72-C73-C74
4	F	1001	CDL	C71-C72-C73-C74
4	A	603	CDL	C77-C78-C79-C80
4	B	1001	CDL	C77-C78-C79-C80
4	C	1001	CDL	C77-C78-C79-C80
4	D	1001	CDL	C77-C78-C79-C80
4	F	1001	CDL	C77-C78-C79-C80
4	E	1001	CDL	C77-C78-C79-C80
4	A	603	CDL	OA5-CA3-CA4-CA6
4	B	1001	CDL	OA5-CA3-CA4-CA6
4	C	1001	CDL	OA5-CA3-CA4-CA6
4	D	1001	CDL	OA5-CA3-CA4-CA6
4	E	1001	CDL	OA5-CA3-CA4-CA6
4	F	1001	CDL	OA5-CA3-CA4-CA6
2	A	604	LMN	CCF-CCQ-OCB-CCS
2	F	1002	LMN	CCF-CCQ-OCB-CCS
2	B	1002	LMN	CCF-CCQ-OCB-CCS
2	C	1002	LMN	CCF-CCQ-OCB-CCS
2	D	1002	LMN	CCF-CCQ-OCB-CCS
2	E	1002	LMN	CCF-CCQ-OCB-CCS
4	A	603	CDL	OB5-CB3-CB4-OB6
4	B	1001	CDL	OB5-CB3-CB4-OB6
4	C	1001	CDL	OB5-CB3-CB4-OB6
4	D	1001	CDL	OB5-CB3-CB4-OB6
4	E	1001	CDL	OB5-CB3-CB4-OB6
4	F	1001	CDL	OB5-CB3-CB4-OB6
4	A	603	CDL	C79-C80-C81-C82
4	C	1001	CDL	C79-C80-C81-C82
4	E	1001	CDL	C79-C80-C81-C82
4	B	1001	CDL	C79-C80-C81-C82
4	D	1001	CDL	C79-C80-C81-C82

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Mol	Chain	Res	Type	Atoms
4	F	1001	CDL	C79-C80-C81-C82
3	A	602	P5S	O-C-CA-N
3	B	1004	P5S	O-C-CA-N
3	C	1004	P5S	O-C-CA-N
3	D	1004	P5S	O-C-CA-N
3	E	1004	P5S	O-C-CA-N
3	F	1004	P5S	O-C-CA-N
4	A	603	CDL	OB5-CB3-CB4-CB6
4	B	1001	CDL	OB5-CB3-CB4-CB6
4	C	1001	CDL	OB5-CB3-CB4-CB6
4	D	1001	CDL	OB5-CB3-CB4-CB6
4	E	1001	CDL	OB5-CB3-CB4-CB6
4	F	1001	CDL	OB5-CB3-CB4-CB6
4	A	603	CDL	OA5-CA3-CA4-OA6
4	B	1001	CDL	OA5-CA3-CA4-OA6
4	C	1001	CDL	OA5-CA3-CA4-OA6
4	D	1001	CDL	OA5-CA3-CA4-OA6
4	E	1001	CDL	OA5-CA3-CA4-OA6
4	F	1001	CDL	OA5-CA3-CA4-OA6
4	A	603	CDL	C11-CA5-OA6-CA4
4	E	1001	CDL	C11-CA5-OA6-CA4
4	B	1001	CDL	C11-CA5-OA6-CA4
4	C	1001	CDL	C11-CA5-OA6-CA4
4	D	1001	CDL	C11-CA5-OA6-CA4
4	F	1001	CDL	C11-CA5-OA6-CA4
3	A	602	P5S	CB-OG-P12-O13
3	B	1004	P5S	CB-OG-P12-O13
3	C	1004	P5S	CB-OG-P12-O13
3	D	1004	P5S	CB-OG-P12-O13
3	E	1004	P5S	CB-OG-P12-O13
3	F	1004	P5S	CB-OG-P12-O13
4	A	603	CDL	CA3-OA5-PA1-OA4
4	B	1001	CDL	CA3-OA5-PA1-OA4
4	C	1001	CDL	CA3-OA5-PA1-OA4
4	D	1001	CDL	CA3-OA5-PA1-OA4
4	E	1001	CDL	CA3-OA5-PA1-OA4
4	F	1001	CDL	CA3-OA5-PA1-OA4
4	A	603	CDL	CA2-OA2-PA1-OA5
4	B	1001	CDL	CA2-OA2-PA1-OA5
4	C	1001	CDL	CA2-OA2-PA1-OA5
4	D	1001	CDL	CA2-OA2-PA1-OA5
4	E	1001	CDL	CA2-OA2-PA1-OA5

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Mol	Chain	Res	Type	Atoms
4	F	1001	CDL	CA2-OA2-PA1-OA5
2	D	1003	LMN	CBH-CBJ-CBL-CBR
2	F	1003	LMN	CBH-CBJ-CBL-CBR
2	B	1003	LMN	CBH-CBJ-CBL-CBR
2	E	1003	LMN	CBH-CBJ-CBL-CBR
2	A	601	LMN	CBH-CBJ-CBL-CBR
2	C	1003	LMN	CBH-CBJ-CBL-CBR
4	B	1001	CDL	OA7-CA5-OA6-CA4
4	C	1001	CDL	OA7-CA5-OA6-CA4
4	D	1001	CDL	OA7-CA5-OA6-CA4
4	F	1001	CDL	OA7-CA5-OA6-CA4
4	A	603	CDL	OA7-CA5-OA6-CA4
4	E	1001	CDL	OA7-CA5-OA6-CA4
4	A	603	CDL	CA6-CA4-OA6-CA5
4	B	1001	CDL	CA6-CA4-OA6-CA5
4	C	1001	CDL	CA6-CA4-OA6-CA5
4	D	1001	CDL	CA6-CA4-OA6-CA5
4	E	1001	CDL	CA6-CA4-OA6-CA5
4	F	1001	CDL	CA6-CA4-OA6-CA5
2	F	1003	LMN	CBA-CBC-CBE-CBG
2	A	601	LMN	CBA-CBC-CBE-CBG
2	B	1003	LMN	CBA-CBC-CBE-CBG
2	D	1003	LMN	CBA-CBC-CBE-CBG
2	C	1003	LMN	CBA-CBC-CBE-CBG
2	E	1003	LMN	CBA-CBC-CBE-CBG
4	E	1001	CDL	C83-C84-C85-C86
4	C	1001	CDL	C83-C84-C85-C86
4	F	1001	CDL	C83-C84-C85-C86
4	B	1001	CDL	C83-C84-C85-C86
4	A	603	CDL	C83-C84-C85-C86
4	D	1001	CDL	C83-C84-C85-C86
4	D	1001	CDL	C11-C12-C13-C14
4	B	1001	CDL	C11-C12-C13-C14
4	C	1001	CDL	C11-C12-C13-C14
4	E	1001	CDL	C11-C12-C13-C14
4	A	603	CDL	C11-C12-C13-C14
4	F	1001	CDL	C11-C12-C13-C14
4	A	603	CDL	C72-C71-CB7-OB8
4	C	1001	CDL	C72-C71-CB7-OB8
4	D	1001	CDL	C72-C71-CB7-OB8
4	F	1001	CDL	C72-C71-CB7-OB8
4	B	1001	CDL	C72-C71-CB7-OB8

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Mol	Chain	Res	Type	Atoms
4	E	1001	CDL	C72-C71-CB7-OB8
2	A	601	LMN	OBV-CBT-CCM-CBQ
2	B	1003	LMN	OBV-CBT-CCM-CBQ
2	C	1003	LMN	OBV-CBT-CCM-CBQ
2	D	1003	LMN	OBV-CBT-CCM-CBQ
2	E	1003	LMN	OBV-CBT-CCM-CBQ
2	F	1003	LMN	OBV-CBT-CCM-CBQ
4	A	603	CDL	C12-C11-CA5-OA6
4	B	1001	CDL	C12-C11-CA5-OA6
4	C	1001	CDL	C12-C11-CA5-OA6
4	D	1001	CDL	C12-C11-CA5-OA6
4	E	1001	CDL	C12-C11-CA5-OA6
4	F	1001	CDL	C12-C11-CA5-OA6
4	B	1001	CDL	C19-C20-C21-C22
4	E	1001	CDL	C19-C20-C21-C22
4	D	1001	CDL	C19-C20-C21-C22
4	A	603	CDL	C19-C20-C21-C22
4	F	1001	CDL	C19-C20-C21-C22
4	E	1001	CDL	C82-C83-C84-C85
4	B	1001	CDL	C82-C83-C84-C85
4	C	1001	CDL	C19-C20-C21-C22
3	A	602	P5S	CB-OG-P12-O16
3	B	1004	P5S	CB-OG-P12-O16
3	C	1004	P5S	CB-OG-P12-O16
3	D	1004	P5S	CB-OG-P12-O16
3	E	1004	P5S	CB-OG-P12-O16
3	F	1004	P5S	CB-OG-P12-O16
4	C	1001	CDL	C82-C83-C84-C85
4	F	1001	CDL	C82-C83-C84-C85
4	A	603	CDL	C82-C83-C84-C85
4	D	1001	CDL	C82-C83-C84-C85
4	A	603	CDL	CB3-OB5-PB2-OB3
4	B	1001	CDL	CB3-OB5-PB2-OB3
4	C	1001	CDL	CB3-OB5-PB2-OB3
4	D	1001	CDL	CB3-OB5-PB2-OB3
4	E	1001	CDL	CB3-OB5-PB2-OB3
4	F	1001	CDL	CB3-OB5-PB2-OB3
4	A	603	CDL	C12-C11-CA5-OA7
4	A	603	CDL	CA3-CA4-OA6-CA5
4	B	1001	CDL	CA3-CA4-OA6-CA5
4	C	1001	CDL	CA3-CA4-OA6-CA5
4	D	1001	CDL	CA3-CA4-OA6-CA5

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Mol	Chain	Res	Type	Atoms
4	E	1001	CDL	CA3-CA4-OA6-CA5
4	F	1001	CDL	CA3-CA4-OA6-CA5
4	A	603	CDL	C72-C71-CB7-OB9
4	B	1001	CDL	C72-C71-CB7-OB9
4	C	1001	CDL	C72-C71-CB7-OB9
4	D	1001	CDL	C72-C71-CB7-OB9
4	E	1001	CDL	C72-C71-CB7-OB9
4	F	1001	CDL	C72-C71-CB7-OB9
4	B	1001	CDL	C12-C11-CA5-OA7
4	C	1001	CDL	C12-C11-CA5-OA7
4	D	1001	CDL	C12-C11-CA5-OA7
4	E	1001	CDL	C12-C11-CA5-OA7
4	F	1001	CDL	C12-C11-CA5-OA7
2	B	1003	LMN	CAZ-CBB-CBD-CBF
2	D	1003	LMN	CAZ-CBB-CBD-CBF
2	A	601	LMN	CAZ-CBB-CBD-CBF
2	C	1003	LMN	CAZ-CBB-CBD-CBF
2	D	1003	LMN	CCF-CCQ-OCB-CCS
2	E	1003	LMN	CAZ-CBB-CBD-CBF
2	E	1003	LMN	CCF-CCQ-OCB-CCS
2	F	1003	LMN	CAZ-CBB-CBD-CBF
2	A	601	LMN	CCF-CCQ-OCB-CCS
2	C	1003	LMN	CCF-CCQ-OCB-CCS
2	F	1003	LMN	CCF-CCQ-OCB-CCS
2	B	1003	LMN	CCF-CCQ-OCB-CCS

There are no ring outliers.

24 monomers are involved in 147 short contacts:

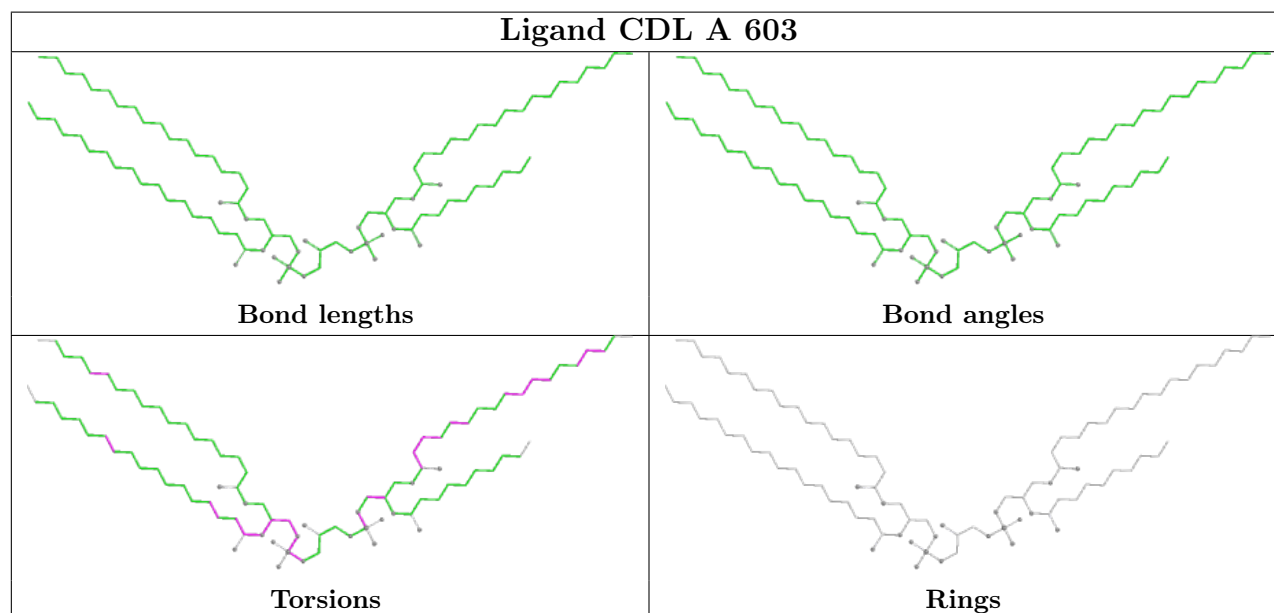
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4	A	603	CDL	3	0
2	B	1002	LMN	16	0
3	A	602	P5S	1	0
3	D	1004	P5S	1	0
2	E	1003	LMN	4	0
2	A	604	LMN	17	0
4	B	1001	CDL	3	0
2	D	1003	LMN	3	0
3	B	1004	P5S	1	0
4	F	1001	CDL	3	0
4	E	1001	CDL	4	0
2	F	1003	LMN	3	0

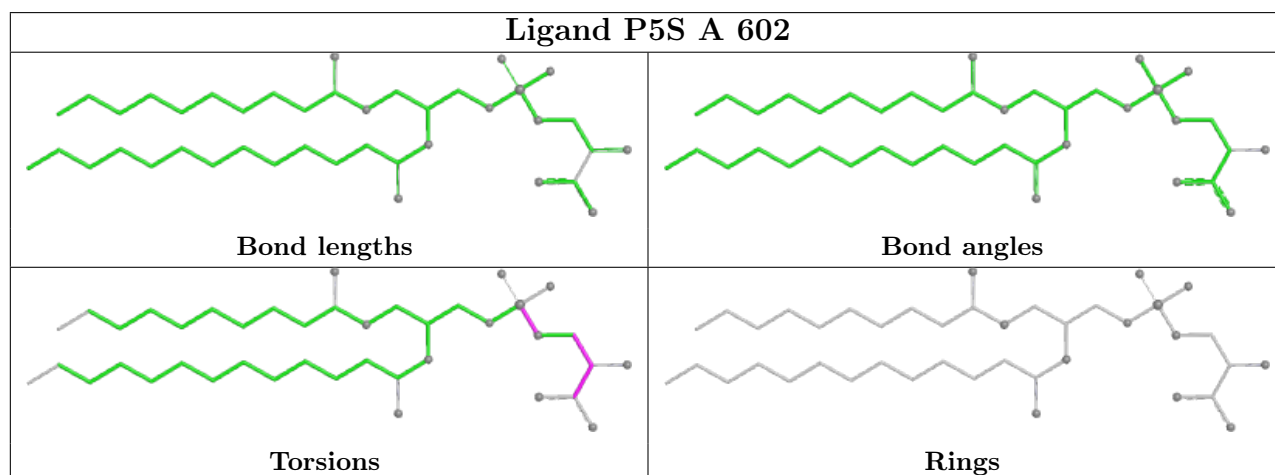
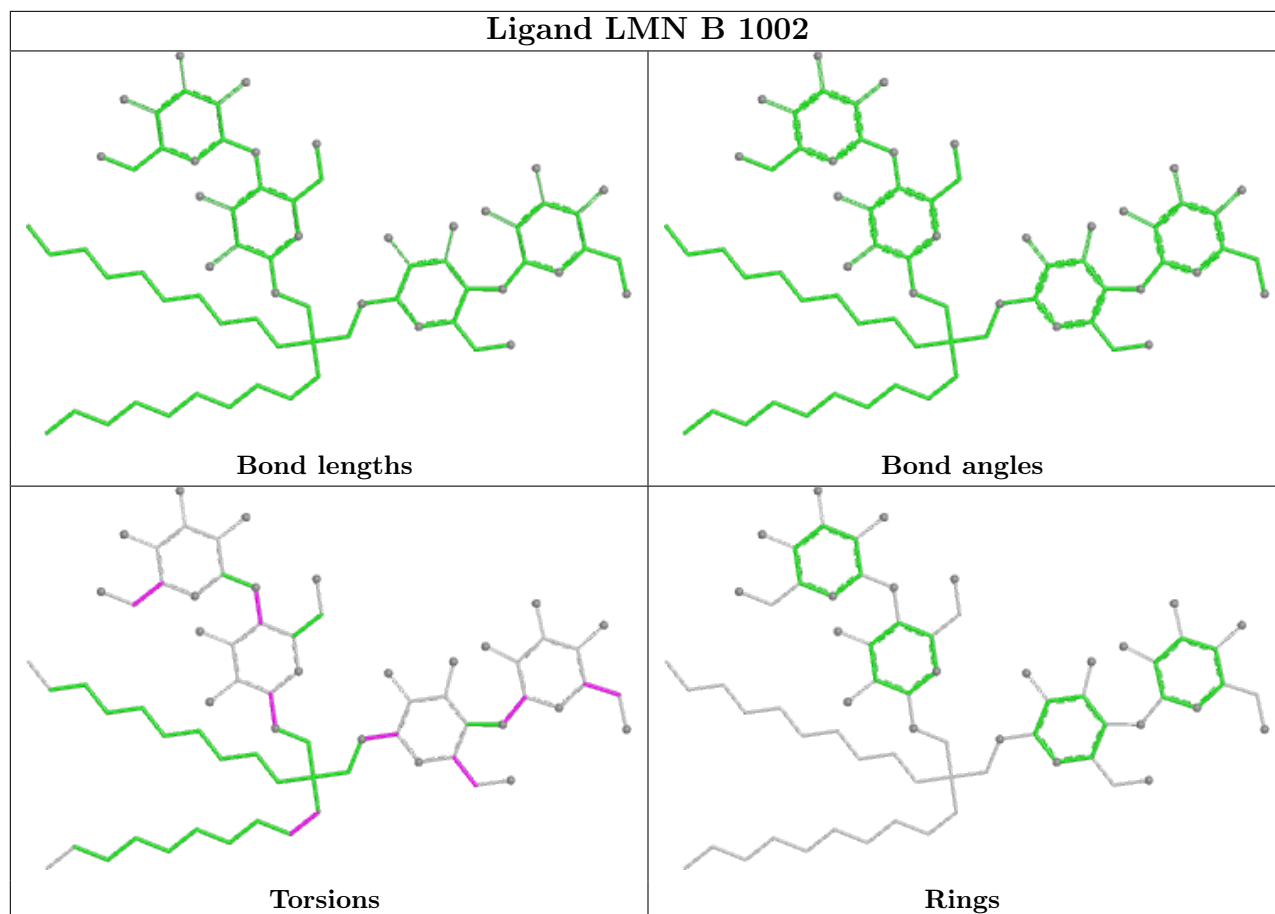
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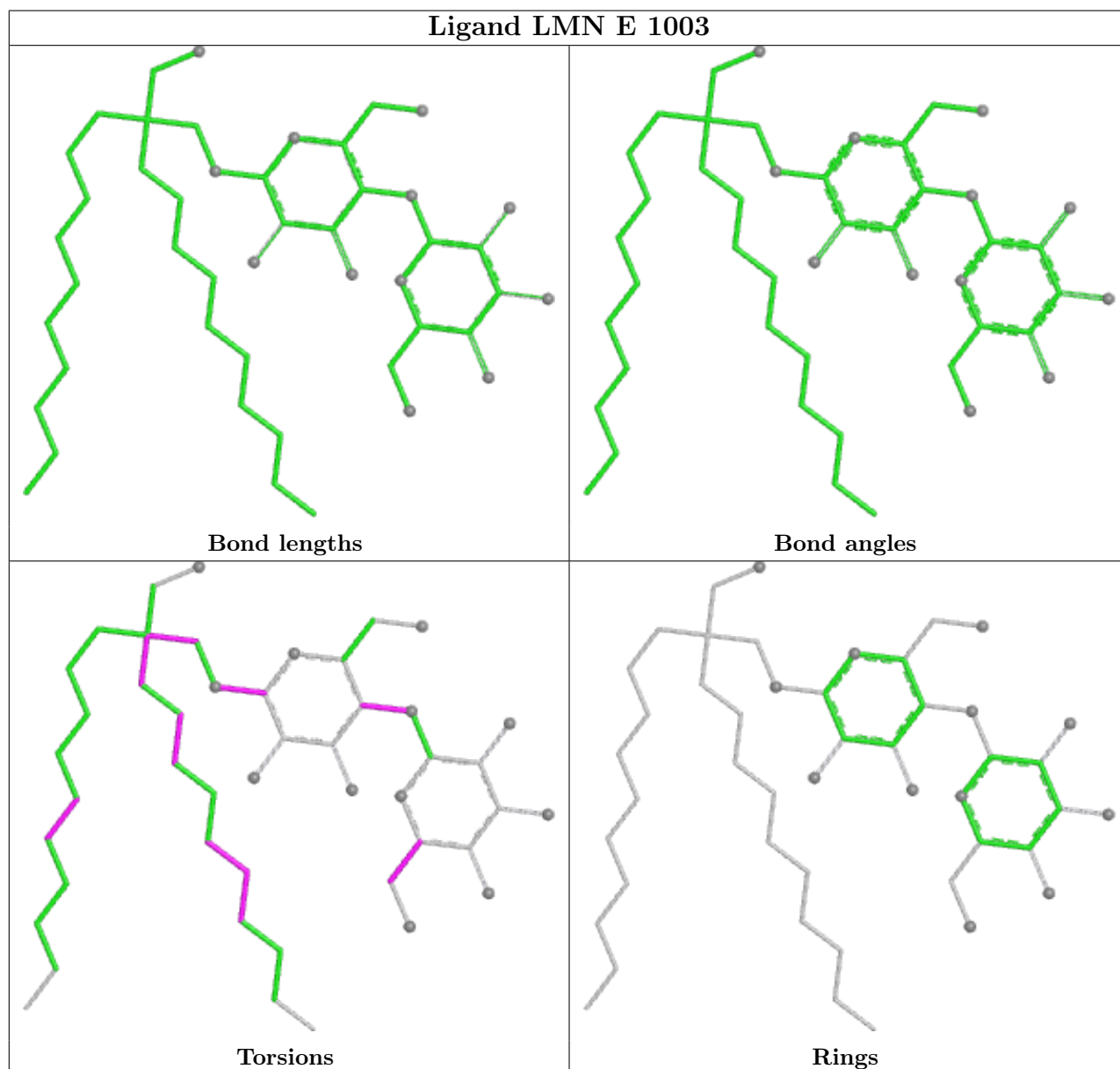
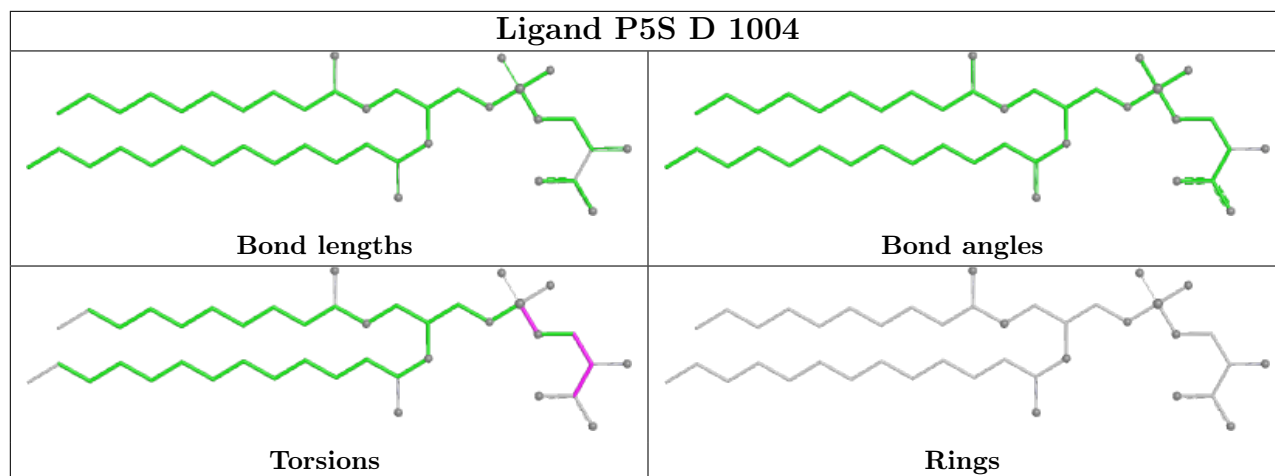
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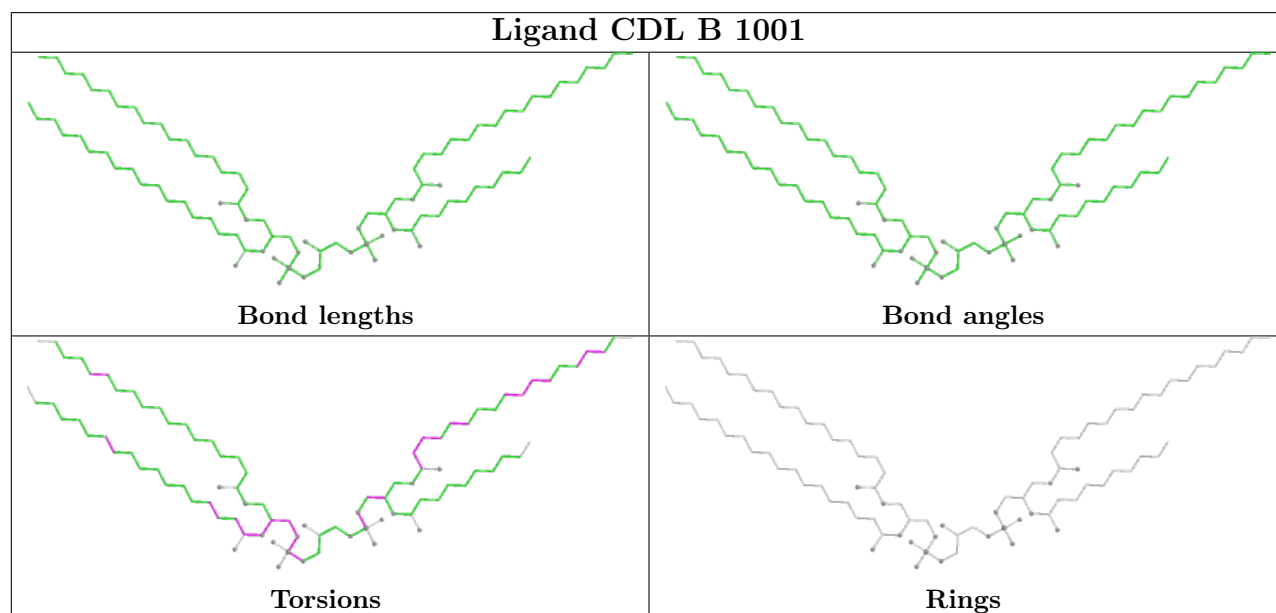
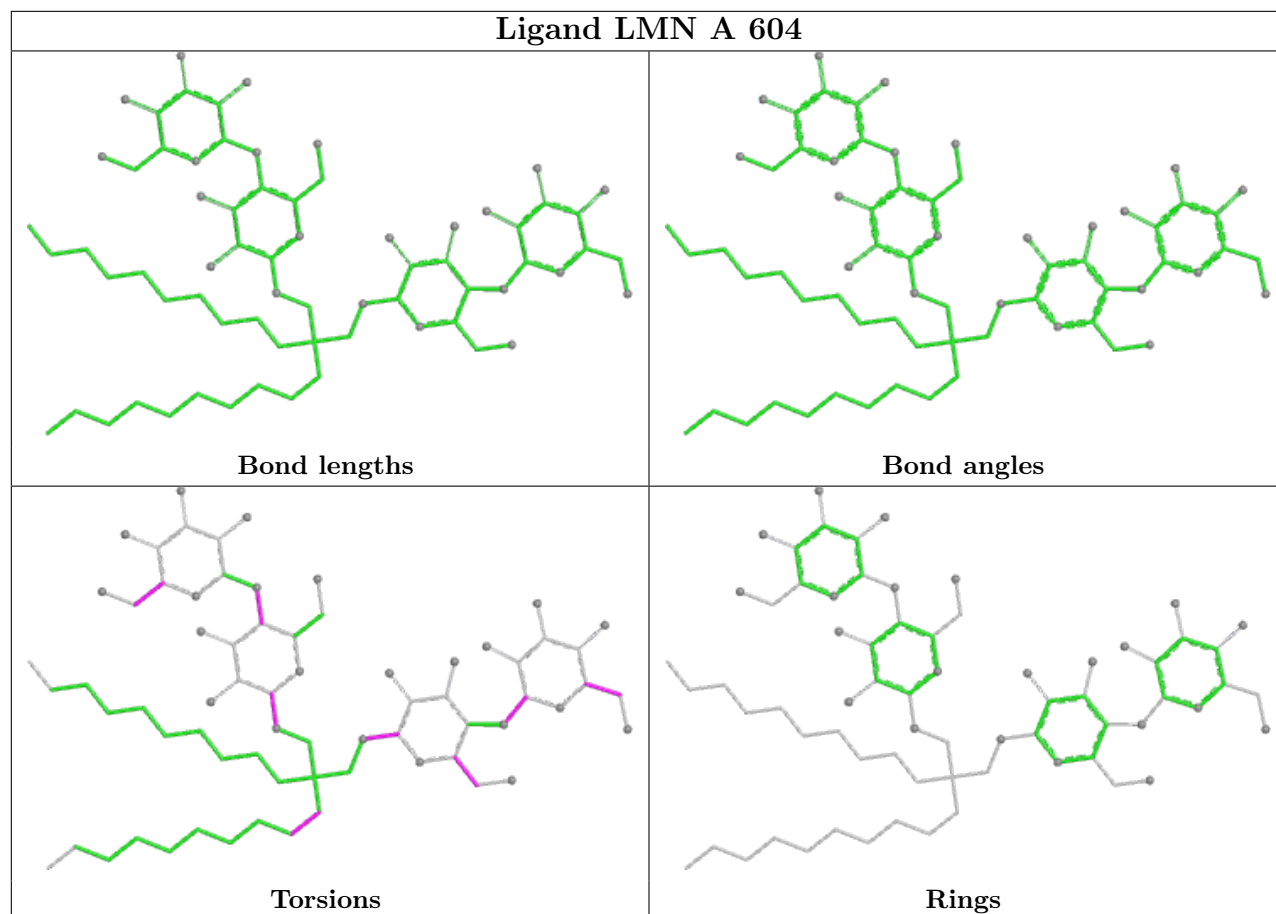
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	601	LMN	4	0
2	C	1003	LMN	4	0
2	F	1002	LMN	17	0
3	E	1004	P5S	1	0
3	F	1004	P5S	1	0
2	C	1002	LMN	14	0
2	E	1002	LMN	17	0
3	C	1004	P5S	1	0
2	B	1003	LMN	4	0
4	D	1001	CDL	4	0
2	D	1002	LMN	17	0
4	C	1001	CDL	4	0

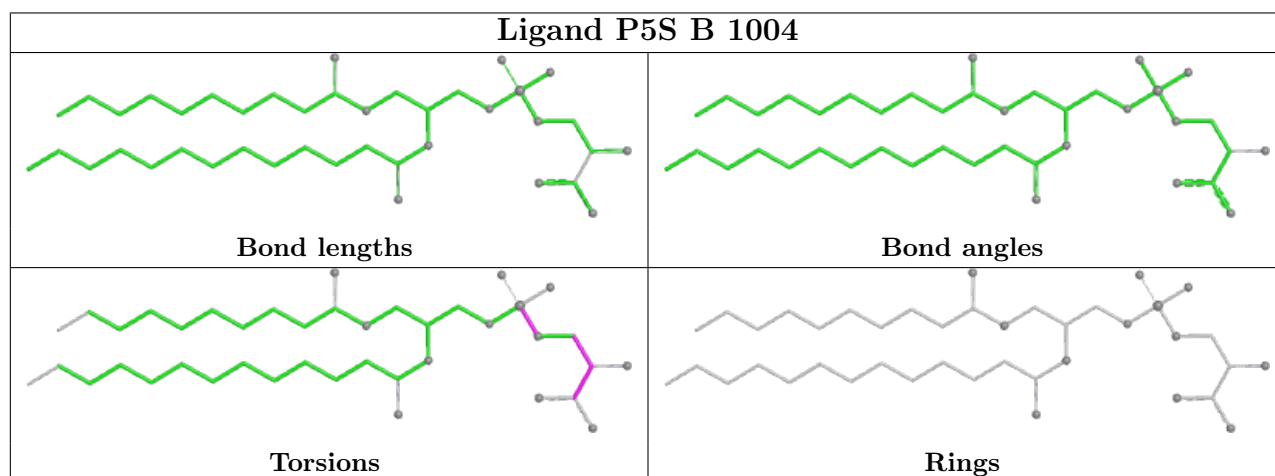
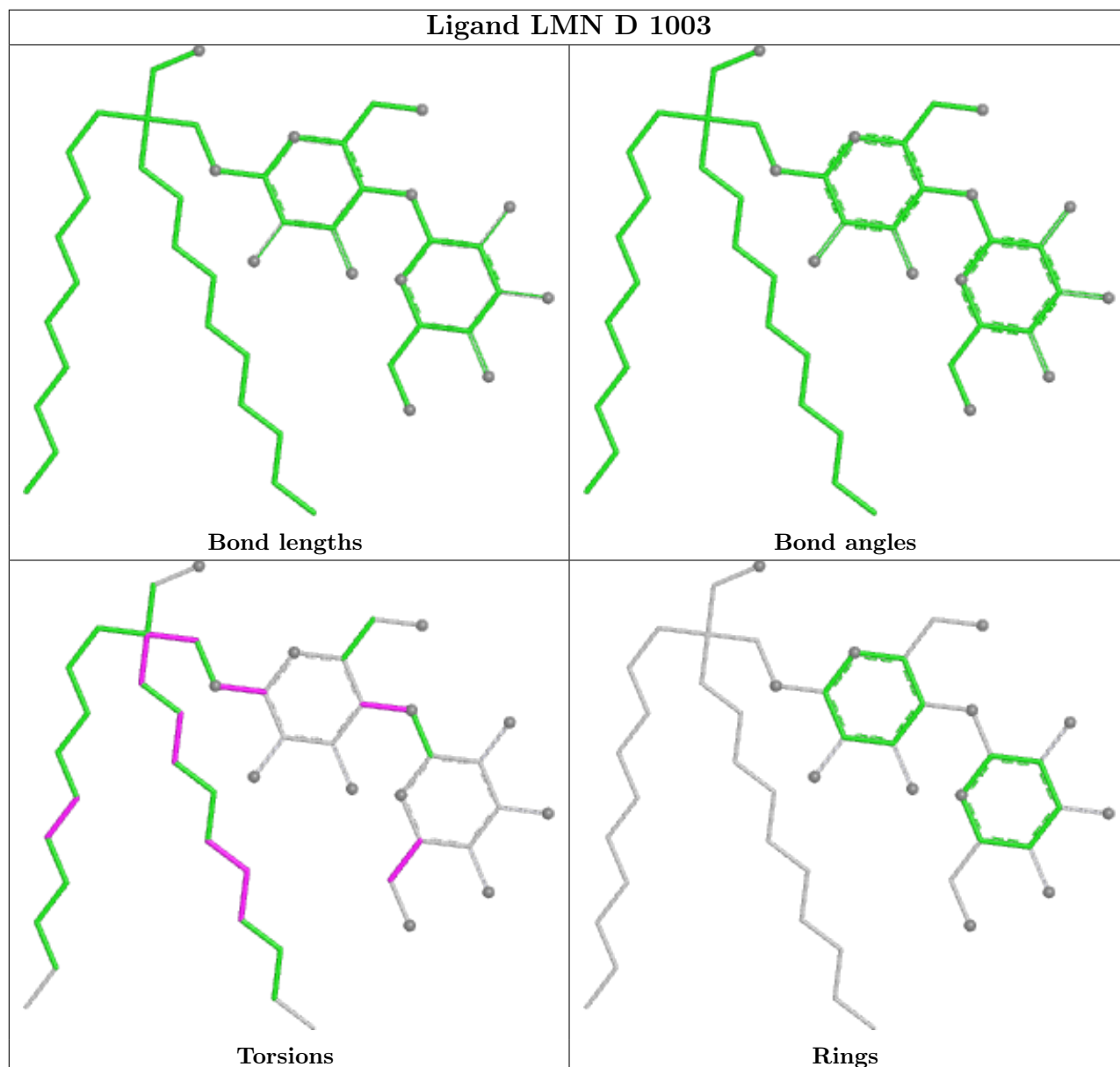
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

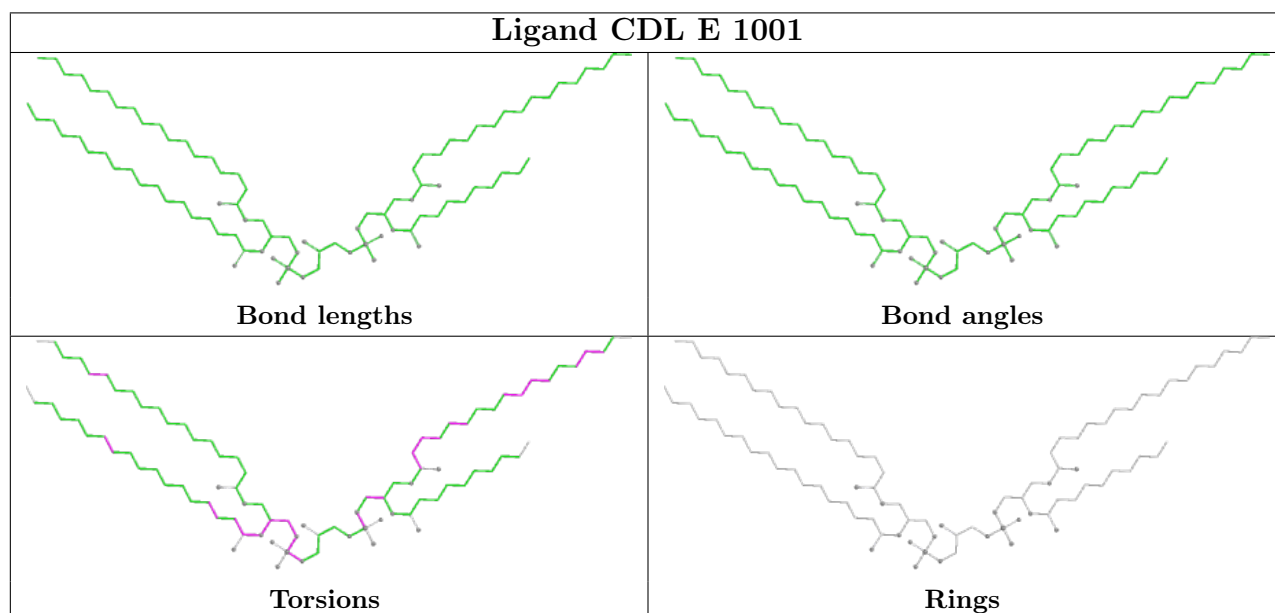
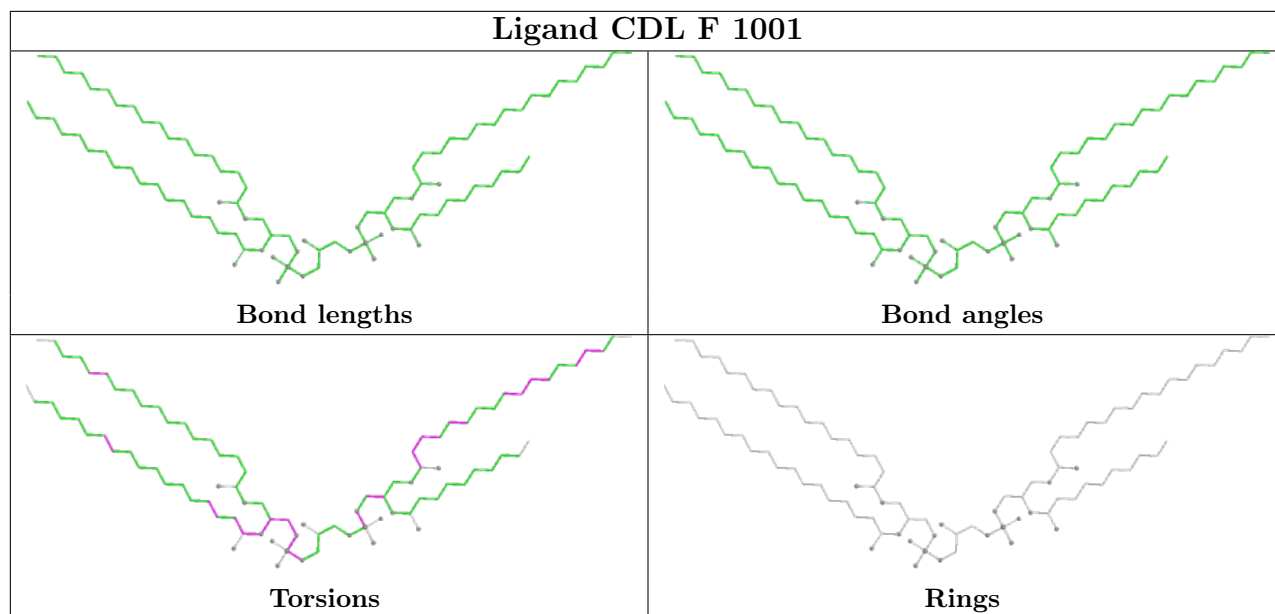




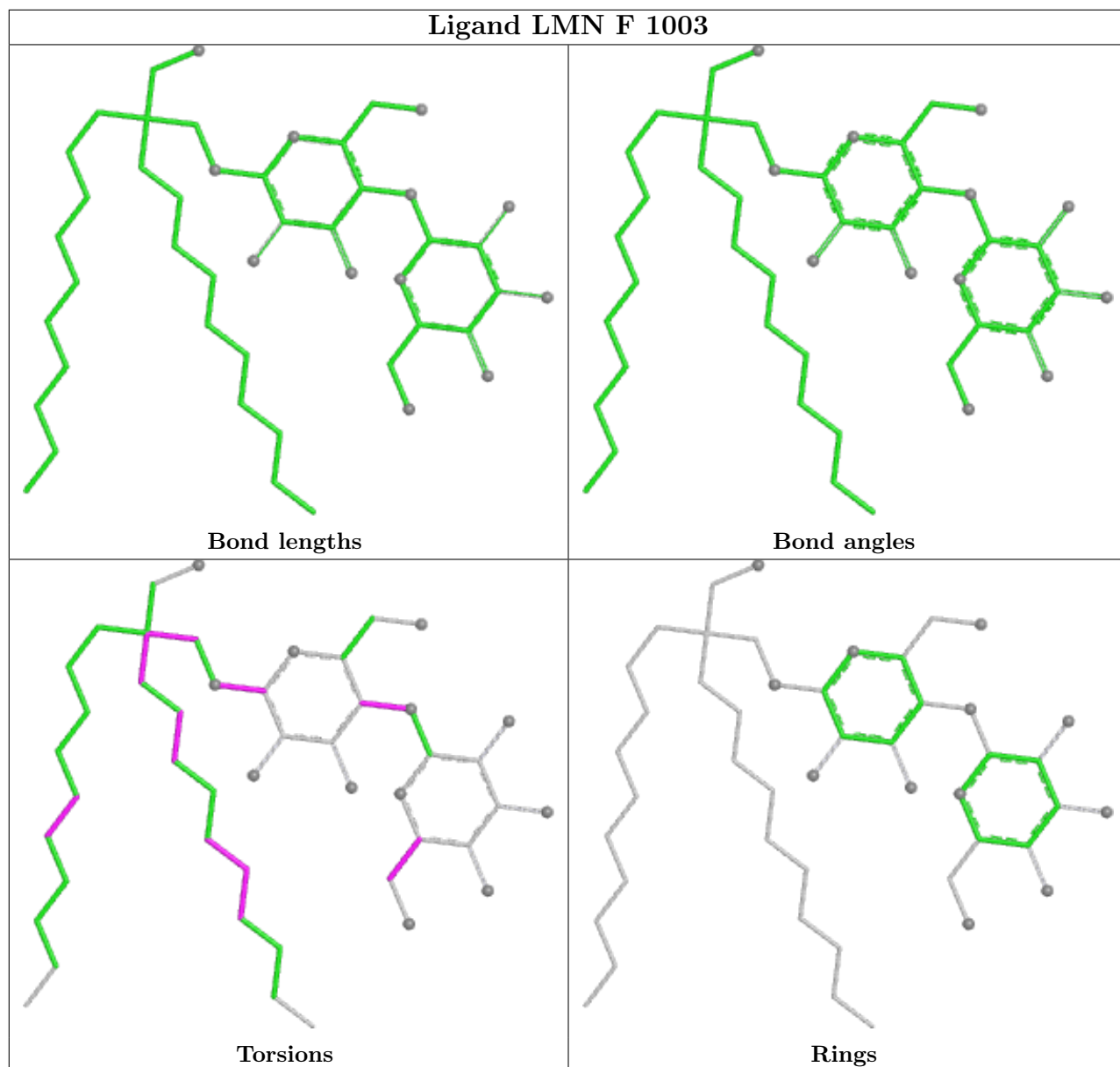


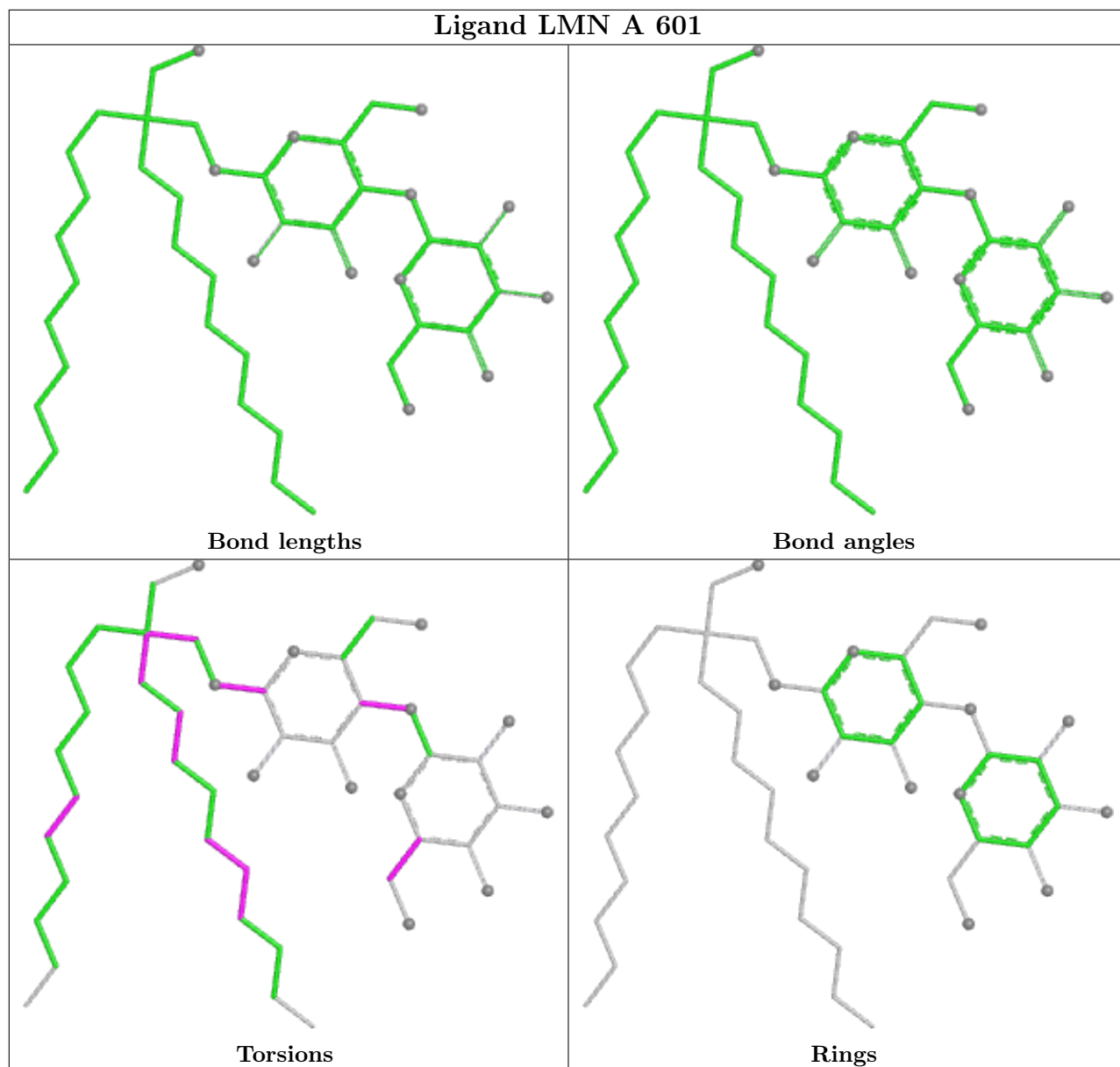


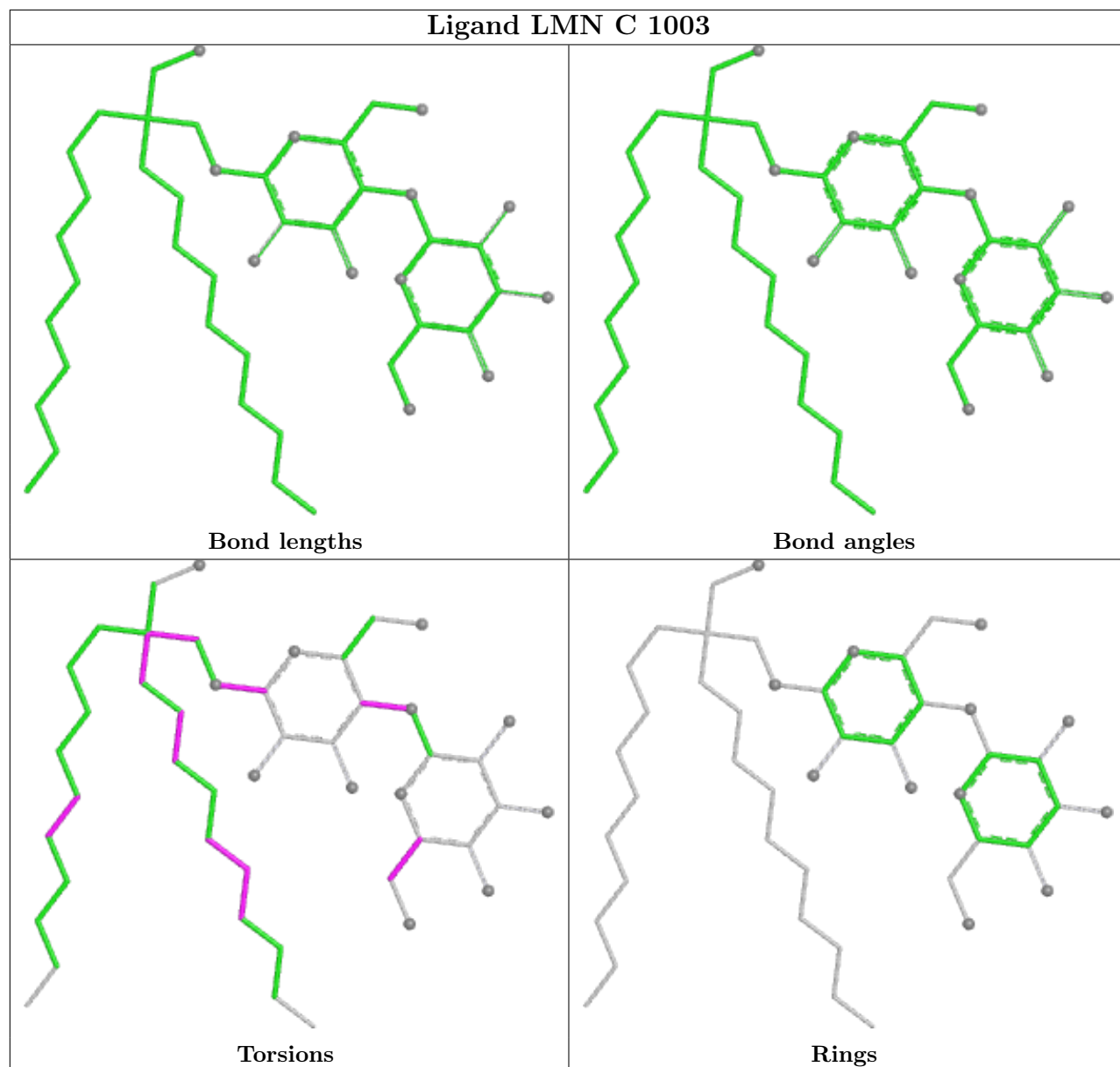


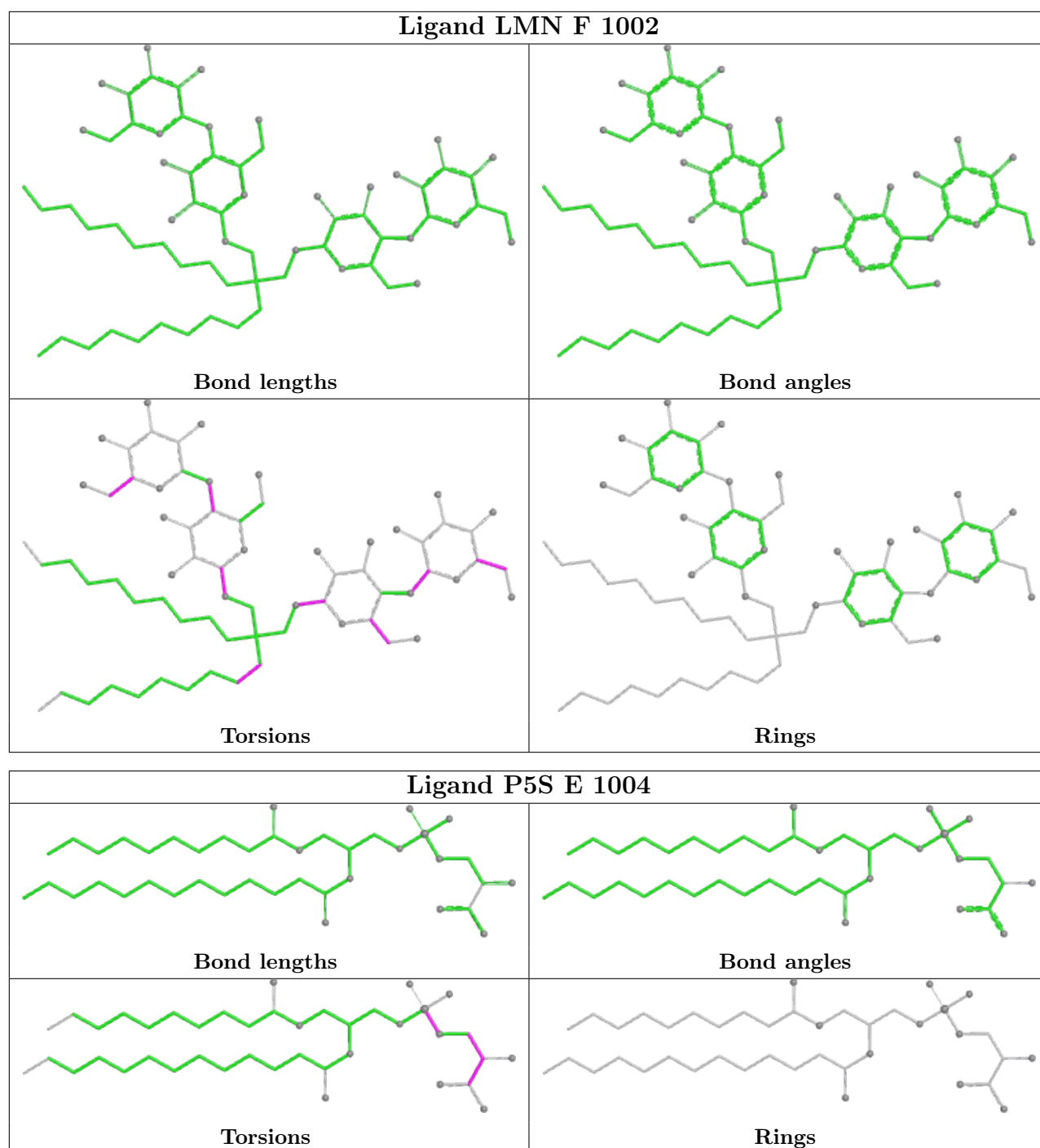


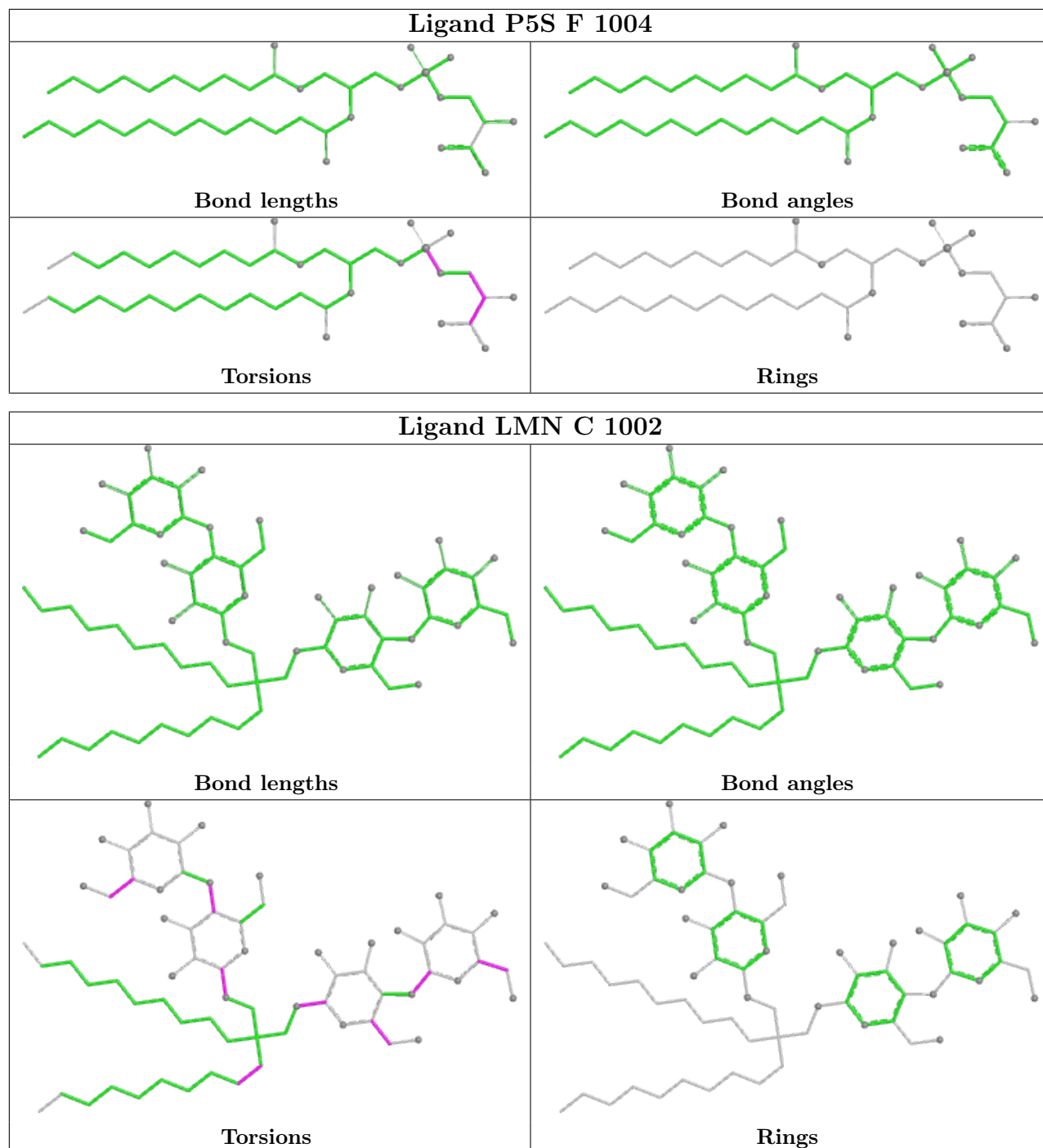


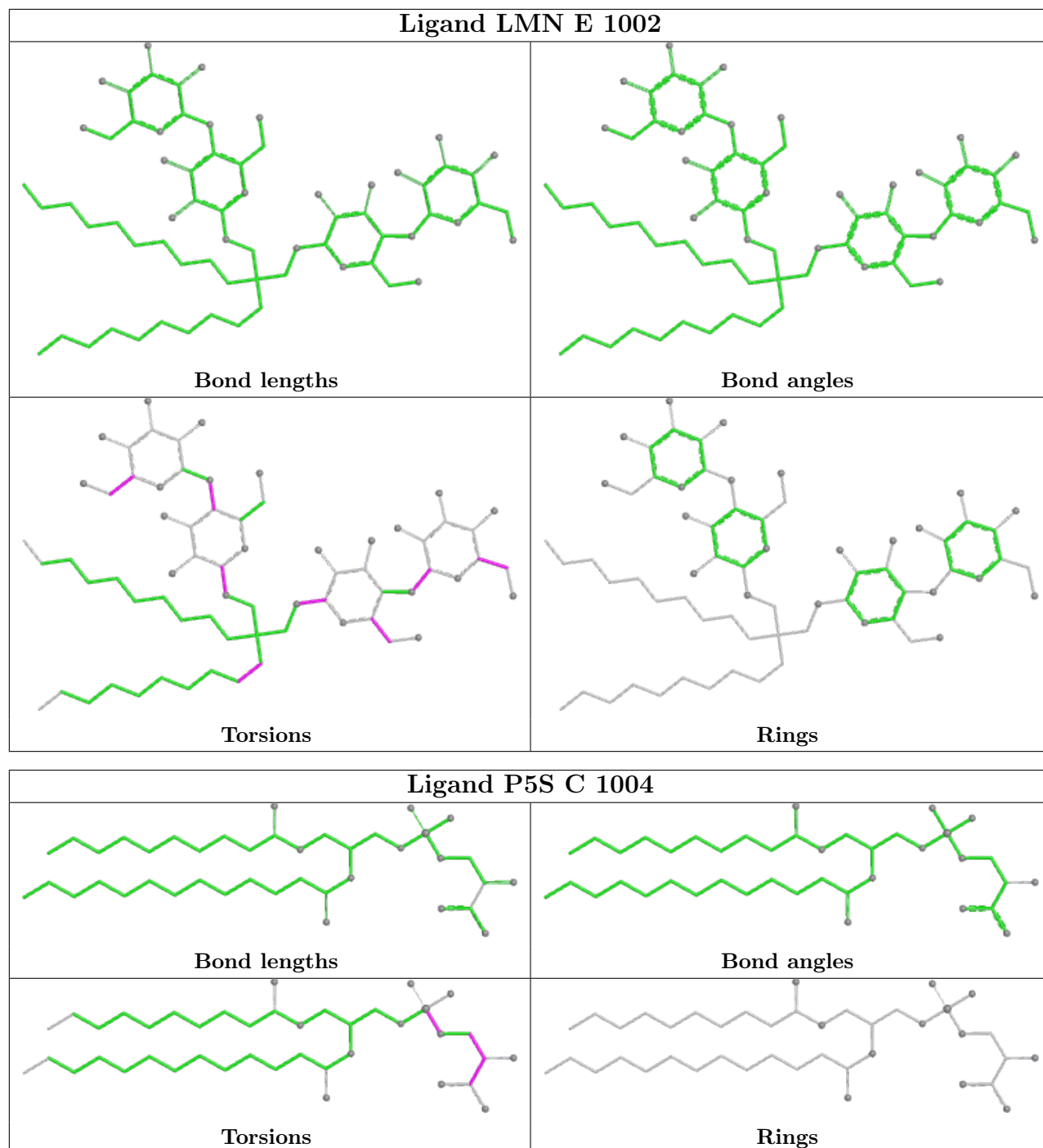


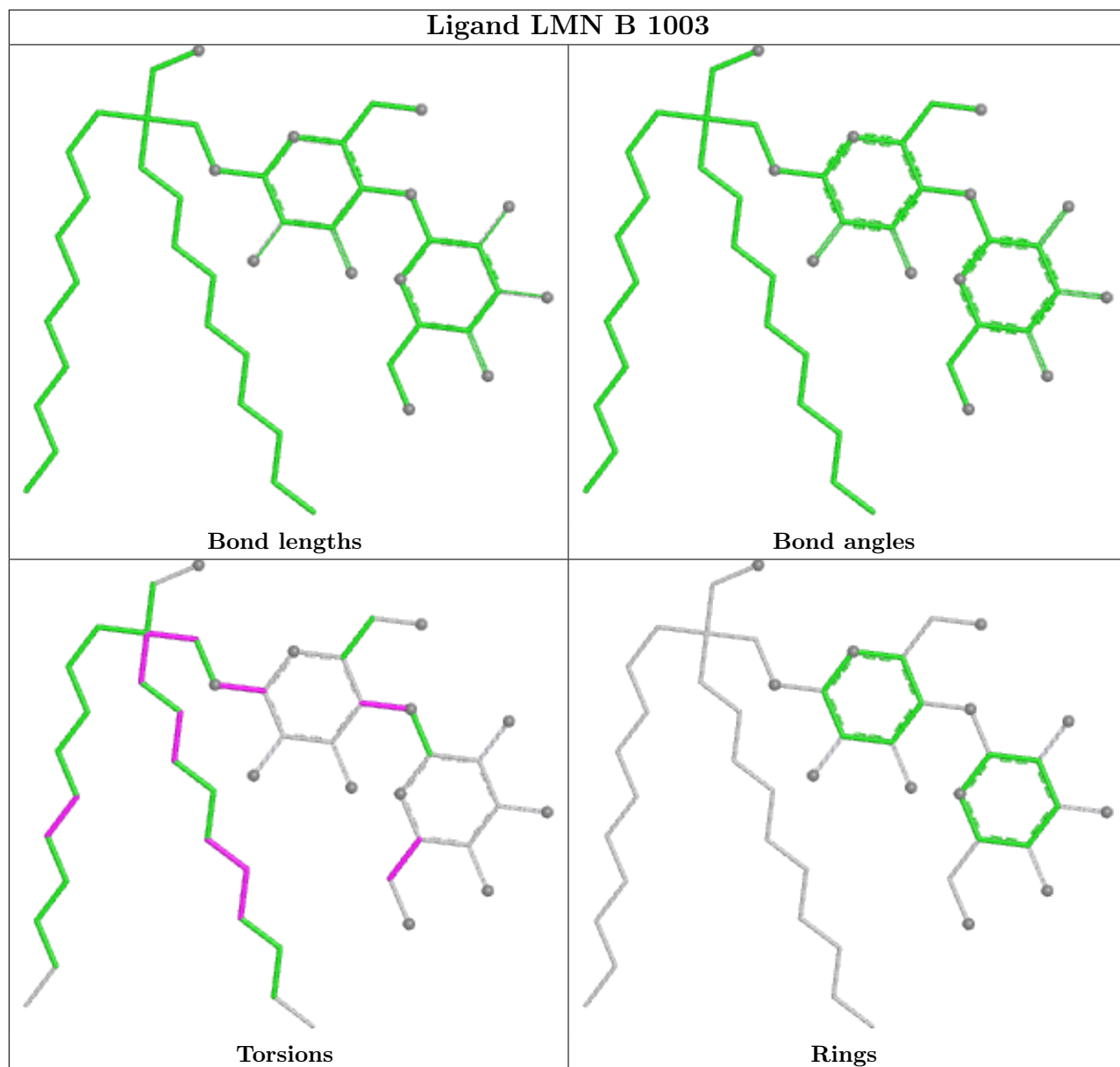


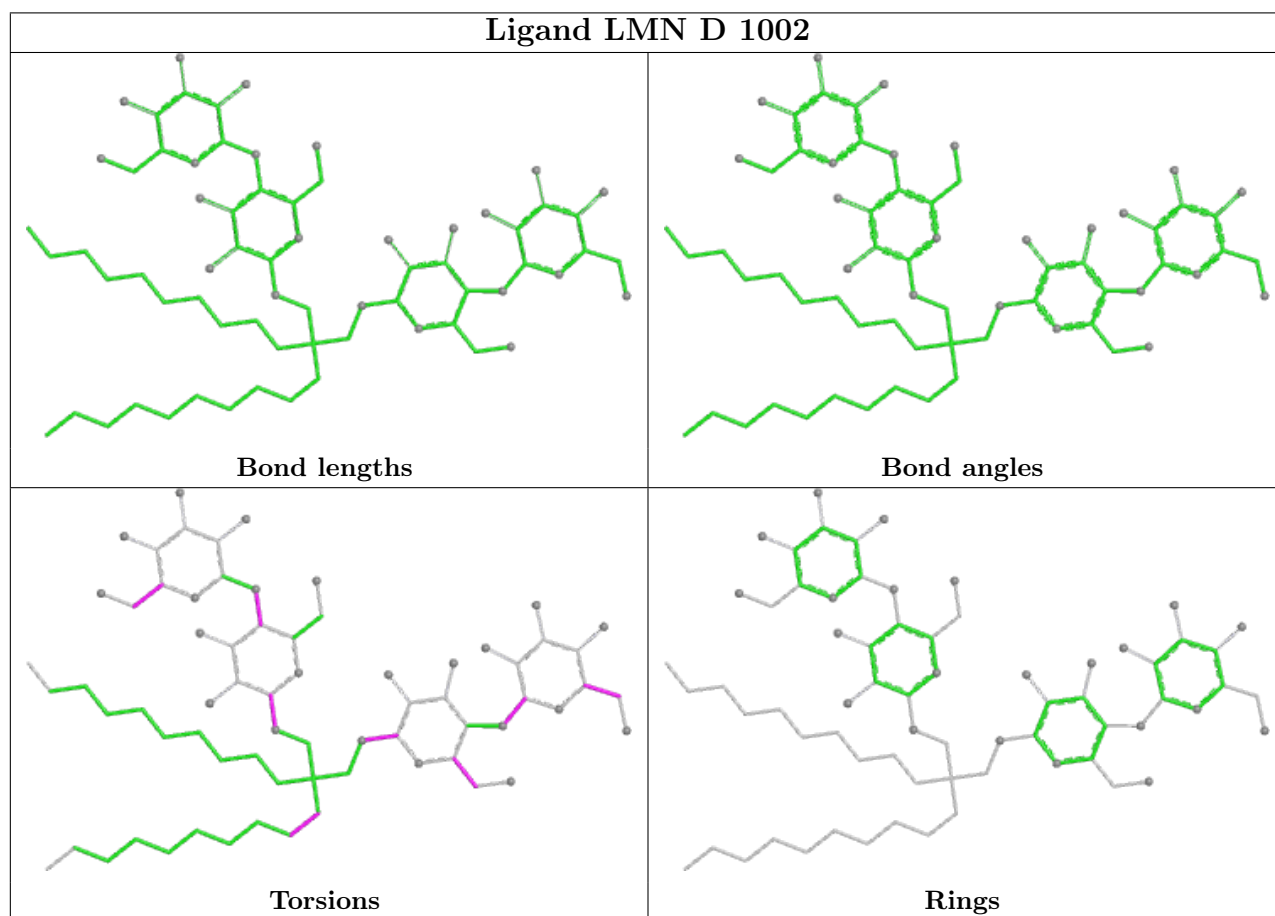
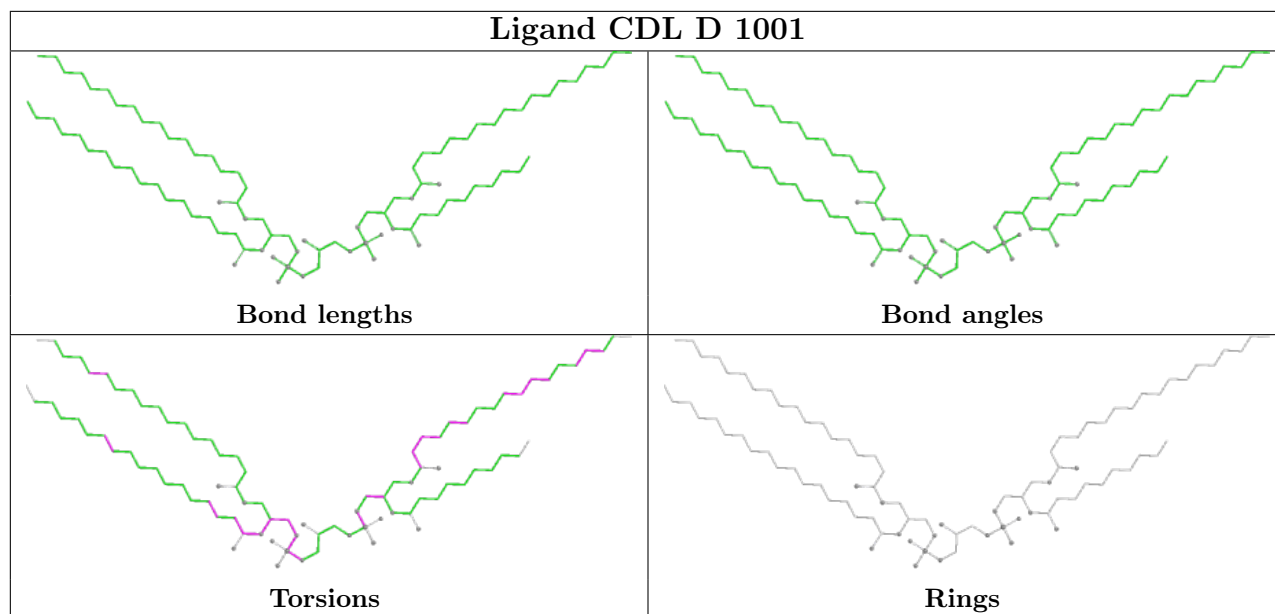




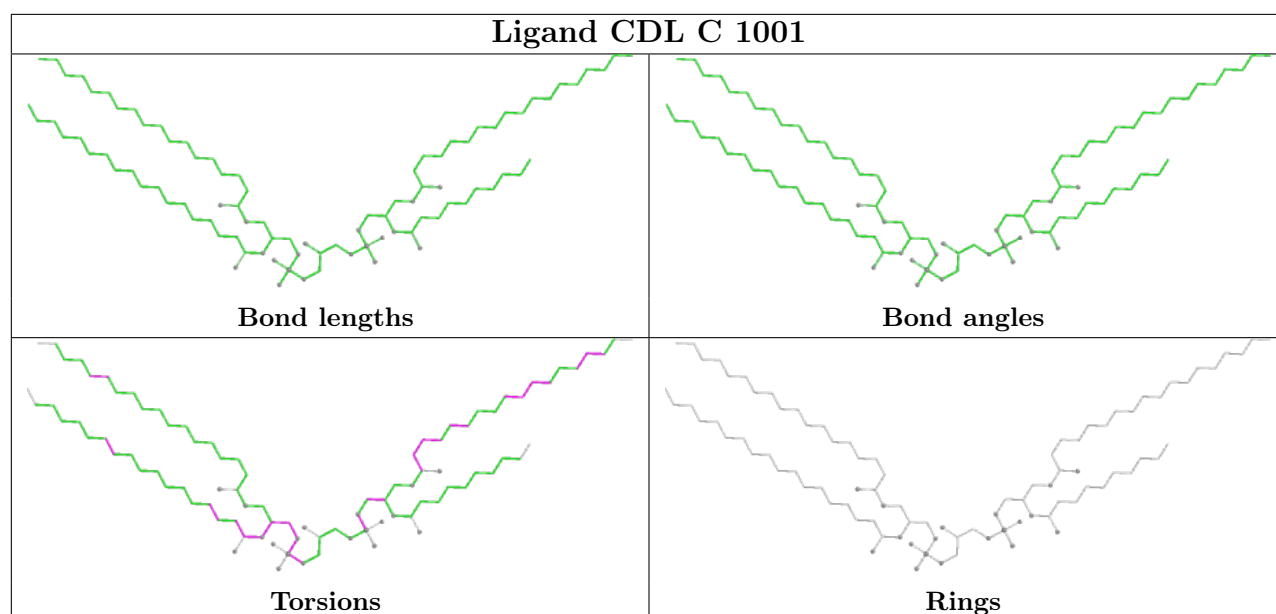












## 5.7 Other polymers [i](#)

There are no such residues in this entry.

## 5.8 Polymer linkage issues [i](#)

There are no chain breaks in this entry.

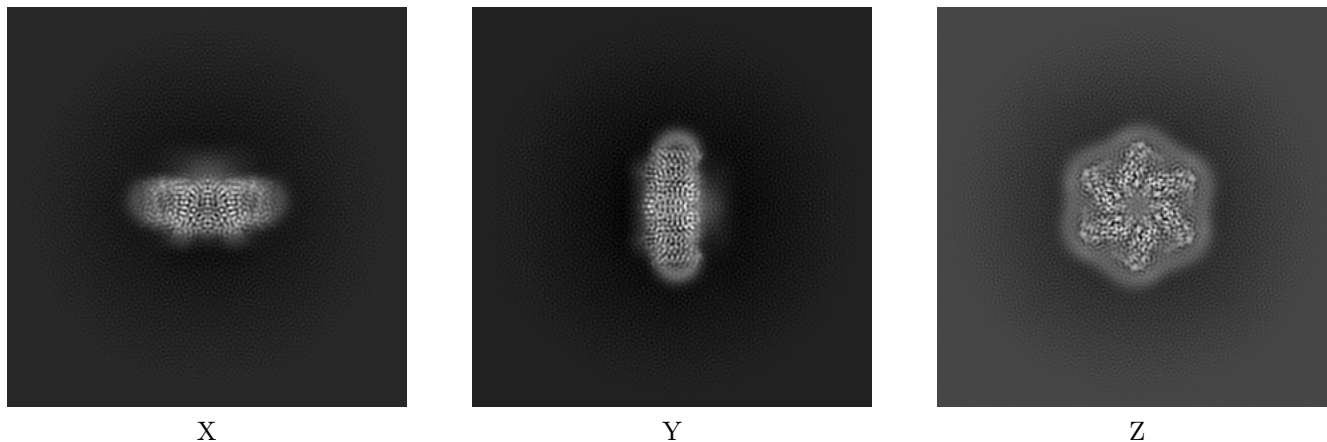
## 6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-10279. These allow visual inspection of the internal detail of the map and identification of artifacts.

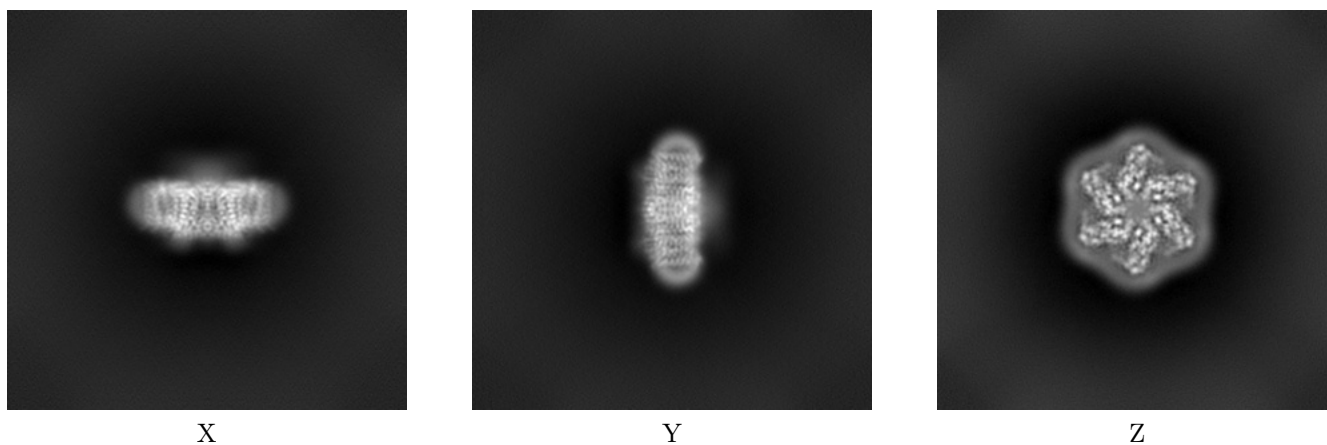
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

### 6.1 Orthogonal projections [i](#)

#### 6.1.1 Primary map



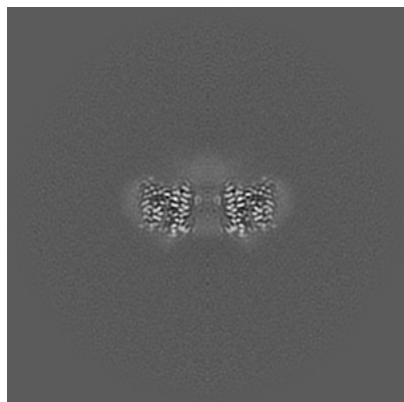
#### 6.1.2 Raw map



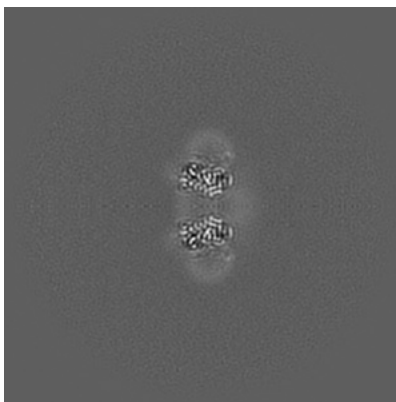
The images above show the map projected in three orthogonal directions.

## 6.2 Central slices [i](#)

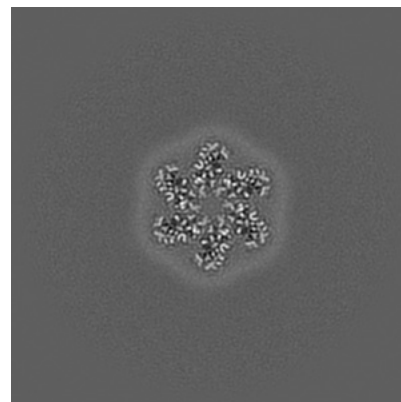
### 6.2.1 Primary map



X Index: 140

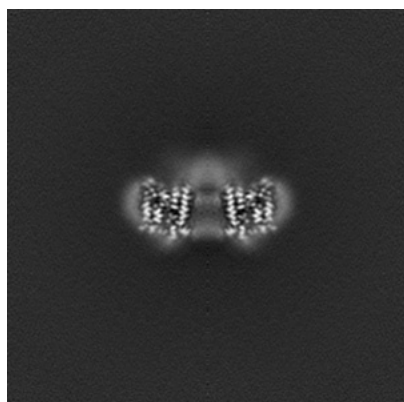


Y Index: 140

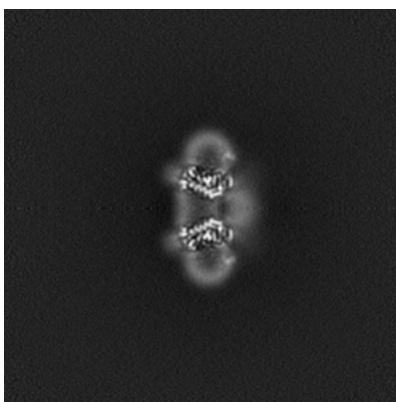


Z Index: 140

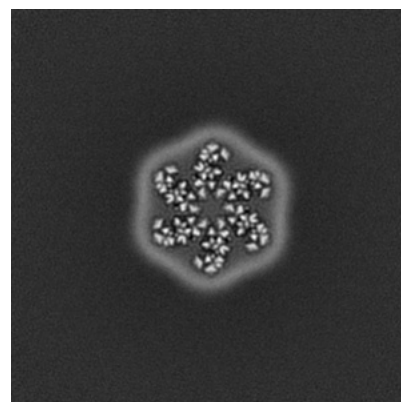
### 6.2.2 Raw map



X Index: 140



Y Index: 140

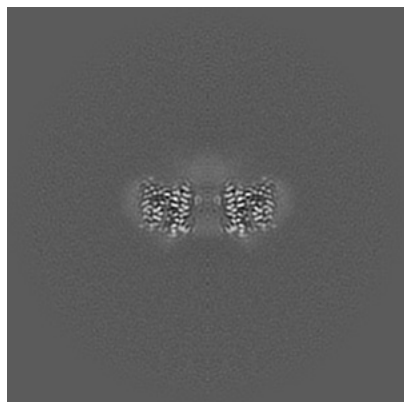


Z Index: 140

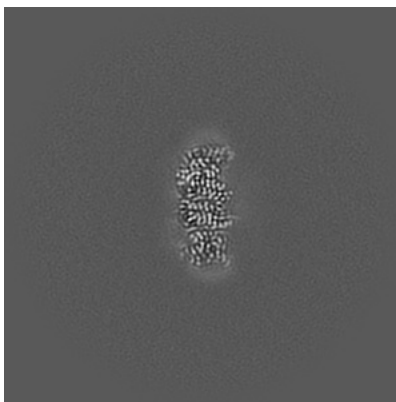
The images above show central slices of the map in three orthogonal directions.

## 6.3 Largest variance slices [i](#)

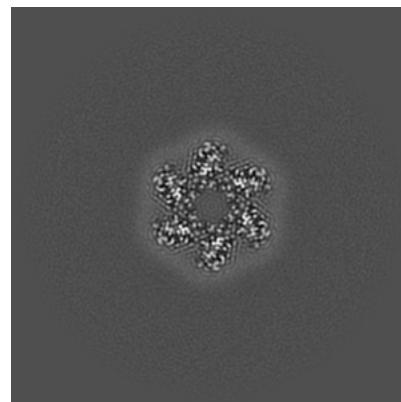
### 6.3.1 Primary map



X Index: 140

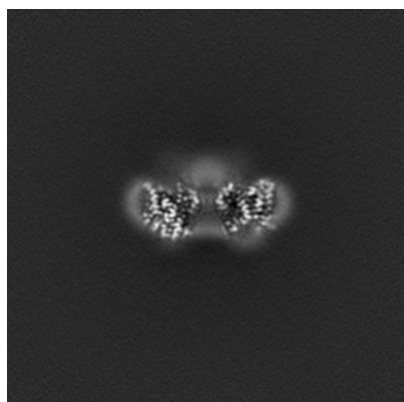


Y Index: 158

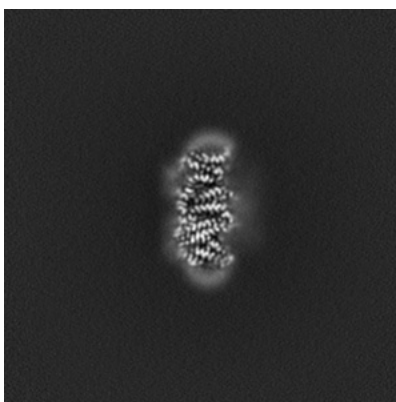


Z Index: 135

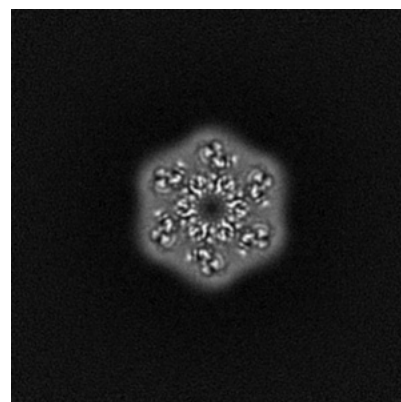
### 6.3.2 Raw map



X Index: 144



Y Index: 124

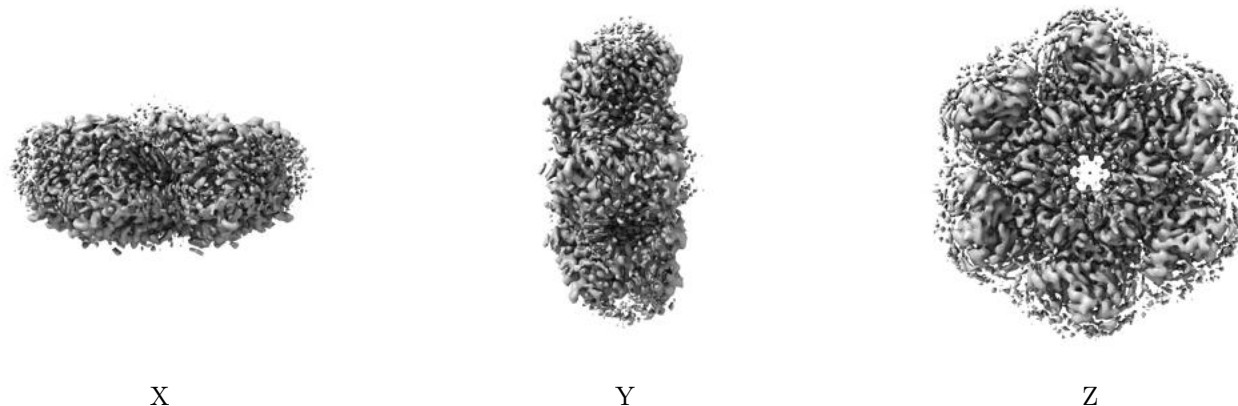


Z Index: 153

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

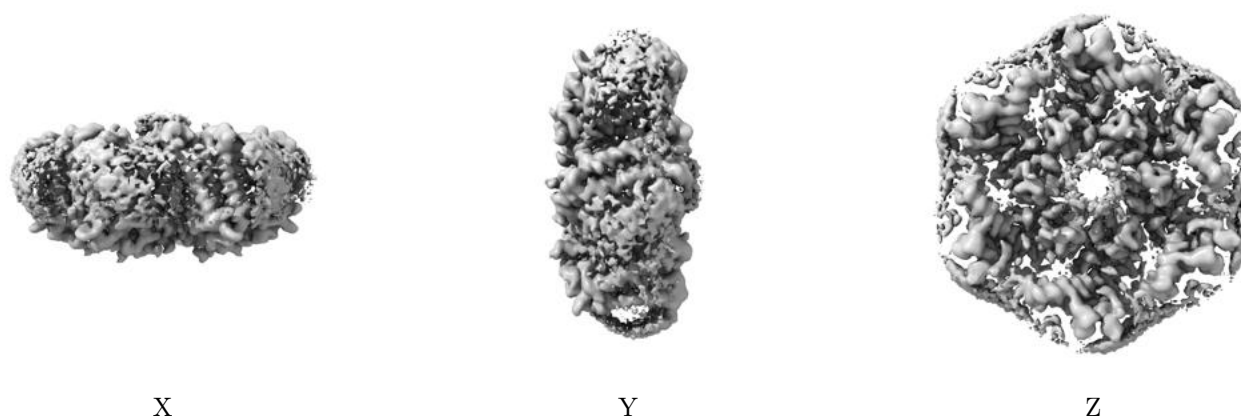
## 6.4 Orthogonal surface views [i](#)

### 6.4.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.7. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

### 6.4.2 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.

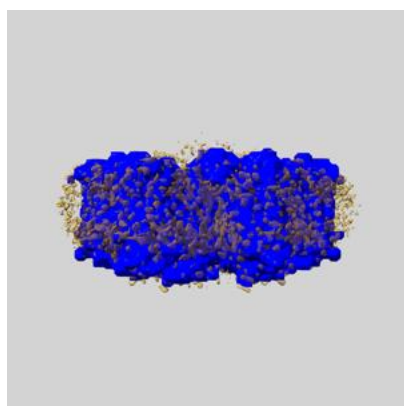
## 6.5 Mask visualisation [i](#)

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

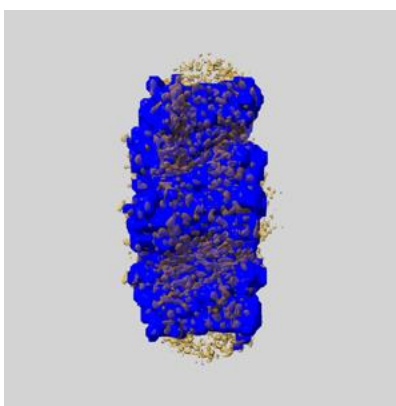
A mask typically either:

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

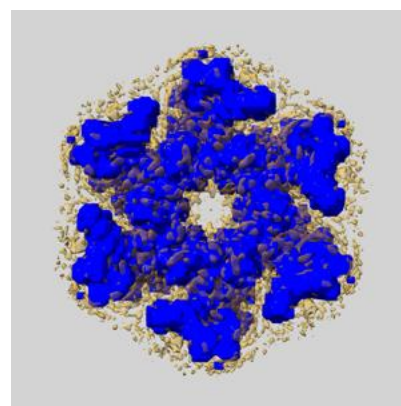
### 6.5.1 emd\_10279\_msk\_1.map [i](#)



X



Y

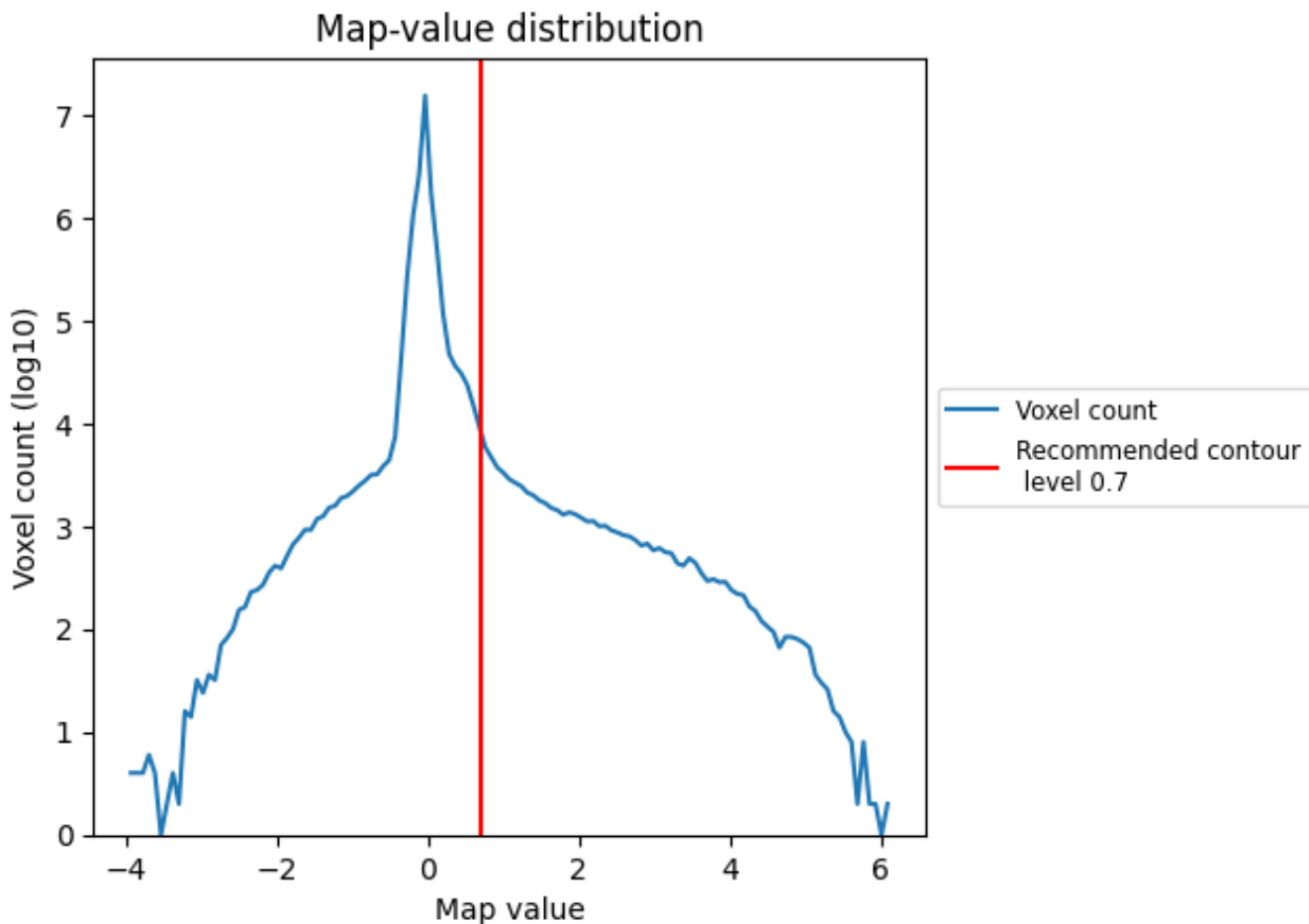


Z

## 7 Map analysis [i](#)

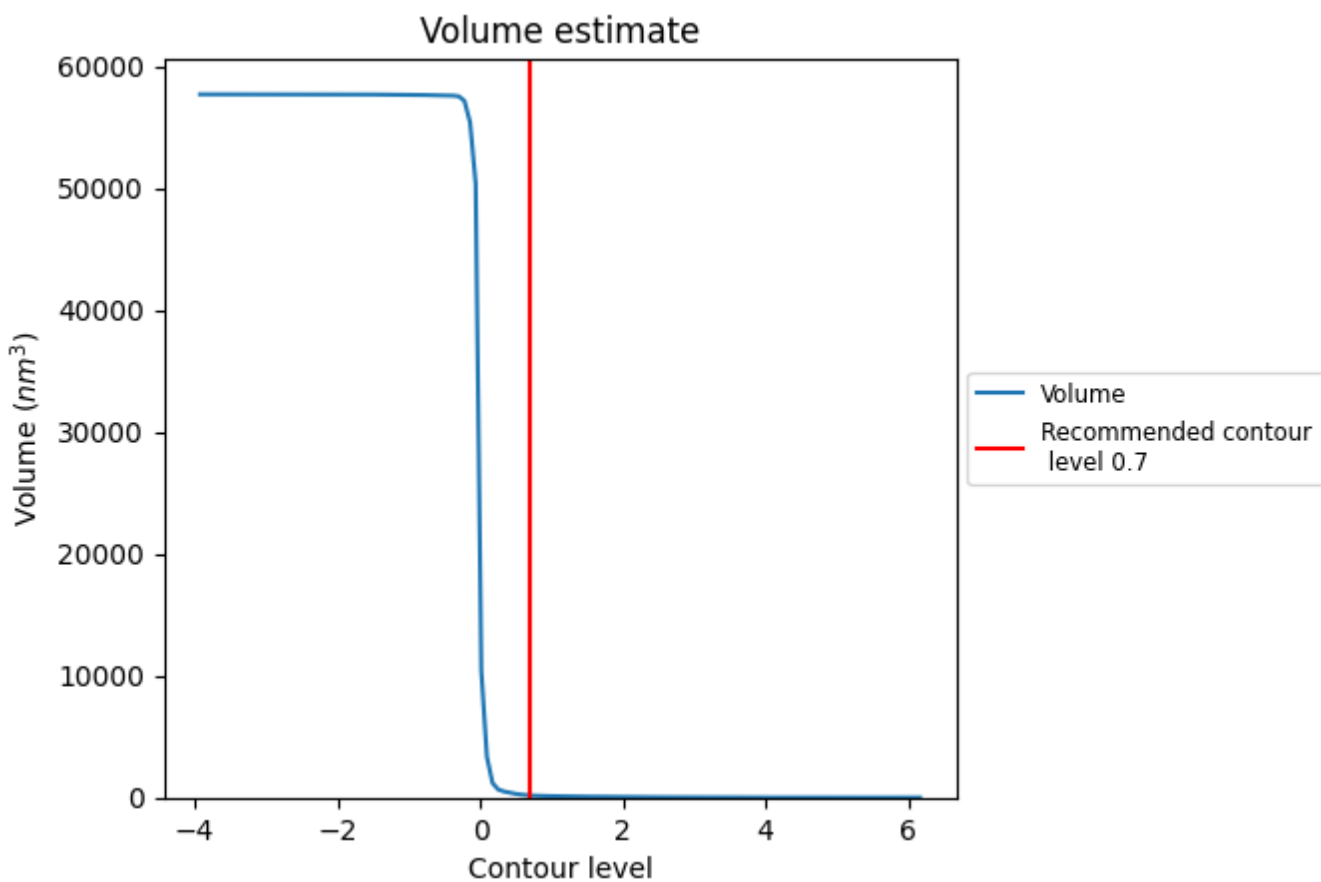
This section contains the results of statistical analysis of the map.

### 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](#)

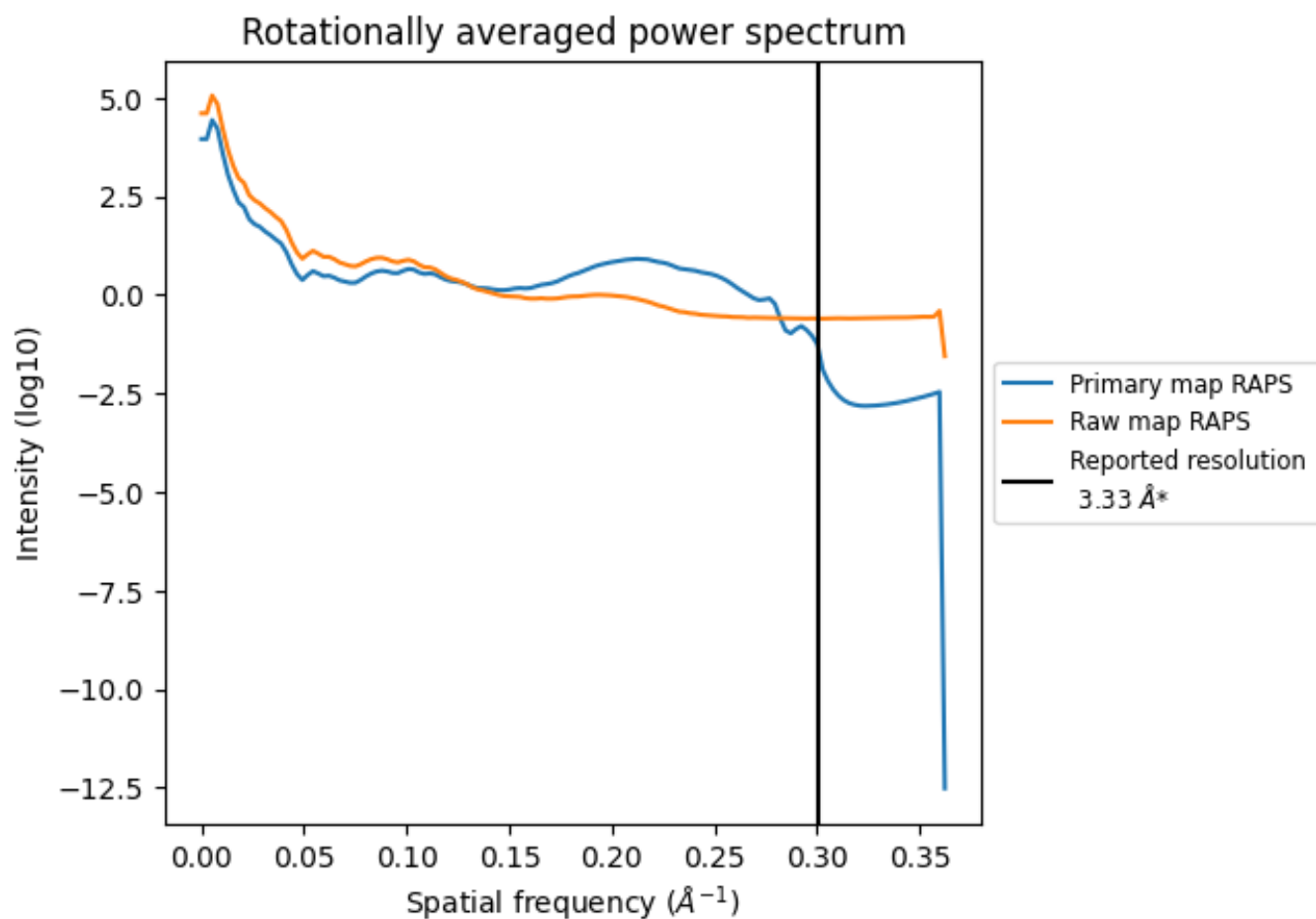


The volume at the recommended contour level is 173 nm<sup>3</sup>; this corresponds to an approximate mass of 157 kDa.

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



### 7.3 Rotationally averaged power spectrum i

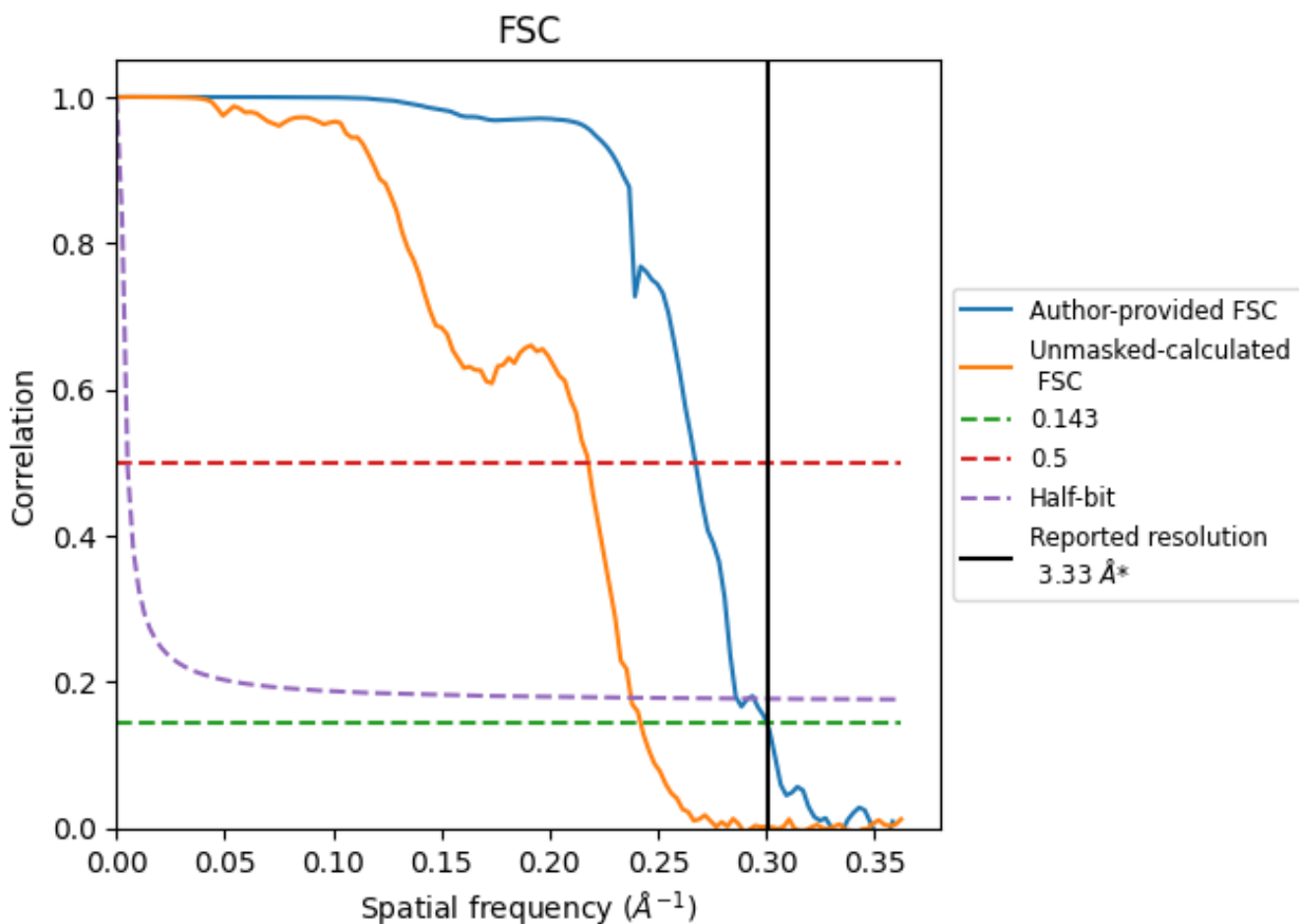


\*Reported resolution corresponds to spatial frequency of  $0.300 \text{ \AA}^{-1}$

## 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 0.300 Å<sup>-1</sup>

## 8.2 Resolution estimates [i](#)

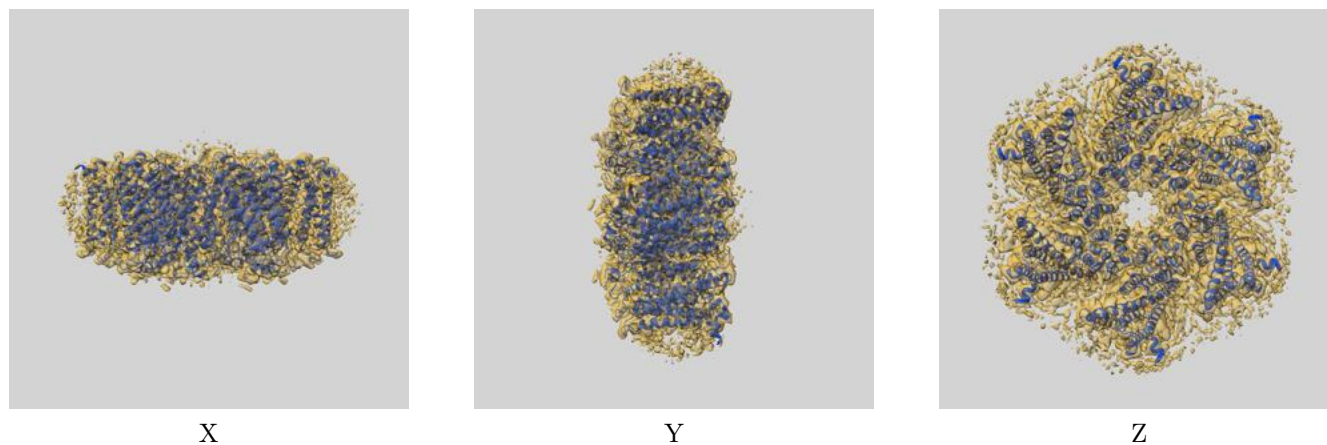
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.33	-	-
Author-provided FSC curve	3.33	3.74	3.49
Unmasked-calculated*	4.13	4.59	4.21

\*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 4.13 differs from the reported value 3.33 by more than 10 %

## 9 Map-model fit [i](#)

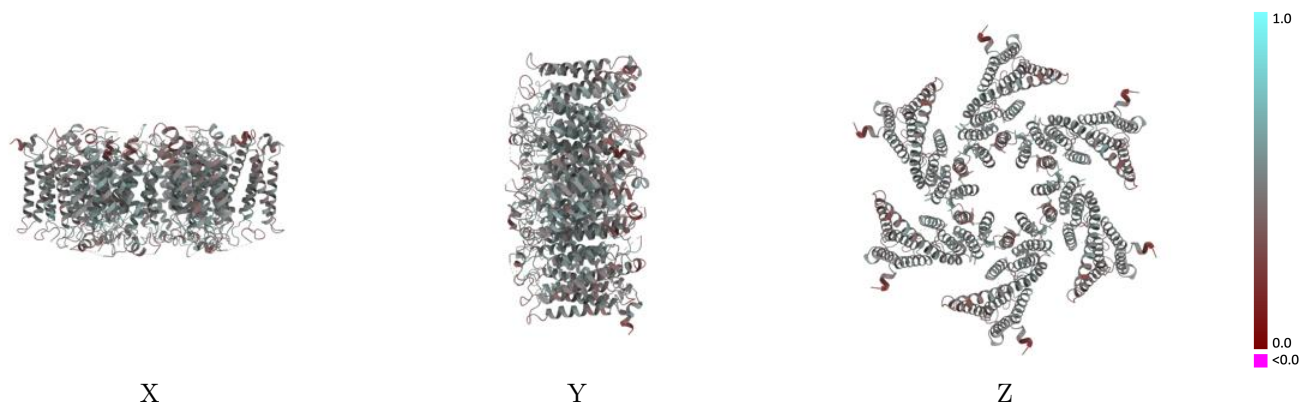
This section contains information regarding the fit between EMDB map EMD-10279 and PDB model 6SP2. Per-residue inclusion information can be found in section 3 on page 13.

### 9.1 Map-model overlay [i](#)



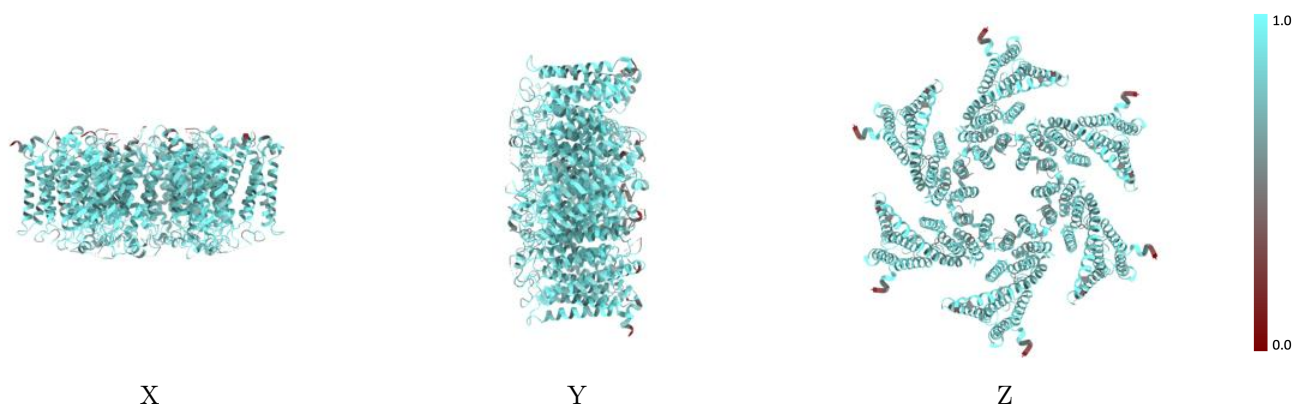
The images above show the 3D surface view of the map at the recommended contour level 0.7 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



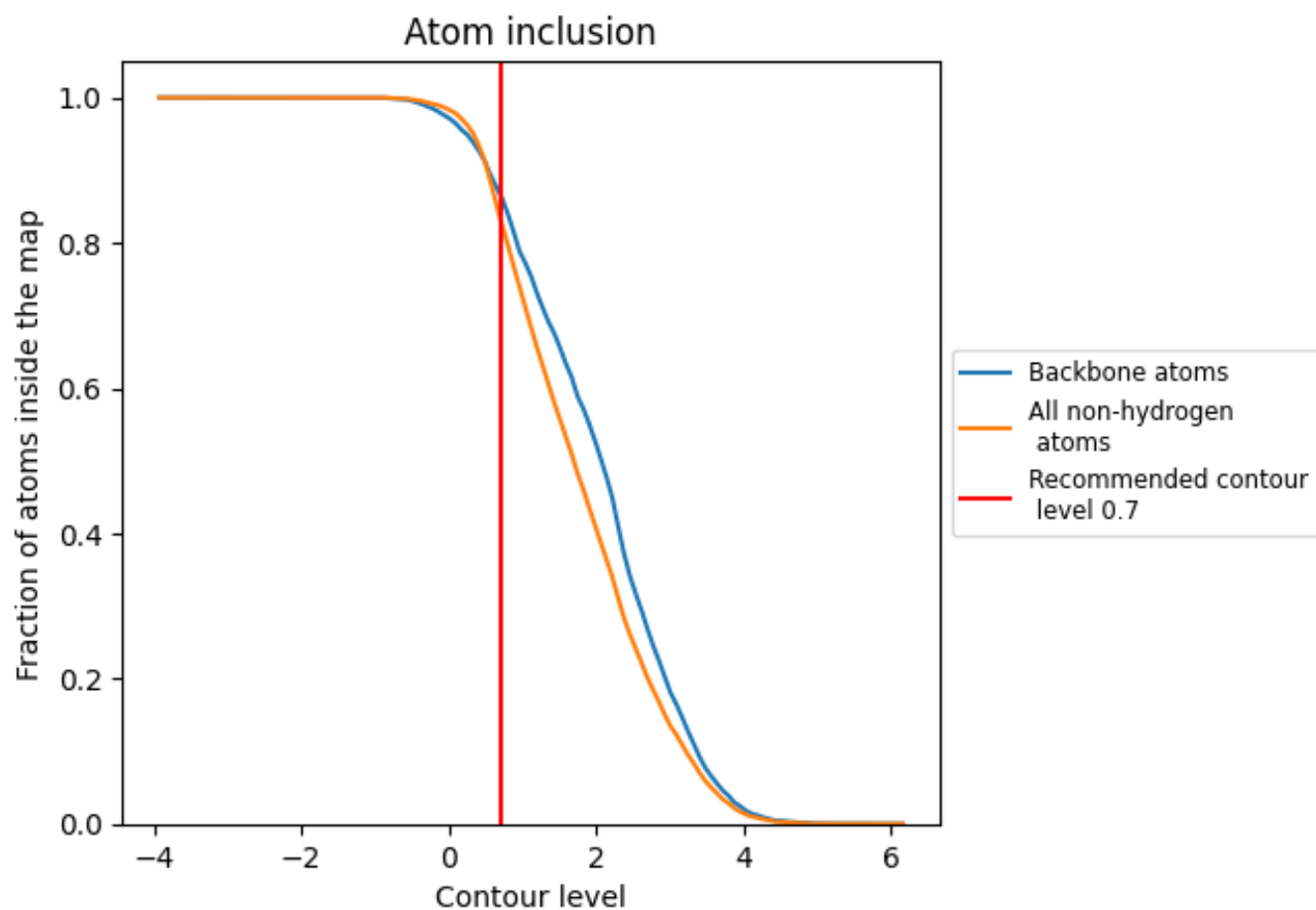
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.7).















## 9.4 Atom inclusion [i](#)



At the recommended contour level, 87% of all backbone atoms, 83% of all non-hydrogen atoms, are inside the map.

## 9.5 Map-model fit summary [i](#)

The table lists the average atom inclusion at the recommended contour level (0.7) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8325	 0.4820
A	 0.8331	 0.4860
B	 0.8311	 0.4830
C	 0.8341	 0.4800
D	 0.8341	 0.4830
E	 0.8295	 0.4800
F	 0.8331	 0.4800

