



wwPDB EM Validation Summary Report ⓘ

Mar 19, 2026 – 10:49 PM UTC

PDB ID : 7XNH / pdb_00007xnh
EMDB ID : EMD-33315
Title : Human Cx36/GJD2 gap junction channel with pore-lining N-terminal helices
in soybean lipids
Authors : Lee, S.N.; Cho, H.J.; Jeong, H.; Ryu, B.; Lee, H.J.; Lee, H.H.; Woo, J.S.
Deposited on : 2022-04-28
Resolution : 3.10 Å(reported)

This is a wwPDB EM Validation Summary Report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>
with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

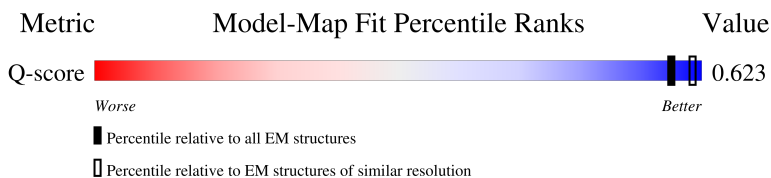
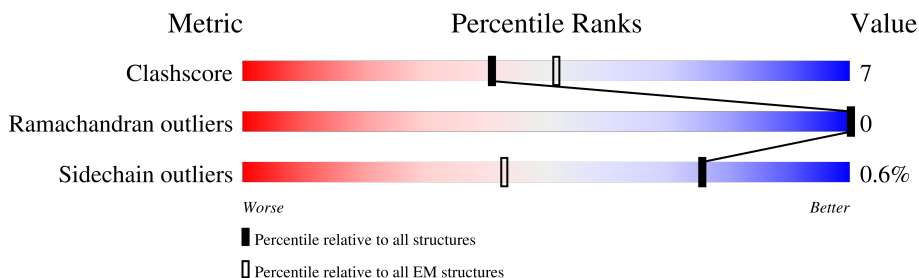
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data in the EMDB archive up until May 2025)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:
ELECTRON MICROSCOPY

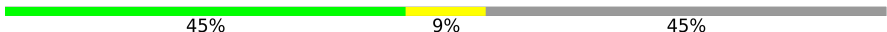
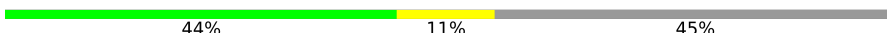
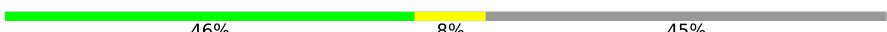
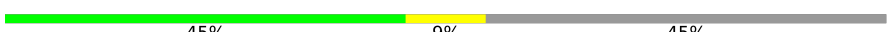
The reported resolution of this entry is 3.10 Å.

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



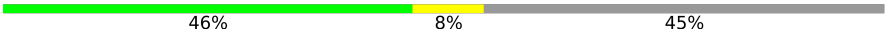
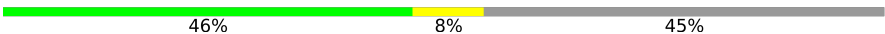
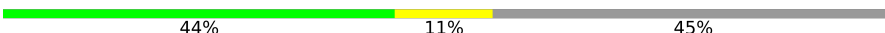
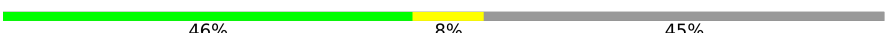
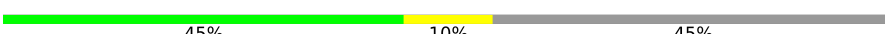
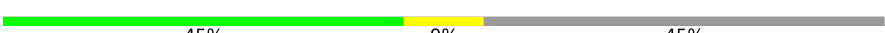
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	14724 (2.60 - 3.60)

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

Mol	Chain	Length	Quality of chain
1	A	332	
1	B	332	
1	C	332	
1	D	332	

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Mol	Chain	Length	Quality of chain
1	E	332	 46%8%45%
1	F	332	 46%8%45%
1	G	332	 44%11%45%
1	H	332	 46%8%45%
1	I	332	 45%10%45%
1	J	332	 45%9%45%
1	K	332	 44%10%45%
1	L	332	 44%10%45%

2 Entry composition [i](#)

There are 2 unique types of molecules in this entry. The entry contains 19224 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 Gap junction delta-2 protein.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	B	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	C	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	D	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	E	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	F	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	G	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	H	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	I	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	J	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	K	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		
1	L	181	Total	C	N	O	S	0	0
			1458	957	235	253	13		

There are 132 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	322	SER	-	expression tag	UNP Q9UKL4
A	323	ARG	-	expression tag	UNP Q9UKL4
A	324	GLY	-	expression tag	UNP Q9UKL4
A	325	ASP	-	expression tag	UNP Q9UKL4
A	326	MET	-	expression tag	UNP Q9UKL4
A	327	LEU	-	expression tag	UNP Q9UKL4

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Chain	Residue	Modelled	Actual	Comment	Reference
A	328	GLU	-	expression tag	UNP Q9UKL4
A	329	VAL	-	expression tag	UNP Q9UKL4
A	330	LEU	-	expression tag	UNP Q9UKL4
A	331	PHE	-	expression tag	UNP Q9UKL4
A	332	GLN	-	expression tag	UNP Q9UKL4
B	322	SER	-	expression tag	UNP Q9UKL4
B	323	ARG	-	expression tag	UNP Q9UKL4
B	324	GLY	-	expression tag	UNP Q9UKL4
B	325	ASP	-	expression tag	UNP Q9UKL4
B	326	MET	-	expression tag	UNP Q9UKL4
B	327	LEU	-	expression tag	UNP Q9UKL4
B	328	GLU	-	expression tag	UNP Q9UKL4
B	329	VAL	-	expression tag	UNP Q9UKL4
B	330	LEU	-	expression tag	UNP Q9UKL4
B	331	PHE	-	expression tag	UNP Q9UKL4
B	332	GLN	-	expression tag	UNP Q9UKL4
C	322	SER	-	expression tag	UNP Q9UKL4
C	323	ARG	-	expression tag	UNP Q9UKL4
C	324	GLY	-	expression tag	UNP Q9UKL4
C	325	ASP	-	expression tag	UNP Q9UKL4
C	326	MET	-	expression tag	UNP Q9UKL4
C	327	LEU	-	expression tag	UNP Q9UKL4
C	328	GLU	-	expression tag	UNP Q9UKL4
C	329	VAL	-	expression tag	UNP Q9UKL4
C	330	LEU	-	expression tag	UNP Q9UKL4
C	331	PHE	-	expression tag	UNP Q9UKL4
C	332	GLN	-	expression tag	UNP Q9UKL4
D	322	SER	-	expression tag	UNP Q9UKL4
D	323	ARG	-	expression tag	UNP Q9UKL4
D	324	GLY	-	expression tag	UNP Q9UKL4
D	325	ASP	-	expression tag	UNP Q9UKL4
D	326	MET	-	expression tag	UNP Q9UKL4
D	327	LEU	-	expression tag	UNP Q9UKL4
D	328	GLU	-	expression tag	UNP Q9UKL4
D	329	VAL	-	expression tag	UNP Q9UKL4
D	330	LEU	-	expression tag	UNP Q9UKL4
D	331	PHE	-	expression tag	UNP Q9UKL4
D	332	GLN	-	expression tag	UNP Q9UKL4
E	322	SER	-	expression tag	UNP Q9UKL4
E	323	ARG	-	expression tag	UNP Q9UKL4
E	324	GLY	-	expression tag	UNP Q9UKL4
E	325	ASP	-	expression tag	UNP Q9UKL4

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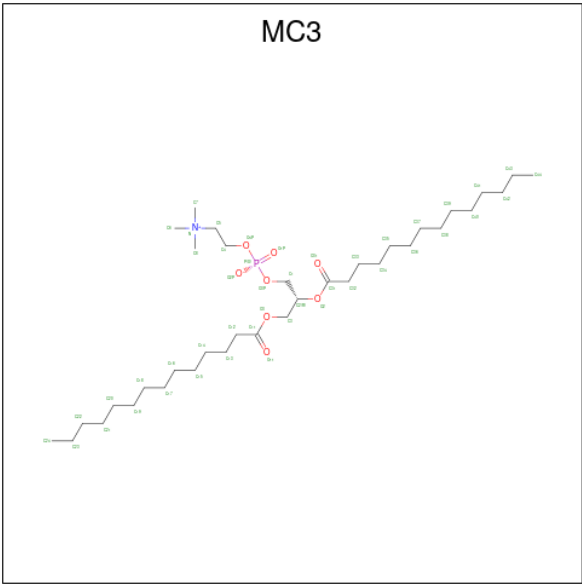
Chain	Residue	Modelled	Actual	Comment	Reference
E	326	MET	-	expression tag	UNP Q9UKL4
E	327	LEU	-	expression tag	UNP Q9UKL4
E	328	GLU	-	expression tag	UNP Q9UKL4
E	329	VAL	-	expression tag	UNP Q9UKL4
E	330	LEU	-	expression tag	UNP Q9UKL4
E	331	PHE	-	expression tag	UNP Q9UKL4
E	332	GLN	-	expression tag	UNP Q9UKL4
F	322	SER	-	expression tag	UNP Q9UKL4
F	323	ARG	-	expression tag	UNP Q9UKL4
F	324	GLY	-	expression tag	UNP Q9UKL4
F	325	ASP	-	expression tag	UNP Q9UKL4
F	326	MET	-	expression tag	UNP Q9UKL4
F	327	LEU	-	expression tag	UNP Q9UKL4
F	328	GLU	-	expression tag	UNP Q9UKL4
F	329	VAL	-	expression tag	UNP Q9UKL4
F	330	LEU	-	expression tag	UNP Q9UKL4
F	331	PHE	-	expression tag	UNP Q9UKL4
F	332	GLN	-	expression tag	UNP Q9UKL4
G	322	SER	-	expression tag	UNP Q9UKL4
G	323	ARG	-	expression tag	UNP Q9UKL4
G	324	GLY	-	expression tag	UNP Q9UKL4
G	325	ASP	-	expression tag	UNP Q9UKL4
G	326	MET	-	expression tag	UNP Q9UKL4
G	327	LEU	-	expression tag	UNP Q9UKL4
G	328	GLU	-	expression tag	UNP Q9UKL4
G	329	VAL	-	expression tag	UNP Q9UKL4
G	330	LEU	-	expression tag	UNP Q9UKL4
G	331	PHE	-	expression tag	UNP Q9UKL4
G	332	GLN	-	expression tag	UNP Q9UKL4
H	322	SER	-	expression tag	UNP Q9UKL4
H	323	ARG	-	expression tag	UNP Q9UKL4
H	324	GLY	-	expression tag	UNP Q9UKL4
H	325	ASP	-	expression tag	UNP Q9UKL4
H	326	MET	-	expression tag	UNP Q9UKL4
H	327	LEU	-	expression tag	UNP Q9UKL4
H	328	GLU	-	expression tag	UNP Q9UKL4
H	329	VAL	-	expression tag	UNP Q9UKL4
H	330	LEU	-	expression tag	UNP Q9UKL4
H	331	PHE	-	expression tag	UNP Q9UKL4
H	332	GLN	-	expression tag	UNP Q9UKL4
I	322	SER	-	expression tag	UNP Q9UKL4
I	323	ARG	-	expression tag	UNP Q9UKL4

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Chain	Residue	Modelled	Actual	Comment	Reference
I	324	GLY	-	expression tag	UNP Q9UKL4
I	325	ASP	-	expression tag	UNP Q9UKL4
I	326	MET	-	expression tag	UNP Q9UKL4
I	327	LEU	-	expression tag	UNP Q9UKL4
I	328	GLU	-	expression tag	UNP Q9UKL4
I	329	VAL	-	expression tag	UNP Q9UKL4
I	330	LEU	-	expression tag	UNP Q9UKL4
I	331	PHE	-	expression tag	UNP Q9UKL4
I	332	GLN	-	expression tag	UNP Q9UKL4
J	322	SER	-	expression tag	UNP Q9UKL4
J	323	ARG	-	expression tag	UNP Q9UKL4
J	324	GLY	-	expression tag	UNP Q9UKL4
J	325	ASP	-	expression tag	UNP Q9UKL4
J	326	MET	-	expression tag	UNP Q9UKL4
J	327	LEU	-	expression tag	UNP Q9UKL4
J	328	GLU	-	expression tag	UNP Q9UKL4
J	329	VAL	-	expression tag	UNP Q9UKL4
J	330	LEU	-	expression tag	UNP Q9UKL4
J	331	PHE	-	expression tag	UNP Q9UKL4
J	332	GLN	-	expression tag	UNP Q9UKL4
K	322	SER	-	expression tag	UNP Q9UKL4
K	323	ARG	-	expression tag	UNP Q9UKL4
K	324	GLY	-	expression tag	UNP Q9UKL4
K	325	ASP	-	expression tag	UNP Q9UKL4
K	326	MET	-	expression tag	UNP Q9UKL4
K	327	LEU	-	expression tag	UNP Q9UKL4
K	328	GLU	-	expression tag	UNP Q9UKL4
K	329	VAL	-	expression tag	UNP Q9UKL4
K	330	LEU	-	expression tag	UNP Q9UKL4
K	331	PHE	-	expression tag	UNP Q9UKL4
K	332	GLN	-	expression tag	UNP Q9UKL4
L	322	SER	-	expression tag	UNP Q9UKL4
L	323	ARG	-	expression tag	UNP Q9UKL4
L	324	GLY	-	expression tag	UNP Q9UKL4
L	325	ASP	-	expression tag	UNP Q9UKL4
L	326	MET	-	expression tag	UNP Q9UKL4
L	327	LEU	-	expression tag	UNP Q9UKL4
L	328	GLU	-	expression tag	UNP Q9UKL4
L	329	VAL	-	expression tag	UNP Q9UKL4
L	330	LEU	-	expression tag	UNP Q9UKL4
L	331	PHE	-	expression tag	UNP Q9UKL4
L	332	GLN	-	expression tag	UNP Q9UKL4

- Molecule 2 is 1,2-DIMYRISTOYL-RAC-GLYCERO-3-PHOSPHOCHOLINE (CCD ID: MC3) (formula: C₃₆H₇₂NO₈P) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms		AltConf
2	A	1	Total	C	0
			14	14	
2	A	1	Total	C	0
			14	14	
2	A	1	Total	C	0
			14	14	
2	A	1	Total	C	0
			10	10	
2	A	1	Total	C	0
			10	10	
2	A	1	Total	C	0
			14	14	
2	A	1	Total	C	0
			14	14	
2	A	1	Total	C	0
			14	14	
2	A	1	Total	C	0
			14	14	
2	A	1	Total	C	0
			14	14	
2	A	1	Total	C	0
			12	12	

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Mol	Chain	Residues	Atoms	AltConf
2	B	1	Total C 14 14	0
2	B	1	Total C 14 14	0
2	B	1	Total C 10 10	0
2	B	1	Total C 10 10	0
2	B	1	Total C 14 14	0
2	B	1	Total C 14 14	0
2	B	1	Total C 14 14	0
2	B	1	Total C 14 14	0
2	B	1	Total C 14 14	0
2	B	1	Total C 12 12	0
2	C	1	Total C 14 14	0
2	C	1	Total C 14 14	0
2	C	1	Total C 10 10	0
2	C	1	Total C 10 10	0
2	C	1	Total C 14 14	0
2	C	1	Total C 14 14	0
2	C	1	Total C 14 14	0
2	C	1	Total C 14 14	0
2	C	1	Total C 14 14	0
2	C	1	Total C 14 14	0

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Mol	Chain	Residues	Atoms	AltConf
2	C	1	Total C 12 12	0
2	D	1	Total C 14 14	0
2	D	1	Total C 14 14	0
2	D	1	Total C 10 10	0
2	D	1	Total C 10 10	0
2	D	1	Total C 14 14	0
2	D	1	Total C 14 14	0
2	D	1	Total C 14 14	0
2	D	1	Total C 14 14	0
2	D	1	Total C 14 14	0
2	D	1	Total C 14 14	0
2	D	1	Total C 12 12	0
2	E	1	Total C 14 14	0
2	E	1	Total C 14 14	0
2	E	1	Total C 10 10	0
2	E	1	Total C 10 10	0
2	E	1	Total C 14 14	0
2	E	1	Total C 14 14	0
2	E	1	Total C 14 14	0
2	E	1	Total C 14 14	0
2	E	1	Total C 14 14	0

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Mol	Chain	Residues	Atoms	AltConf
2	E	1	Total C 14 14	0
2	E	1	Total C 12 12	0
2	F	1	Total C 14 14	0
2	F	1	Total C 14 14	0
2	F	1	Total C 14 14	0
2	F	1	Total C 14 14	0
2	F	1	Total C 14 14	0
2	F	1	Total C 12 12	0
2	F	1	Total C 14 14	0
2	F	1	Total C 14 14	0
2	F	1	Total C 10 10	0
2	F	1	Total C 10 10	0
2	G	1	Total C 14 14	0
2	G	1	Total C 14 14	0
2	G	1	Total C 14 14	0
2	G	1	Total C 10 10	0
2	G	1	Total C 10 10	0
2	G	1	Total C 14 14	0
2	G	1	Total C 14 14	0
2	G	1	Total C 14 14	0
2	G	1	Total C 14 14	0

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Mol	Chain	Residues	Atoms	AltConf
2	G	1	Total C 14 14	0
2	G	1	Total C 14 14	0
2	G	1	Total C 12 12	0
2	H	1	Total C 14 14	0
2	H	1	Total C 14 14	0
2	H	1	Total C 10 10	0
2	H	1	Total C 10 10	0
2	H	1	Total C 14 14	0
2	H	1	Total C 14 14	0
2	H	1	Total C 14 14	0
2	H	1	Total C 14 14	0
2	H	1	Total C 14 14	0
2	H	1	Total C 14 14	0
2	H	1	Total C 12 12	0
2	I	1	Total C 14 14	0
2	I	1	Total C 14 14	0
2	I	1	Total C 10 10	0
2	I	1	Total C 10 10	0
2	I	1	Total C 14 14	0
2	I	1	Total C 14 14	0
2	I	1	Total C 14 14	0

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Mol	Chain	Residues	Atoms	AltConf
2	I	1	Total C 14 14	0
2	I	1	Total C 14 14	0
2	I	1	Total C 14 14	0
2	I	1	Total C 12 12	0
2	J	1	Total C 14 14	0
2	J	1	Total C 14 14	0
2	J	1	Total C 10 10	0
2	J	1	Total C 10 10	0
2	J	1	Total C 14 14	0
2	J	1	Total C 14 14	0
2	J	1	Total C 14 14	0
2	J	1	Total C 14 14	0
2	J	1	Total C 14 14	0
2	J	1	Total C 12 12	0
2	K	1	Total C 14 14	0
2	K	1	Total C 14 14	0
2	K	1	Total C 10 10	0
2	K	1	Total C 10 10	0
2	K	1	Total C 14 14	0
2	K	1	Total C 14 14	0

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Mol	Chain	Residues	Atoms	AltConf
2	K	1	Total C 14 14	0
2	K	1	Total C 14 14	0
2	K	1	Total C 14 14	0
2	K	1	Total C 14 14	0
2	K	1	Total C 12 12	0
2	L	1	Total C 14 14	0
2	L	1	Total C 14 14	0
2	L	1	Total C 14 14	0
2	L	1	Total C 14 14	0
2	L	1	Total C 14 14	0
2	L	1	Total C 12 12	0
2	L	1	Total C 14 14	0
2	L	1	Total C 14 14	0
2	L	1	Total C 10 10	0
2	L	1	Total C 10 10	0

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

● Molecule 1: Gap junction delta-2 protein

Chain J:

45%

9%

45%

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

● Molecule 1: Gap junction delta-2 protein

Chain K:

44%

10%

45%

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4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	55379	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 ($e^-/\text{\AA}^2$)	49	Depositor
Minimum defocus (nm)	1500	Depositor
Maximum defocus (nm)	2750	Depositor
Magnification	Not provided	
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.119	Depositor
Minimum map value	-0.101	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.003	Depositor
Recommended contour level	0.01	Depositor
Map size (Å)	263.2, 263.2, 263.2	wwPDB
Map dimensions	400, 400, 400	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.65800005, 0.65800005, 0.65800005	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: MC3

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	0.24	0/1493	0.42	0/2031
1	B	0.24	0/1493	0.43	0/2031
1	C	0.24	0/1493	0.42	0/2031
1	D	0.24	0/1493	0.42	0/2031
1	E	0.24	0/1493	0.42	0/2031
1	F	0.24	0/1493	0.42	0/2031
1	G	0.24	0/1493	0.42	0/2031
1	H	0.24	0/1493	0.42	0/2031
1	I	0.24	0/1493	0.42	0/2031
1	J	0.24	0/1493	0.42	0/2031
1	K	0.24	0/1493	0.42	0/2031
1	L	0.24	0/1493	0.42	0/2031
All	All	0.24	0/17916	0.42	0/24372

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts

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	1458	0	1456	22	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	B	1458	0	1456	29	0
1	C	1458	0	1456	24	0
1	D	1458	0	1456	23	0
1	E	1458	0	1456	23	0
1	F	1458	0	1456	21	0
1	G	1458	0	1456	30	0
1	H	1458	0	1456	22	0
1	I	1458	0	1456	28	0
1	J	1458	0	1456	24	0
1	K	1458	0	1456	25	0
1	L	1458	0	1456	26	0
2	A	158	0	298	4	0
2	B	144	0	271	9	0
2	C	144	0	271	3	0
2	D	144	0	271	3	0
2	E	144	0	271	2	0
2	F	130	0	244	0	0
2	G	158	0	298	8	0
2	H	144	0	271	6	0
2	I	144	0	271	8	0
2	J	144	0	271	3	0
2	K	144	0	271	5	0
2	L	130	0	244	3	0
All	All	19224	0	20724	274	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 7.

The worst 5 of 274 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:230:GLU:OE1	1:L:238:LYS:NZ	2.14	0.78
1:F:230:GLU:OE1	1:K:238:LYS:NZ	2.16	0.78
1:A:248:THR:HG23	2:A:401:MC3:H352	1.67	0.76
1:G:248:THR:HG23	2:G:401:MC3:H352	1.67	0.76
2:H:405:MC3:H352	1:I:248:THR:HG23	1.67	0.76

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	B	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	C	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	D	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	E	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	F	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	G	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	H	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	I	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	J	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	K	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
1	L	177/332 (53%)	176 (99%)	1 (1%)	0	100	100
All	All	2124/3984 (53%)	2112 (99%)	12 (1%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	163/289 (56%)	162 (99%)	1 (1%)	78	83
1	B	163/289 (56%)	162 (99%)	1 (1%)	78	83

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	C	163/289 (56%)	162 (99%)	1 (1%)	78	83
1	D	163/289 (56%)	162 (99%)	1 (1%)	78	83
1	E	163/289 (56%)	162 (99%)	1 (1%)	78	83
1	F	163/289 (56%)	162 (99%)	1 (1%)	78	83
1	G	163/289 (56%)	162 (99%)	1 (1%)	78	83
1	H	163/289 (56%)	162 (99%)	1 (1%)	78	83
1	I	163/289 (56%)	162 (99%)	1 (1%)	78	83
1	J	163/289 (56%)	162 (99%)	1 (1%)	78	83
1	K	163/289 (56%)	162 (99%)	1 (1%)	78	83
1	L	163/289 (56%)	162 (99%)	1 (1%)	78	83
All	All	1956/3468 (56%)	1944 (99%)	12 (1%)	76	83

5 of 12 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	H	238	LYS
1	I	238	LYS
1	L	238	LYS
1	J	238	LYS
1	D	238	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 19 such sidechains are listed below:

Mol	Chain	Res	Type
1	J	50	GLN
1	L	50	GLN
1	L	207	ASN
1	K	207	ASN
1	F	50	GLN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

132 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	$\# Z > 2$
2	MC3	I	409	-	13,13,45	0.27	0	12,12,53	0.89	0
2	MC3	B	405	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	K	406	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	E	406	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	J	402	-	13,13,45	0.29	0	12,12,53	0.87	0
2	MC3	D	402	-	13,13,45	0.29	0	12,12,53	0.87	0
2	MC3	F	404	-	13,13,45	0.27	0	12,12,53	0.90	0
2	MC3	D	403	-	9,9,45	0.29	0	8,8,53	0.79	0
2	MC3	F	403	-	13,13,45	0.28	0	12,12,53	0.85	0
2	MC3	H	402	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	H	411	-	11,11,45	0.27	0	10,10,53	0.87	0
2	MC3	E	404	-	9,9,45	0.28	0	8,8,53	0.85	0
2	MC3	H	406	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	K	405	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	I	405	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	K	402	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	H	404	-	9,9,45	0.27	0	8,8,53	0.84	0
2	MC3	A	408	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	A	409	-	13,13,45	0.28	0	12,12,53	0.85	0
2	MC3	F	402	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	I	403	-	9,9,45	0.29	0	8,8,53	0.79	0
2	MC3	E	403	-	9,9,45	0.29	0	8,8,53	0.79	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MC3	L	410	-	9,9,45	0.27	0	8,8,53	0.84	0
2	MC3	J	404	-	9,9,45	0.27	0	8,8,53	0.84	0
2	MC3	J	411	-	11,11,45	0.27	0	10,10,53	0.87	0
2	MC3	J	403	-	9,9,45	0.29	0	8,8,53	0.79	0
2	MC3	K	403	-	9,9,45	0.30	0	8,8,53	0.80	0
2	MC3	G	409	-	13,13,45	0.28	0	12,12,53	0.85	0
2	MC3	L	405	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	D	408	-	13,13,45	0.28	0	12,12,53	0.85	0
2	MC3	B	411	-	11,11,45	0.27	0	10,10,53	0.88	0
2	MC3	B	407	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	A	405	-	9,9,45	0.27	0	8,8,53	0.85	0
2	MC3	F	407	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	E	411	-	11,11,45	0.27	0	10,10,53	0.88	0
2	MC3	I	410	-	13,13,45	0.27	0	12,12,53	0.91	0
2	MC3	H	409	-	13,13,45	0.27	0	12,12,53	0.90	0
2	MC3	K	411	-	11,11,45	0.27	0	10,10,53	0.87	0
2	MC3	B	406	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	F	405	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	H	403	-	9,9,45	0.29	0	8,8,53	0.79	0
2	MC3	G	407	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	J	405	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	A	412	-	11,11,45	0.27	0	10,10,53	0.87	0
2	MC3	D	404	-	9,9,45	0.27	0	8,8,53	0.84	0
2	MC3	B	404	-	9,9,45	0.27	0	8,8,53	0.84	0
2	MC3	E	410	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	A	410	-	13,13,45	0.27	0	12,12,53	0.90	0
2	MC3	K	404	-	9,9,45	0.27	0	8,8,53	0.84	0
2	MC3	G	410	-	13,13,45	0.27	0	12,12,53	0.89	0
2	MC3	L	404	-	13,13,45	0.27	0	12,12,53	0.89	0
2	MC3	H	405	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	C	402	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	I	404	-	9,9,45	0.27	0	8,8,53	0.84	0
2	MC3	B	402	-	13,13,45	0.29	0	12,12,53	0.88	0
2	MC3	G	405	-	9,9,45	0.27	0	8,8,53	0.85	0
2	MC3	D	405	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	G	401	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	I	402	-	13,13,45	0.29	0	12,12,53	0.87	0
2	MC3	A	407	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	G	412	-	11,11,45	0.27	0	10,10,53	0.87	0
2	MC3	K	401	-	13,13,45	0.29	0	12,12,53	0.87	0
2	MC3	C	406	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	J	410	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	J	407	-	13,13,45	0.26	0	12,12,53	0.91	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MC3	A	404	-	9,9,45	0.29	0	8,8,53	0.79	0
2	MC3	F	410	-	9,9,45	0.27	0	8,8,53	0.85	0
2	MC3	D	409	-	13,13,45	0.27	0	12,12,53	0.90	0
2	MC3	J	406	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	H	401	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	C	401	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	E	405	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	L	407	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	G	404	-	9,9,45	0.29	0	8,8,53	0.79	0
2	MC3	D	407	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	F	409	-	9,9,45	0.29	0	8,8,53	0.79	0
2	MC3	J	409	-	13,13,45	0.27	0	12,12,53	0.89	0
2	MC3	I	401	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	F	408	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	C	405	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	G	406	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	E	409	-	13,13,45	0.27	0	12,12,53	0.89	0
2	MC3	H	410	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	G	402	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	A	401	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	C	410	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	K	409	-	13,13,45	0.27	0	12,12,53	0.90	0
2	MC3	B	403	-	9,9,45	0.30	0	8,8,53	0.80	0
2	MC3	D	410	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	G	403	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	K	407	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	I	407	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	L	406	-	11,11,45	0.27	0	10,10,53	0.88	0
2	MC3	A	411	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	E	401	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	H	408	-	13,13,45	0.28	0	12,12,53	0.85	0
2	MC3	C	408	-	13,13,45	0.28	0	12,12,53	0.85	0
2	MC3	G	411	-	13,13,45	0.27	0	12,12,53	0.91	0
2	MC3	C	403	-	9,9,45	0.30	0	8,8,53	0.80	0
2	MC3	B	410	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	D	411	-	11,11,45	0.27	0	10,10,53	0.87	0
2	MC3	H	407	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	J	401	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	G	408	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	A	402	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	J	408	-	13,13,45	0.28	0	12,12,53	0.85	0
2	MC3	L	402	-	13,13,45	0.25	0	12,12,53	0.90	0
2	MC3	C	404	-	9,9,45	0.27	0	8,8,53	0.84	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	MC3	C	411	-	11,11,45	0.27	0	10,10,53	0.88	0
2	MC3	A	406	-	13,13,45	0.29	0	12,12,53	0.84	0
2	MC3	C	409	-	13,13,45	0.27	0	12,12,53	0.89	0
2	MC3	B	409	-	13,13,45	0.28	0	12,12,53	0.90	0
2	MC3	E	402	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	L	401	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	F	401	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	B	408	-	13,13,45	0.28	0	12,12,53	0.85	0
2	MC3	D	401	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	K	410	-	13,13,45	0.27	0	12,12,53	0.91	0
2	MC3	C	407	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	I	411	-	11,11,45	0.27	0	10,10,53	0.87	0
2	MC3	A	403	-	13,13,45	0.29	0	12,12,53	0.87	0
2	MC3	E	408	-	13,13,45	0.28	0	12,12,53	0.85	0
2	MC3	K	408	-	13,13,45	0.29	0	12,12,53	0.85	0
2	MC3	B	401	-	13,13,45	0.28	0	12,12,53	0.87	0
2	MC3	L	408	-	13,13,45	0.29	0	12,12,53	0.87	0
2	MC3	L	409	-	9,9,45	0.30	0	8,8,53	0.80	0
2	MC3	L	403	-	13,13,45	0.29	0	12,12,53	0.86	0
2	MC3	E	407	-	13,13,45	0.26	0	12,12,53	0.91	0
2	MC3	I	406	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	D	406	-	13,13,45	0.26	0	12,12,53	0.90	0
2	MC3	F	406	-	11,11,45	0.27	0	10,10,53	0.87	0
2	MC3	I	408	-	13,13,45	0.28	0	12,12,53	0.85	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. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MC3	I	409	-	-	2/11/11/49	-
2	MC3	B	405	-	-	8/11/11/49	-
2	MC3	K	406	-	-	3/11/11/49	-
2	MC3	E	406	-	-	3/11/11/49	-
2	MC3	J	402	-	-	6/11/11/49	-
2	MC3	D	402	-	-	6/11/11/49	-
2	MC3	F	404	-	-	2/11/11/49	-
2	MC3	D	403	-	-	3/7/7/49	-
2	MC3	F	403	-	-	5/11/11/49	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MC3	H	402	-	-	6/11/11/49	-
2	MC3	H	411	-	-	4/9/9/49	-
2	MC3	E	404	-	-	1/7/7/49	-
2	MC3	H	406	-	-	3/11/11/49	-
2	MC3	K	405	-	-	8/11/11/49	-
2	MC3	I	405	-	-	8/11/11/49	-
2	MC3	K	402	-	-	6/11/11/49	-
2	MC3	H	404	-	-	1/7/7/49	-
2	MC3	A	408	-	-	4/11/11/49	-
2	MC3	A	409	-	-	5/11/11/49	-
2	MC3	F	402	-	-	4/11/11/49	-
2	MC3	I	403	-	-	3/7/7/49	-
2	MC3	E	403	-	-	3/7/7/49	-
2	MC3	L	410	-	-	1/7/7/49	-
2	MC3	J	404	-	-	1/7/7/49	-
2	MC3	J	411	-	-	4/9/9/49	-
2	MC3	J	403	-	-	3/7/7/49	-
2	MC3	K	403	-	-	3/7/7/49	-
2	MC3	G	409	-	-	5/11/11/49	-
2	MC3	L	405	-	-	3/11/11/49	-
2	MC3	D	408	-	-	5/11/11/49	-
2	MC3	B	411	-	-	4/9/9/49	-
2	MC3	B	407	-	-	4/11/11/49	-
2	MC3	A	405	-	-	1/7/7/49	-
2	MC3	F	407	-	-	5/11/11/49	-
2	MC3	E	411	-	-	4/9/9/49	-
2	MC3	I	410	-	-	3/11/11/49	-
2	MC3	H	409	-	-	2/11/11/49	-
2	MC3	K	411	-	-	4/9/9/49	-
2	MC3	B	406	-	-	3/11/11/49	-
2	MC3	F	405	-	-	3/11/11/49	-
2	MC3	H	403	-	-	3/7/7/49	-
2	MC3	G	407	-	-	3/11/11/49	-
2	MC3	J	405	-	-	8/11/11/49	-
2	MC3	A	412	-	-	4/9/9/49	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MC3	D	404	-	-	1/7/7/49	-
2	MC3	B	404	-	-	1/7/7/49	-
2	MC3	E	410	-	-	3/11/11/49	-
2	MC3	A	410	-	-	2/11/11/49	-
2	MC3	K	404	-	-	1/7/7/49	-
2	MC3	G	410	-	-	2/11/11/49	-
2	MC3	L	404	-	-	2/11/11/49	-
2	MC3	H	405	-	-	8/11/11/49	-
2	MC3	C	402	-	-	6/11/11/49	-
2	MC3	I	404	-	-	1/7/7/49	-
2	MC3	B	402	-	-	6/11/11/49	-
2	MC3	G	405	-	-	1/7/7/49	-
2	MC3	D	405	-	-	8/11/11/49	-
2	MC3	G	401	-	-	8/11/11/49	-
2	MC3	I	402	-	-	6/11/11/49	-
2	MC3	A	407	-	-	3/11/11/49	-
2	MC3	G	412	-	-	4/9/9/49	-
2	MC3	K	401	-	-	5/11/11/49	-
2	MC3	C	406	-	-	3/11/11/49	-
2	MC3	J	410	-	-	3/11/11/49	-
2	MC3	J	407	-	-	4/11/11/49	-
2	MC3	A	404	-	-	3/7/7/49	-
2	MC3	F	410	-	-	1/7/7/49	-
2	MC3	D	409	-	-	2/11/11/49	-
2	MC3	J	406	-	-	3/11/11/49	-
2	MC3	H	401	-	-	5/11/11/49	-
2	MC3	C	401	-	-	5/11/11/49	-
2	MC3	E	405	-	-	8/11/11/49	-
2	MC3	L	407	-	-	5/11/11/49	-
2	MC3	G	404	-	-	3/7/7/49	-
2	MC3	D	407	-	-	4/11/11/49	-
2	MC3	F	409	-	-	3/7/7/49	-
2	MC3	J	409	-	-	2/11/11/49	-
2	MC3	I	401	-	-	5/11/11/49	-
2	MC3	F	408	-	-	6/11/11/49	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MC3	C	405	-	-	8/11/11/49	-
2	MC3	G	406	-	-	8/11/11/49	-
2	MC3	E	409	-	-	2/11/11/49	-
2	MC3	H	410	-	-	3/11/11/49	-
2	MC3	G	402	-	-	5/11/11/49	-
2	MC3	A	401	-	-	8/11/11/49	-
2	MC3	C	410	-	-	3/11/11/49	-
2	MC3	K	409	-	-	2/11/11/49	-
2	MC3	B	403	-	-	3/7/7/49	-
2	MC3	D	410	-	-	3/11/11/49	-
2	MC3	G	403	-	-	6/11/11/49	-
2	MC3	K	407	-	-	4/11/11/49	-
2	MC3	I	407	-	-	4/11/11/49	-
2	MC3	L	406	-	-	4/9/9/49	-
2	MC3	A	411	-	-	3/11/11/49	-
2	MC3	E	401	-	-	5/11/11/49	-
2	MC3	H	408	-	-	5/11/11/49	-
2	MC3	C	408	-	-	5/11/11/49	-
2	MC3	G	411	-	-	3/11/11/49	-
2	MC3	C	403	-	-	3/7/7/49	-
2	MC3	B	410	-	-	3/11/11/49	-
2	MC3	D	411	-	-	4/9/9/49	-
2	MC3	H	407	-	-	4/11/11/49	-
2	MC3	J	401	-	-	5/11/11/49	-
2	MC3	G	408	-	-	4/11/11/49	-
2	MC3	A	402	-	-	5/11/11/49	-
2	MC3	J	408	-	-	5/11/11/49	-
2	MC3	L	402	-	-	4/11/11/49	-
2	MC3	C	404	-	-	1/7/7/49	-
2	MC3	C	411	-	-	4/9/9/49	-
2	MC3	A	406	-	-	8/11/11/49	-
2	MC3	C	409	-	-	2/11/11/49	-
2	MC3	B	409	-	-	2/11/11/49	-
2	MC3	E	402	-	-	6/11/11/49	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	MC3	L	401	-	-	3/11/11/49	-
2	MC3	F	401	-	-	3/11/11/49	-
2	MC3	B	408	-	-	4/11/11/49	-
2	MC3	D	401	-	-	5/11/11/49	-
2	MC3	K	410	-	-	3/11/11/49	-
2	MC3	C	407	-	-	4/11/11/49	-
2	MC3	I	411	-	-	4/9/9/49	-
2	MC3	A	403	-	-	6/11/11/49	-
2	MC3	E	408	-	-	4/11/11/49	-
2	MC3	K	408	-	-	4/11/11/49	-
2	MC3	B	401	-	-	5/11/11/49	-
2	MC3	L	408	-	-	6/11/11/49	-
2	MC3	L	409	-	-	3/7/7/49	-
2	MC3	L	403	-	-	4/11/11/49	-
2	MC3	E	407	-	-	4/11/11/49	-
2	MC3	I	406	-	-	3/11/11/49	-
2	MC3	D	406	-	-	3/11/11/49	-
2	MC3	F	406	-	-	4/9/9/49	-
2	MC3	I	408	-	-	5/11/11/49	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

5 of 524 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	402	MC3	C37-C38-C39-C40
2	F	407	MC3	C37-C38-C39-C40
2	F	409	MC3	C33-C34-C35-C36
2	I	401	MC3	C37-C38-C39-C40
2	J	401	MC3	C37-C38-C39-C40

There are no ring outliers.

44 monomers are involved in 54 short contacts:

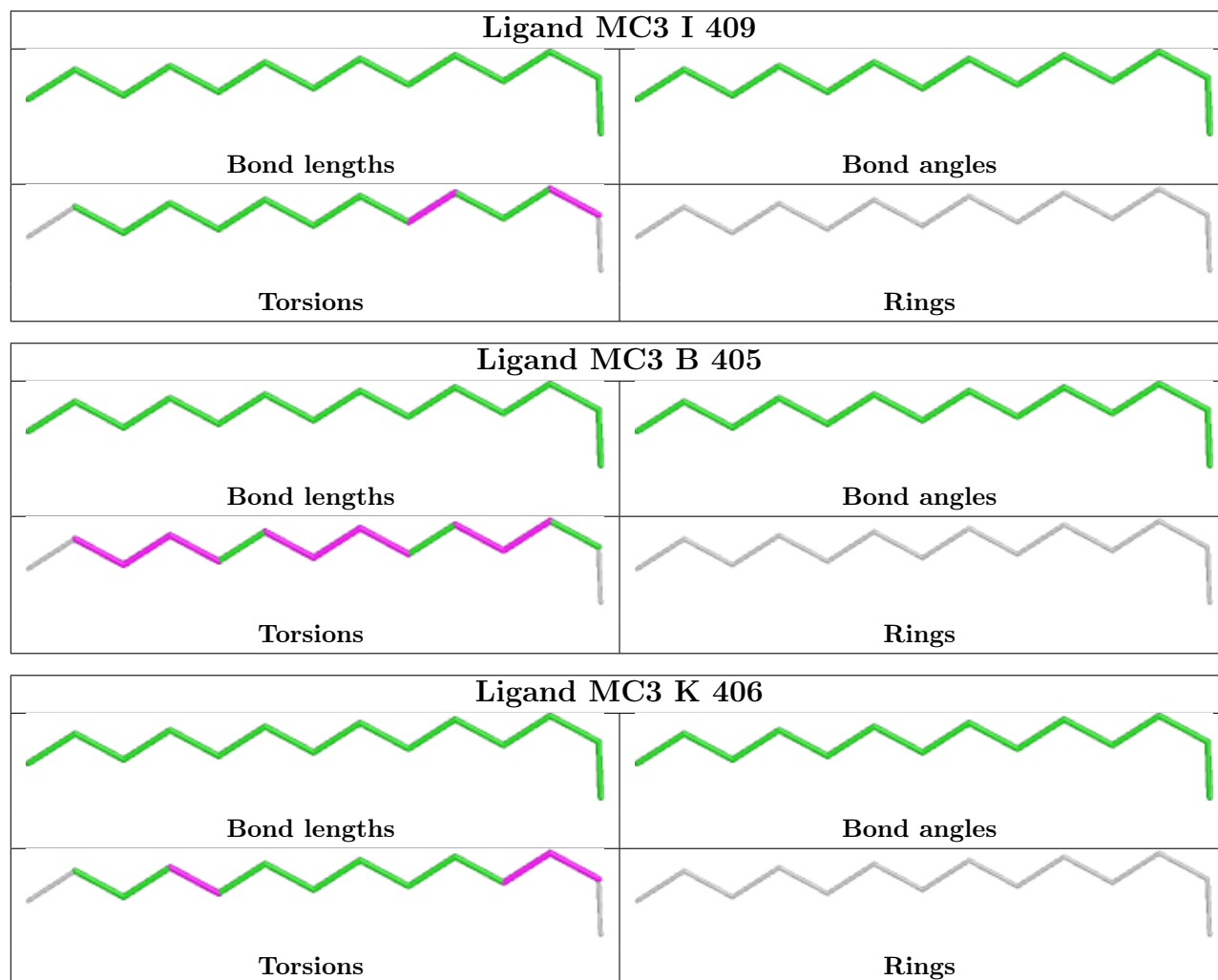
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	405	MC3	1	0
2	K	406	MC3	1	0
2	E	406	MC3	1	0
2	J	402	MC3	1	0
2	D	402	MC3	1	0
2	H	402	MC3	1	0
2	H	406	MC3	1	0
2	K	405	MC3	1	0
2	I	405	MC3	1	0
2	H	404	MC3	1	0
2	G	409	MC3	1	0
2	L	405	MC3	2	0
2	B	411	MC3	1	0
2	A	405	MC3	1	0
2	I	410	MC3	4	0
2	J	405	MC3	1	0
2	D	404	MC3	1	0
2	B	404	MC3	1	0
2	H	405	MC3	1	0
2	C	402	MC3	1	0
2	B	402	MC3	1	0
2	G	405	MC3	1	0
2	D	405	MC3	1	0
2	G	401	MC3	1	0
2	C	406	MC3	1	0
2	J	406	MC3	1	0
2	E	405	MC3	1	0
2	C	405	MC3	1	0
2	G	406	MC3	1	0
2	H	410	MC3	1	0
2	A	401	MC3	1	0
2	G	403	MC3	1	0
2	H	408	MC3	1	0
2	G	411	MC3	3	0
2	B	410	MC3	4	0
2	A	406	MC3	1	0
2	B	408	MC3	1	0
2	K	410	MC3	2	0
2	I	411	MC3	1	0
2	A	403	MC3	1	0
2	K	408	MC3	1	0
2	L	403	MC3	1	0
2	I	406	MC3	1	0

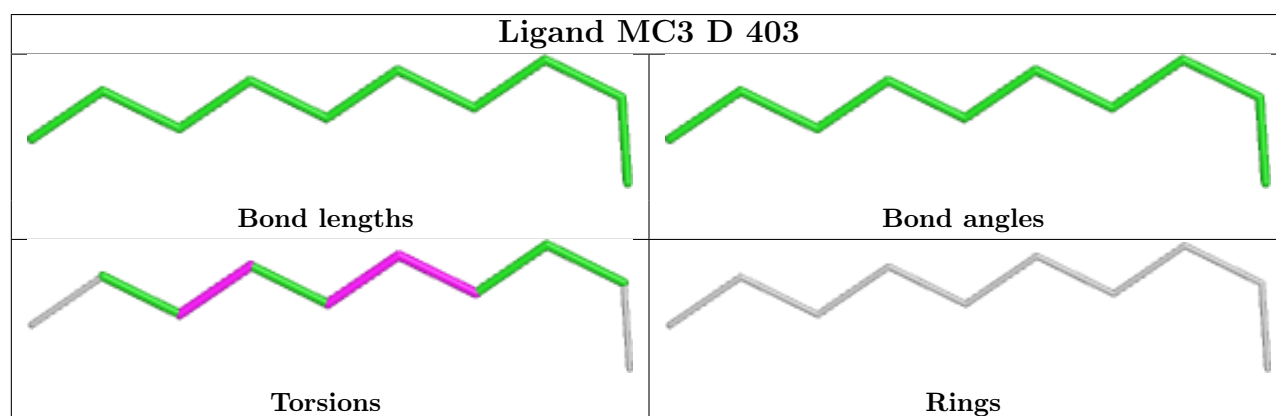
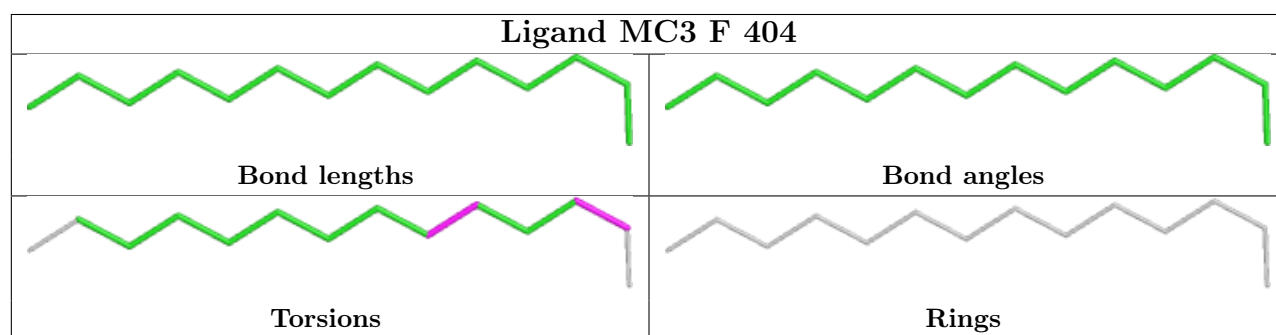
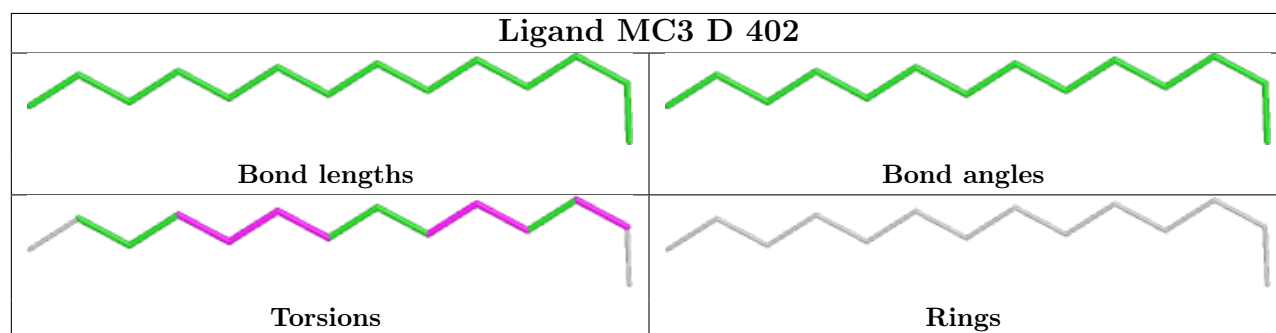
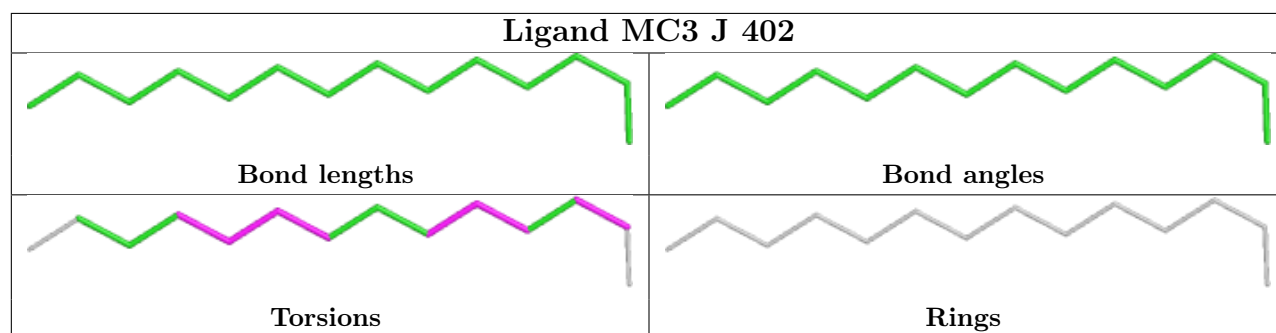
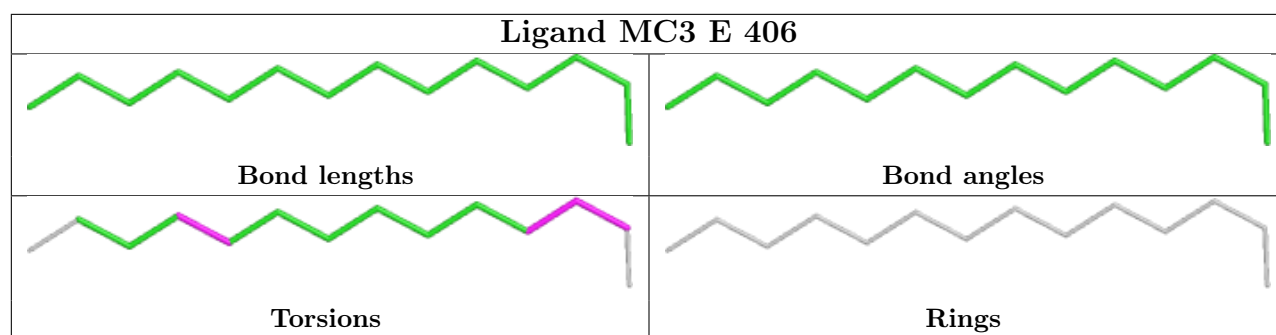
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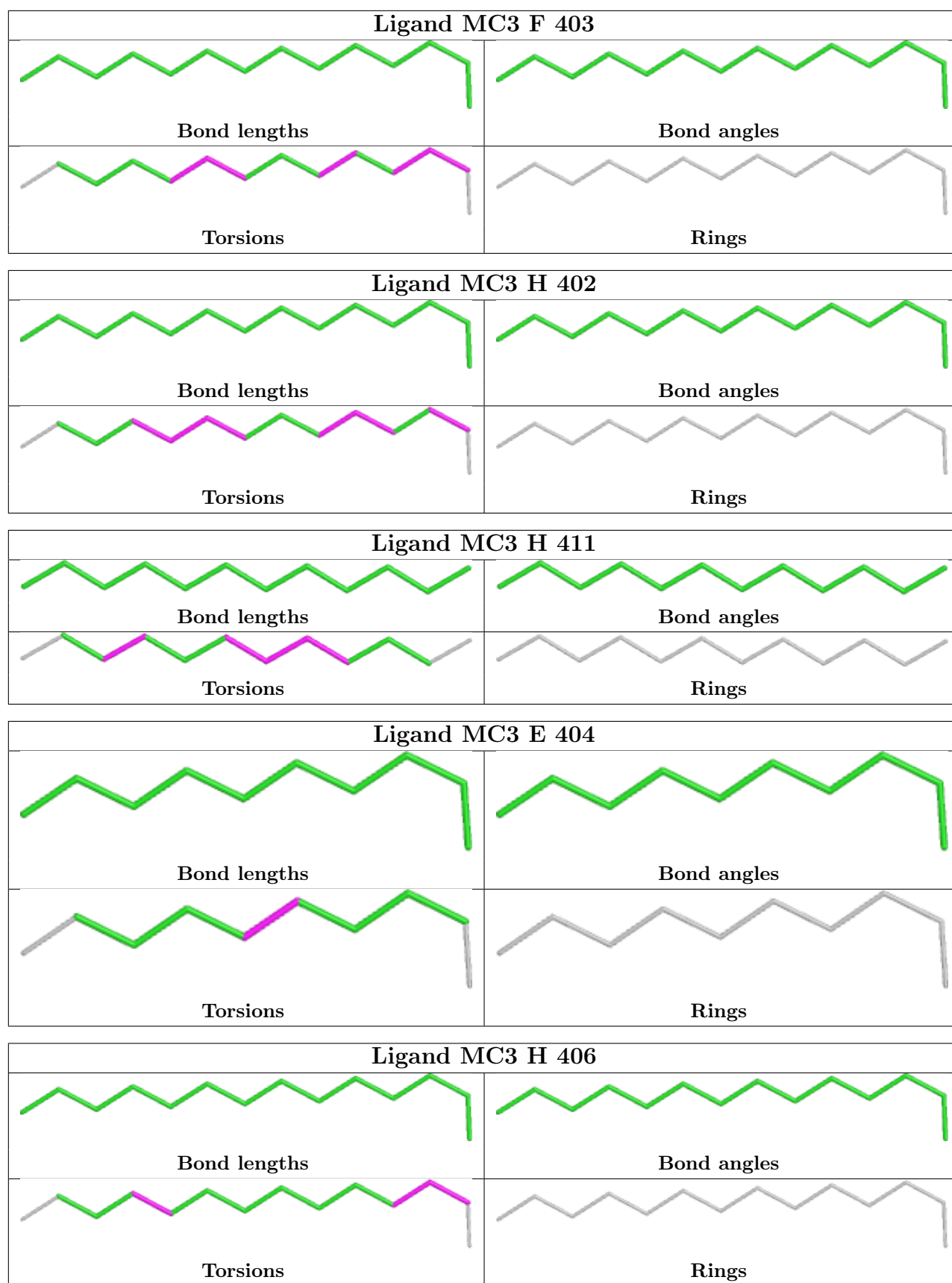
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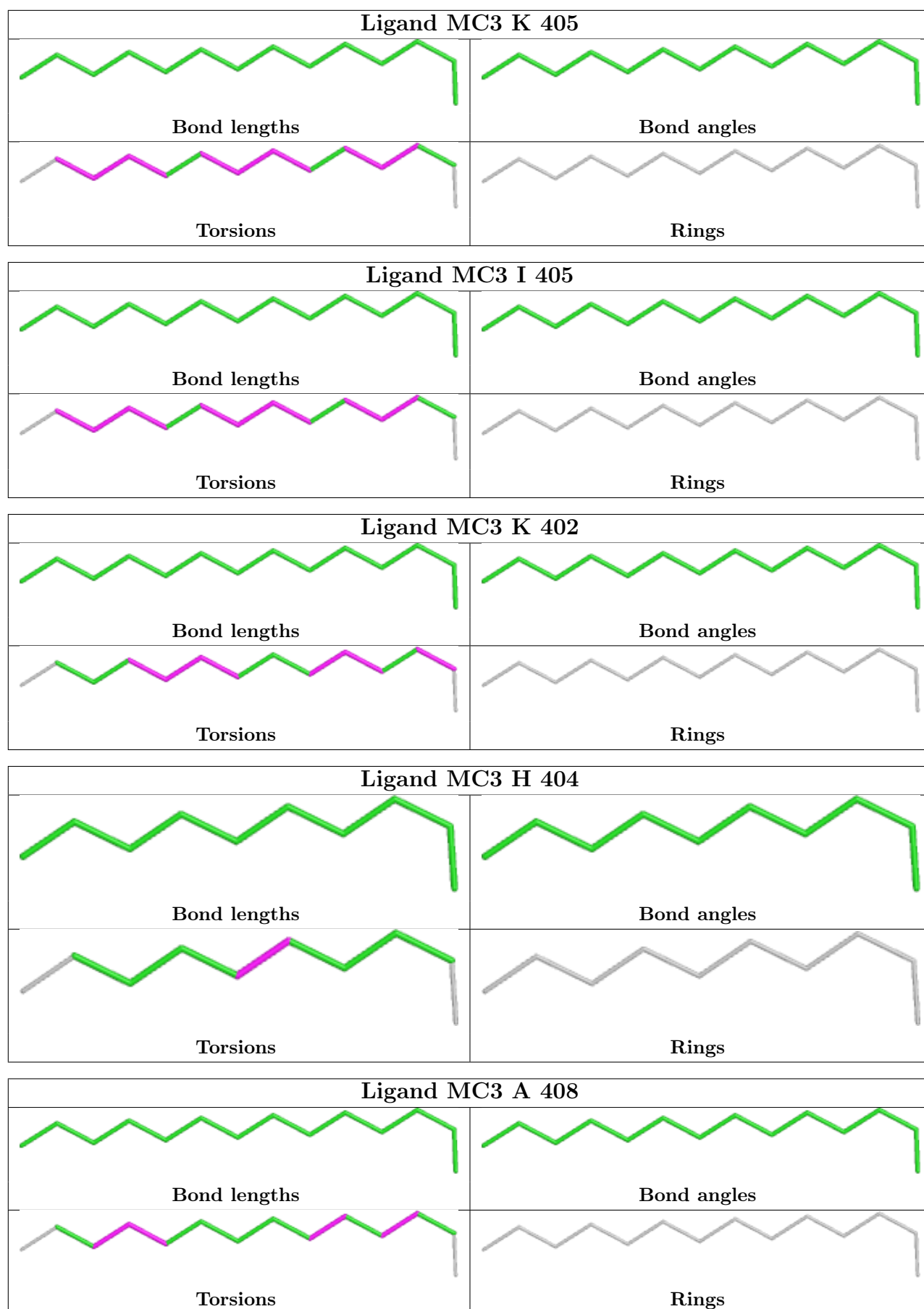
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	I	408	MC3	1	0

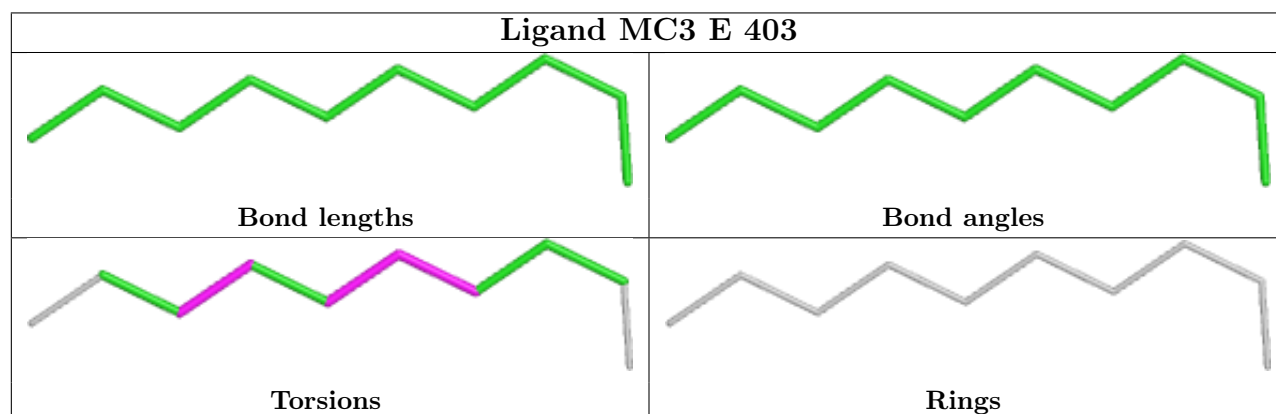
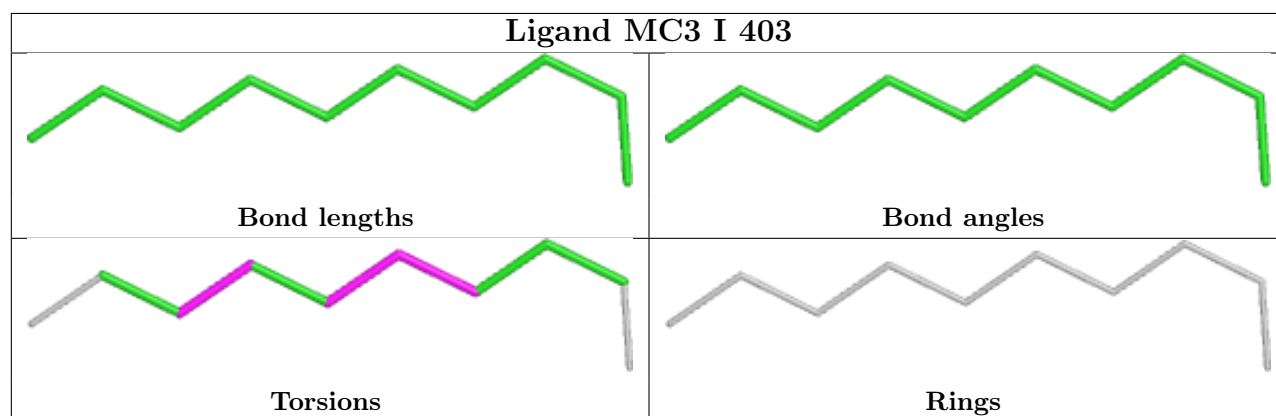
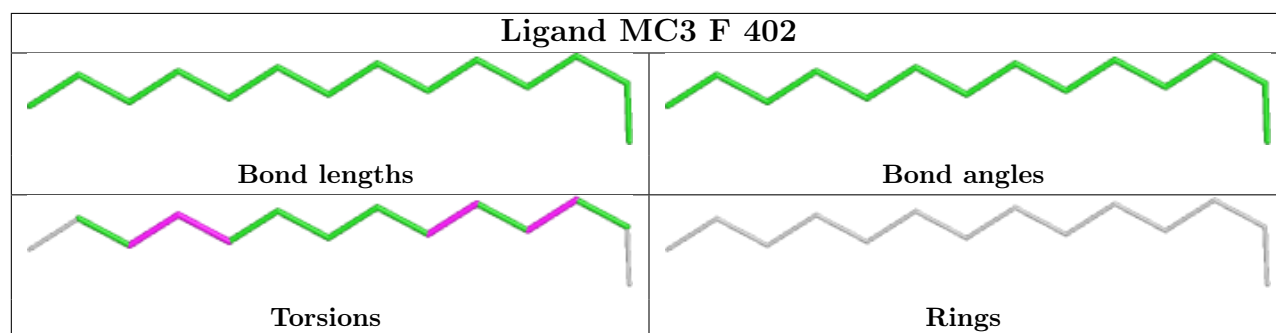
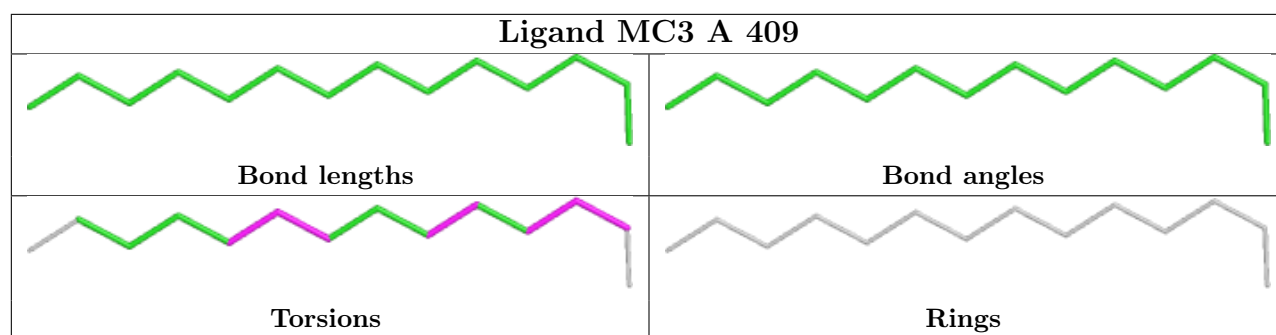
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

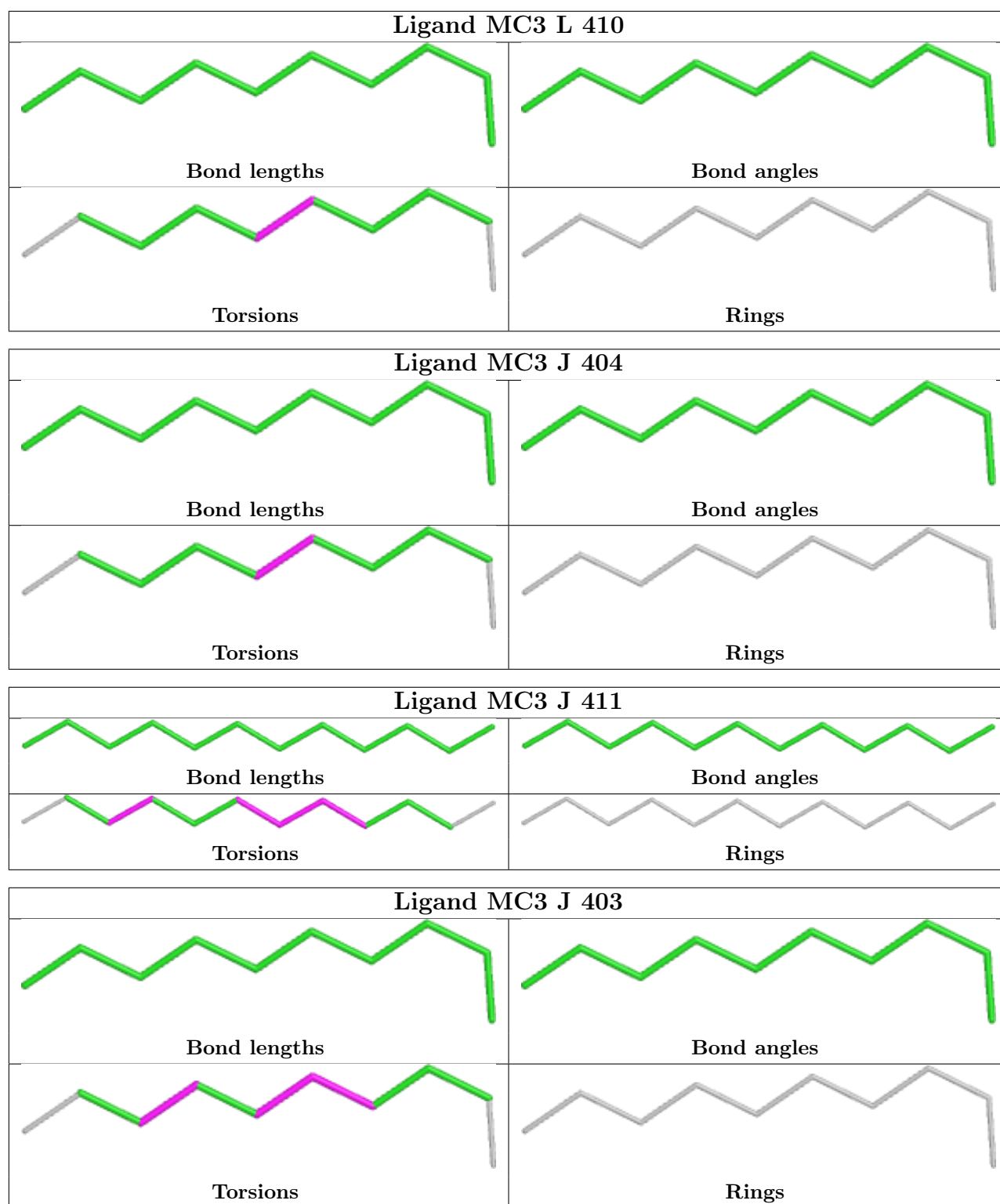


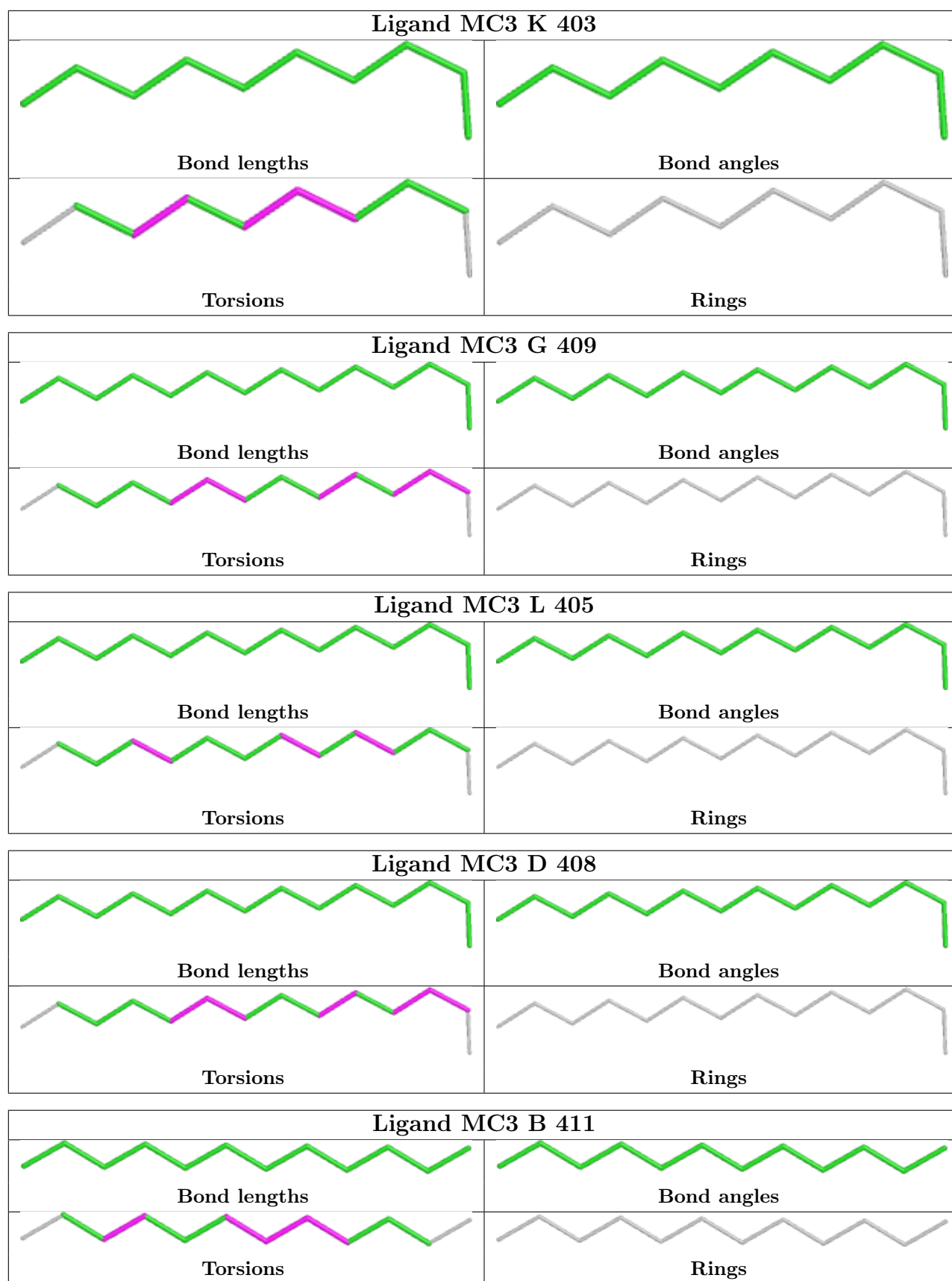


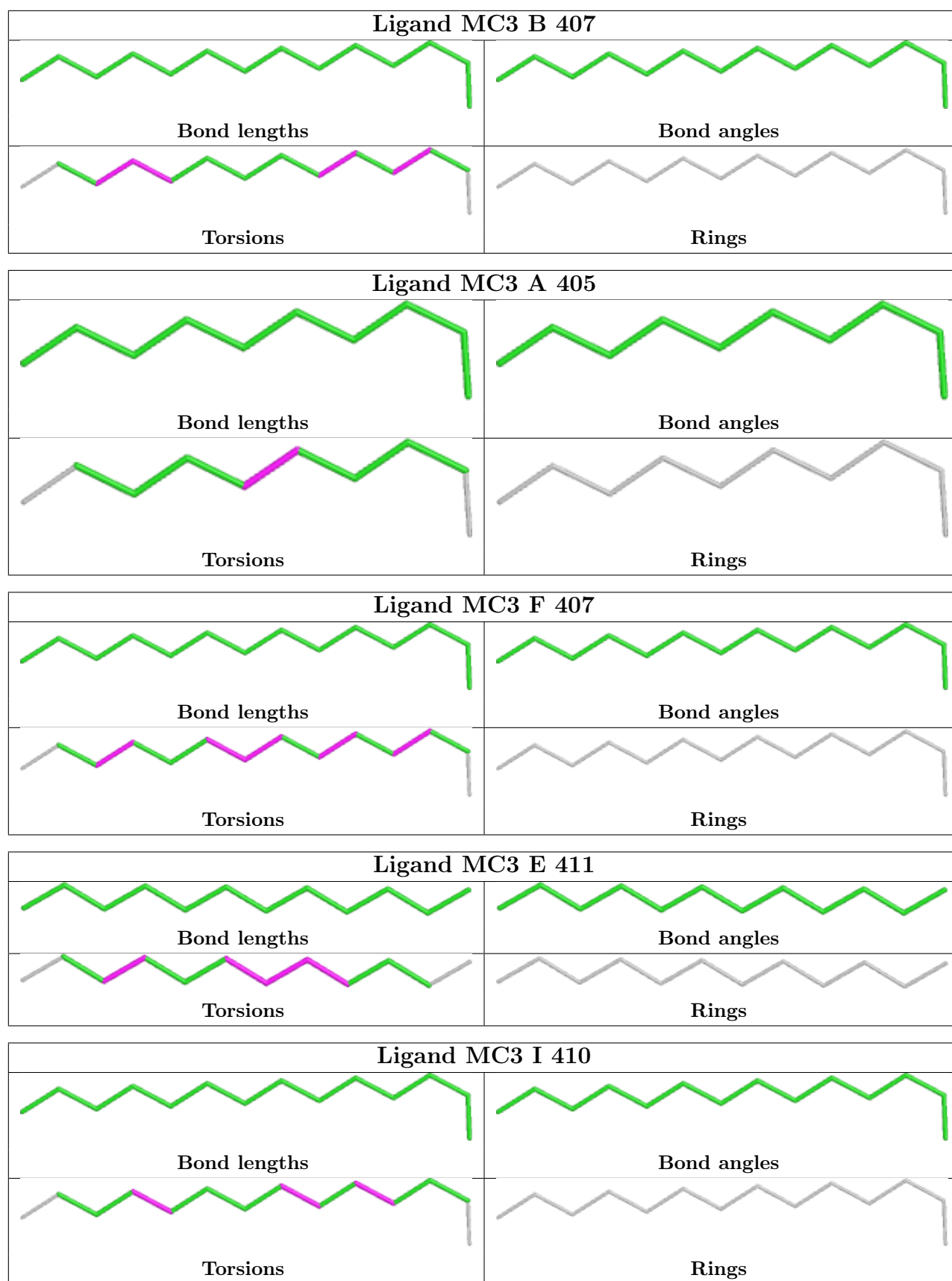


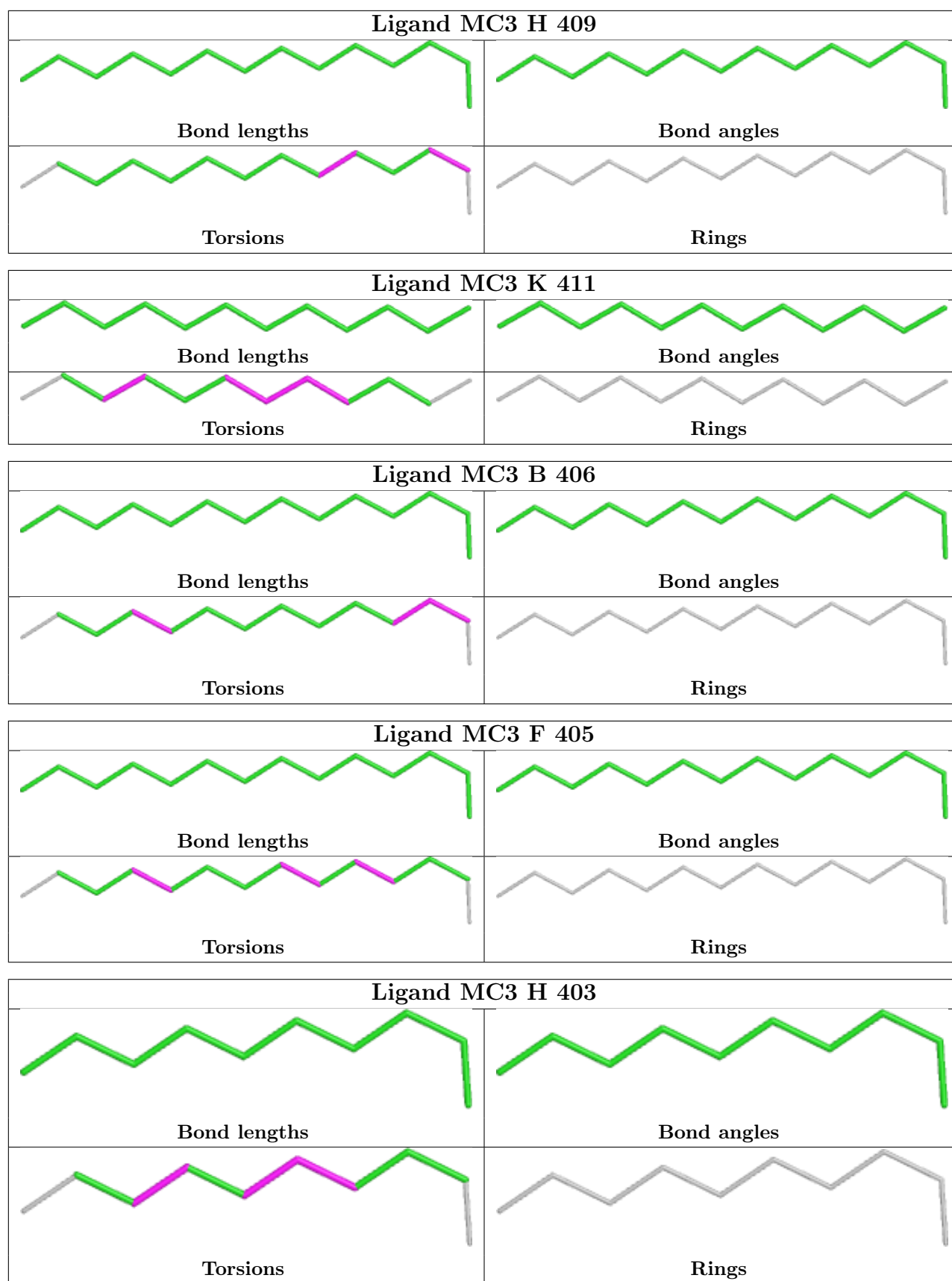


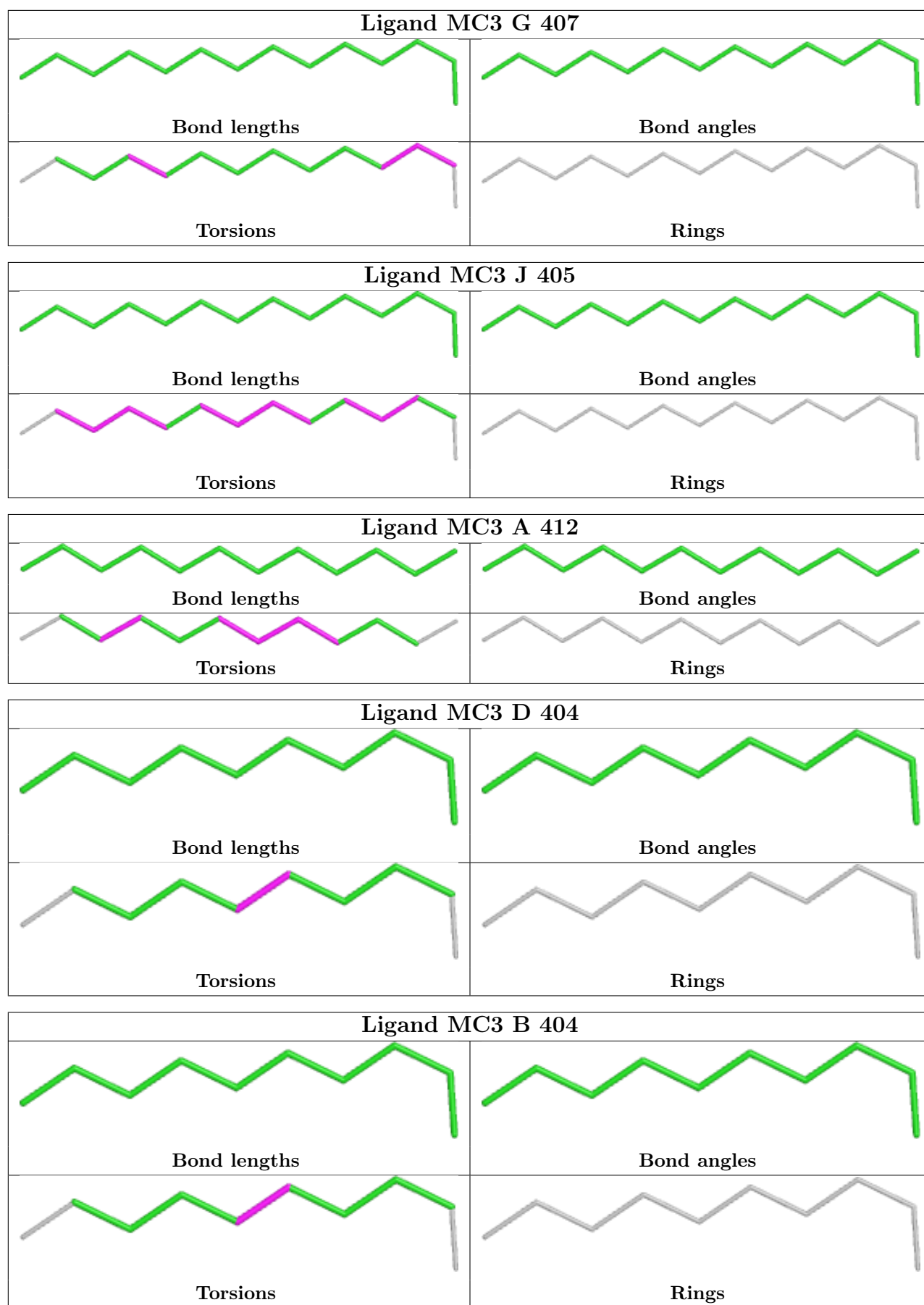


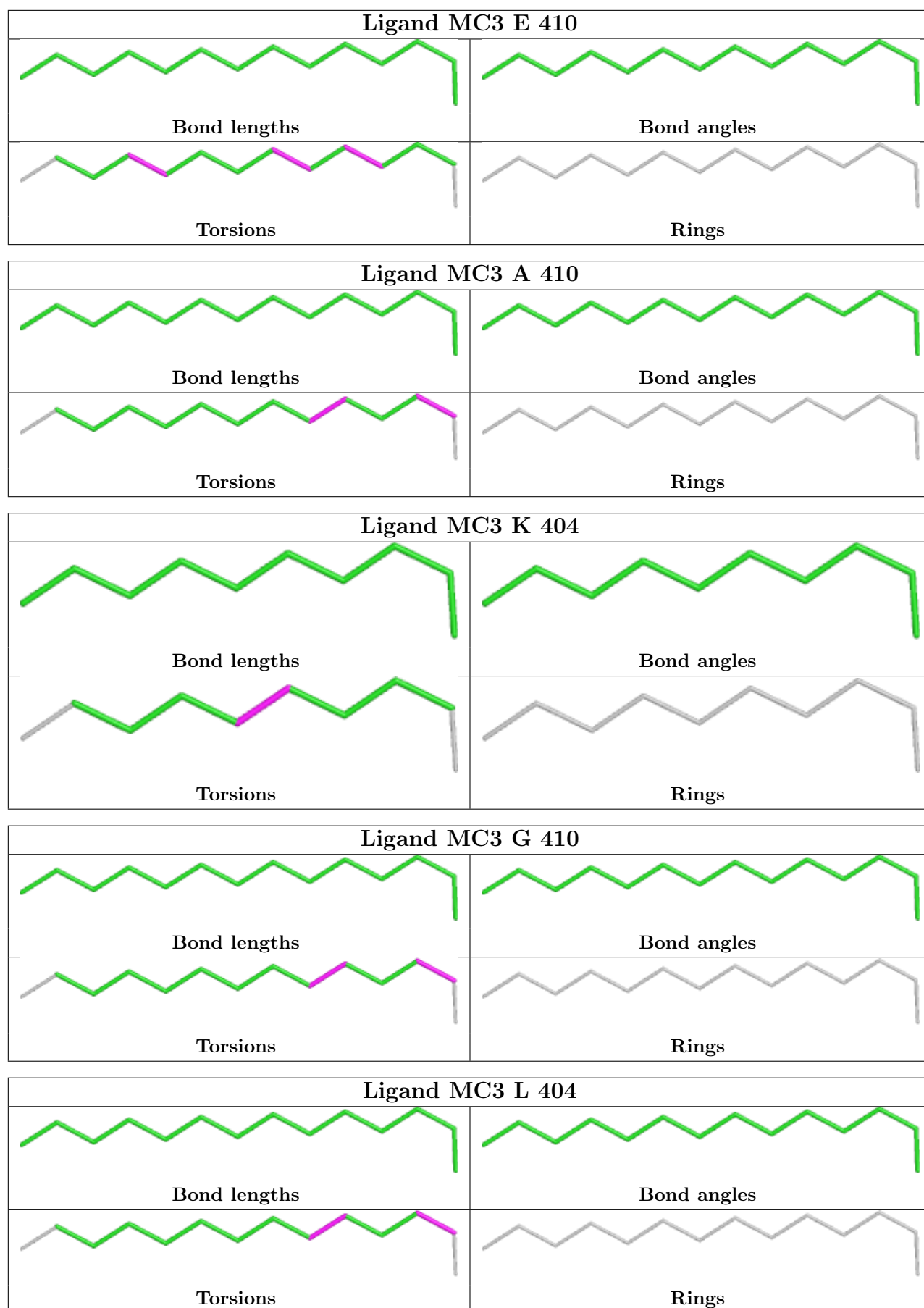


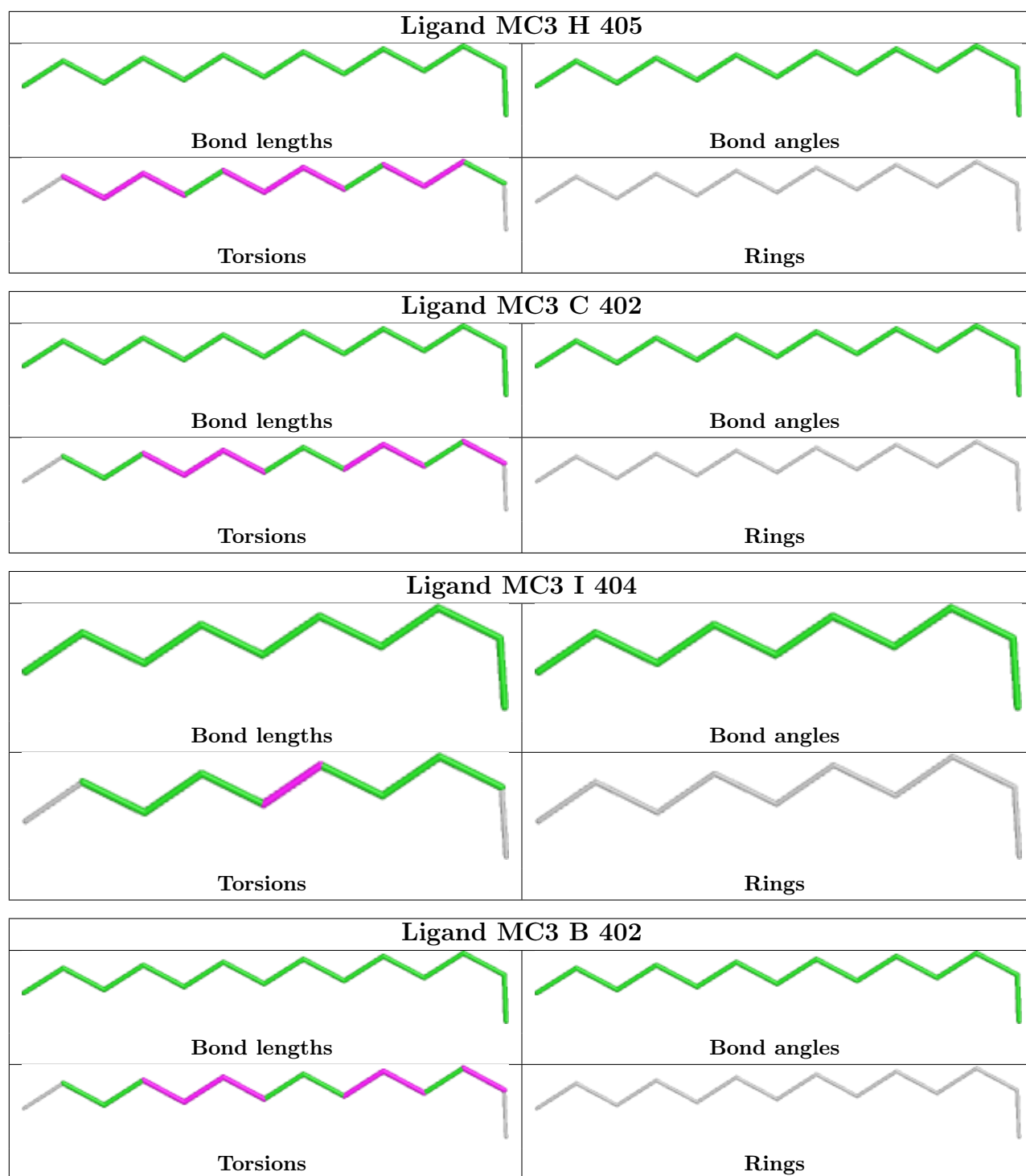


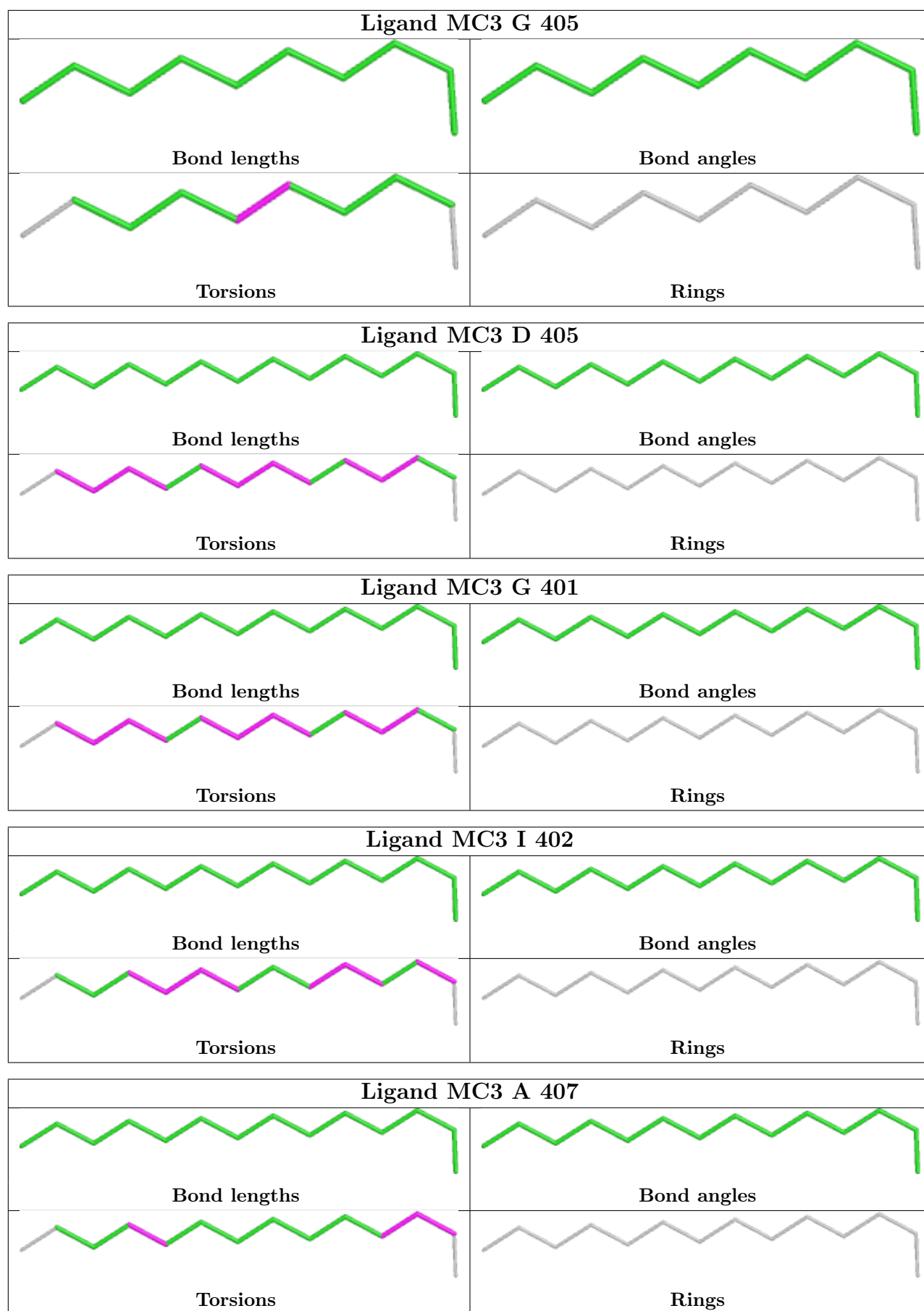


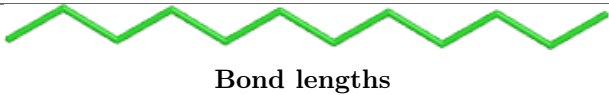


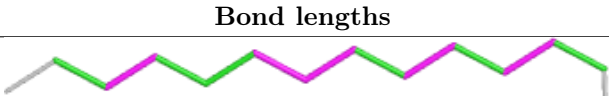
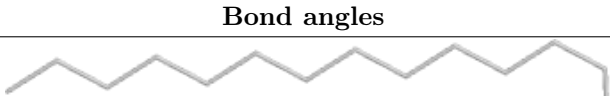


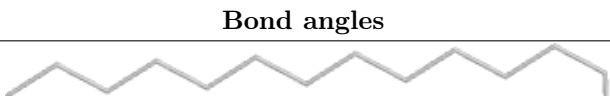


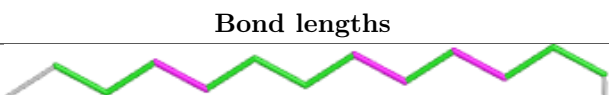
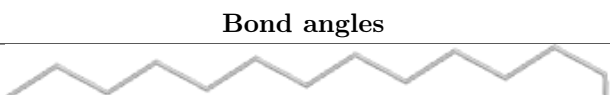


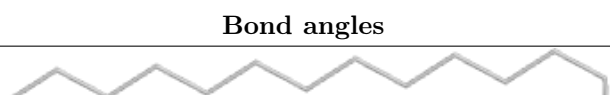


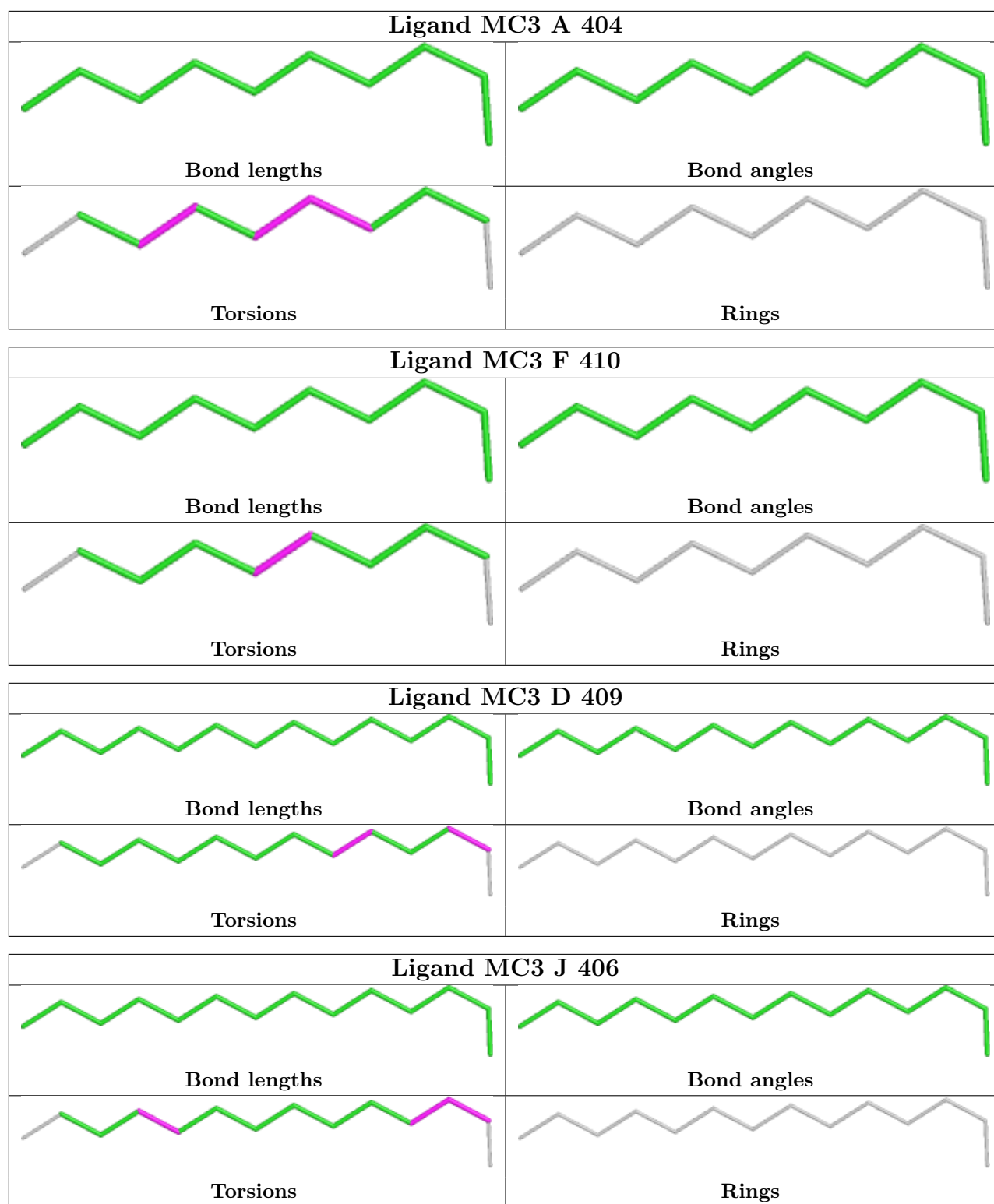
Ligand MC3 G 412	
 Bond lengths	 Bond angles
 Torsions	 Rings

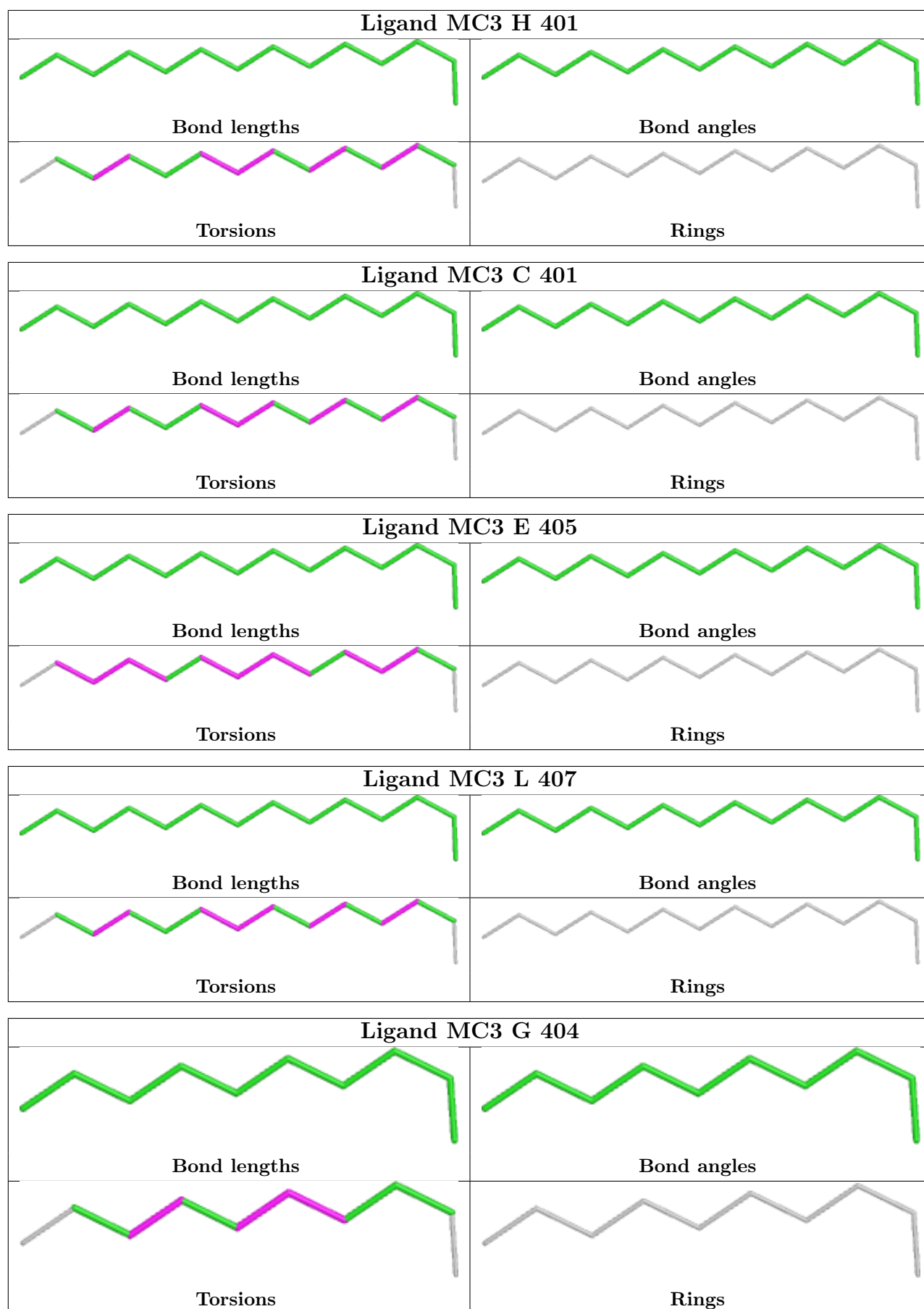
Ligand MC3 K 401	
 Bond lengths	 Bond angles
 Torsions	 Rings

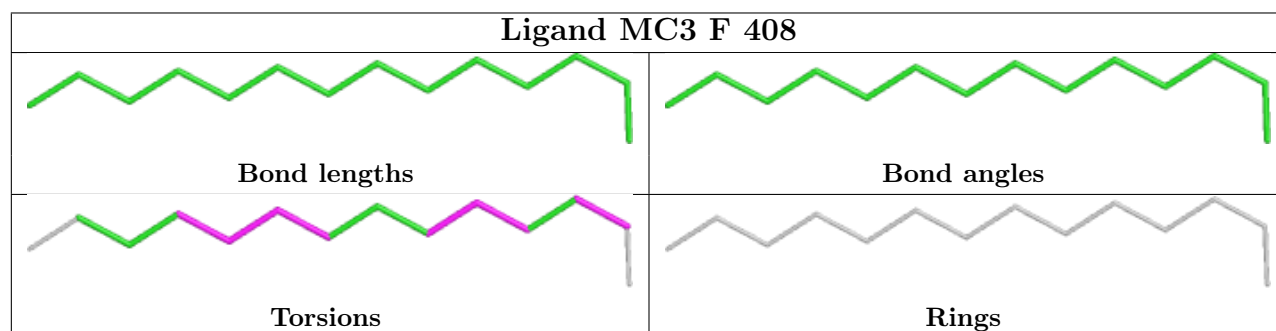
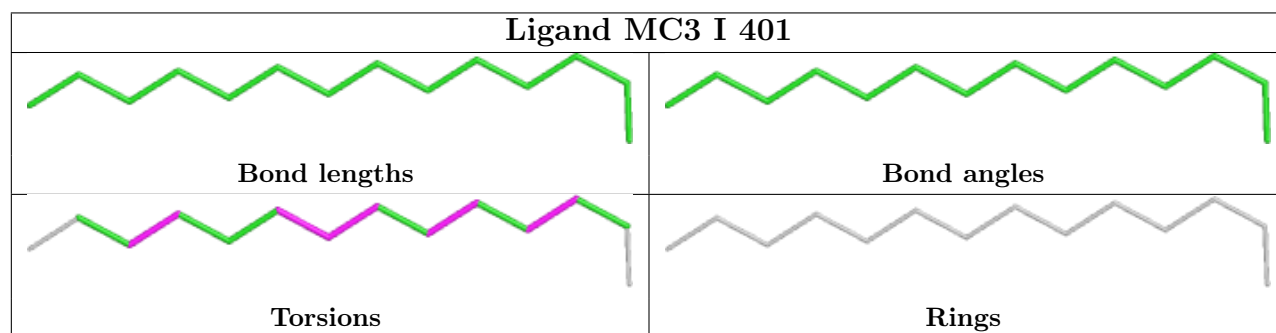
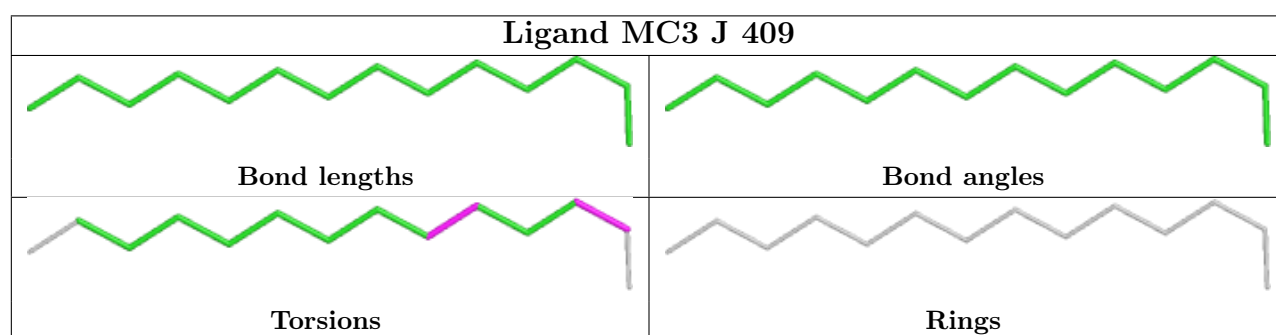
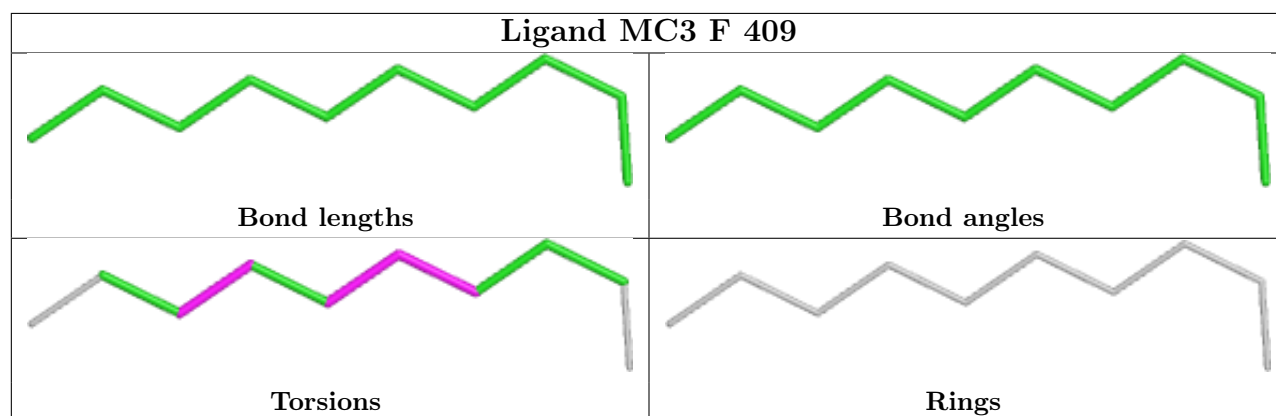
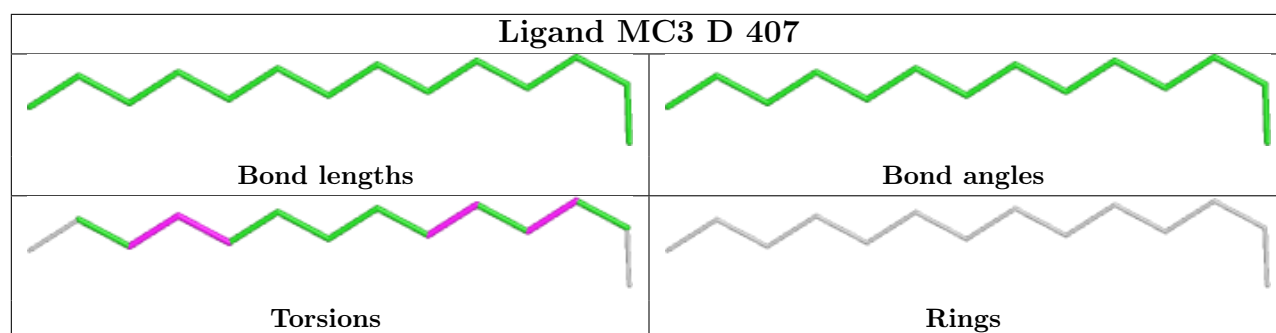
Ligand MC3 C 406	
 Bond lengths	 Bond angles
 Torsions	 Rings

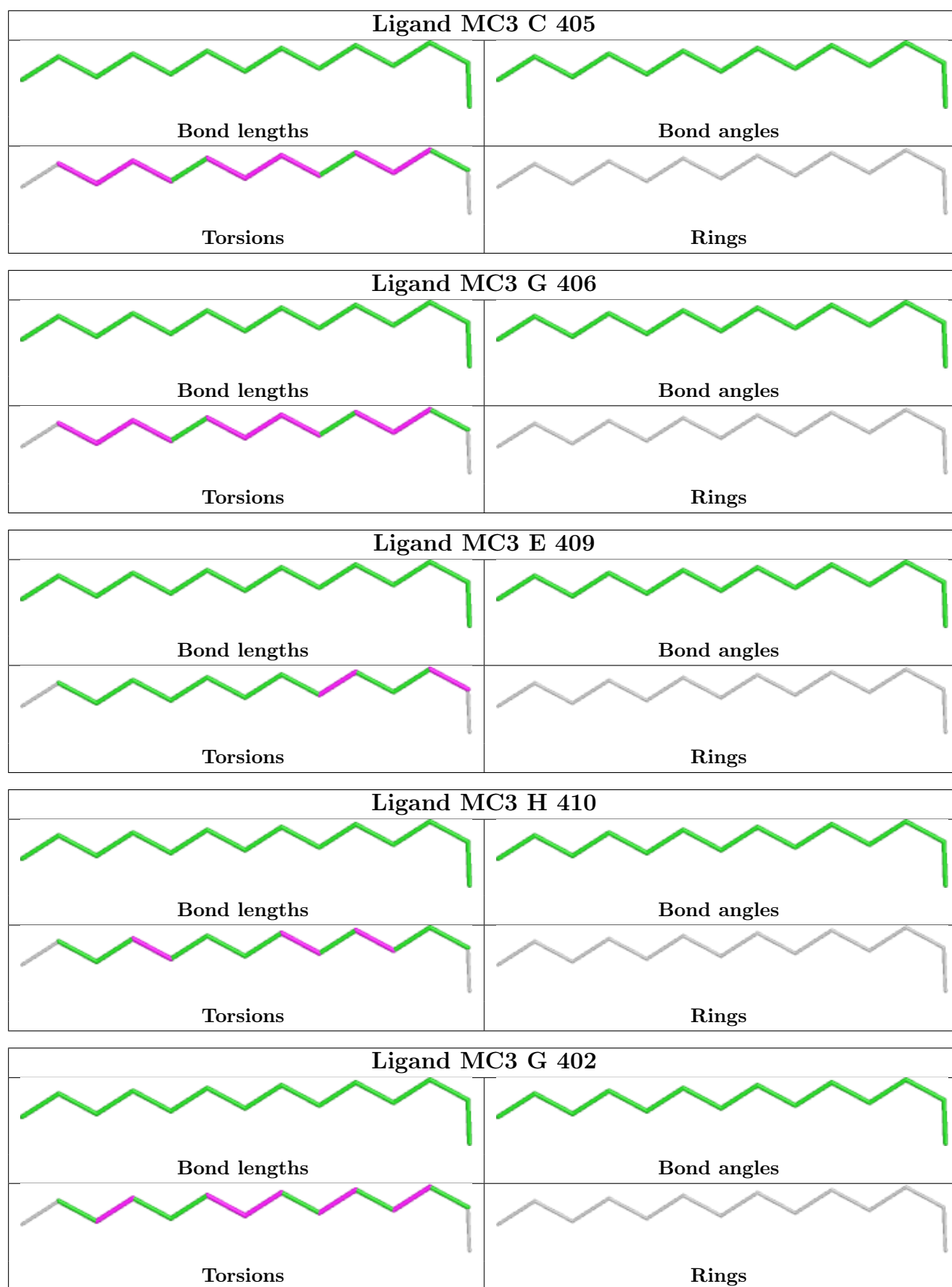
Ligand MC3 J 410	
 Bond lengths	 Bond angles
 Torsions	 Rings

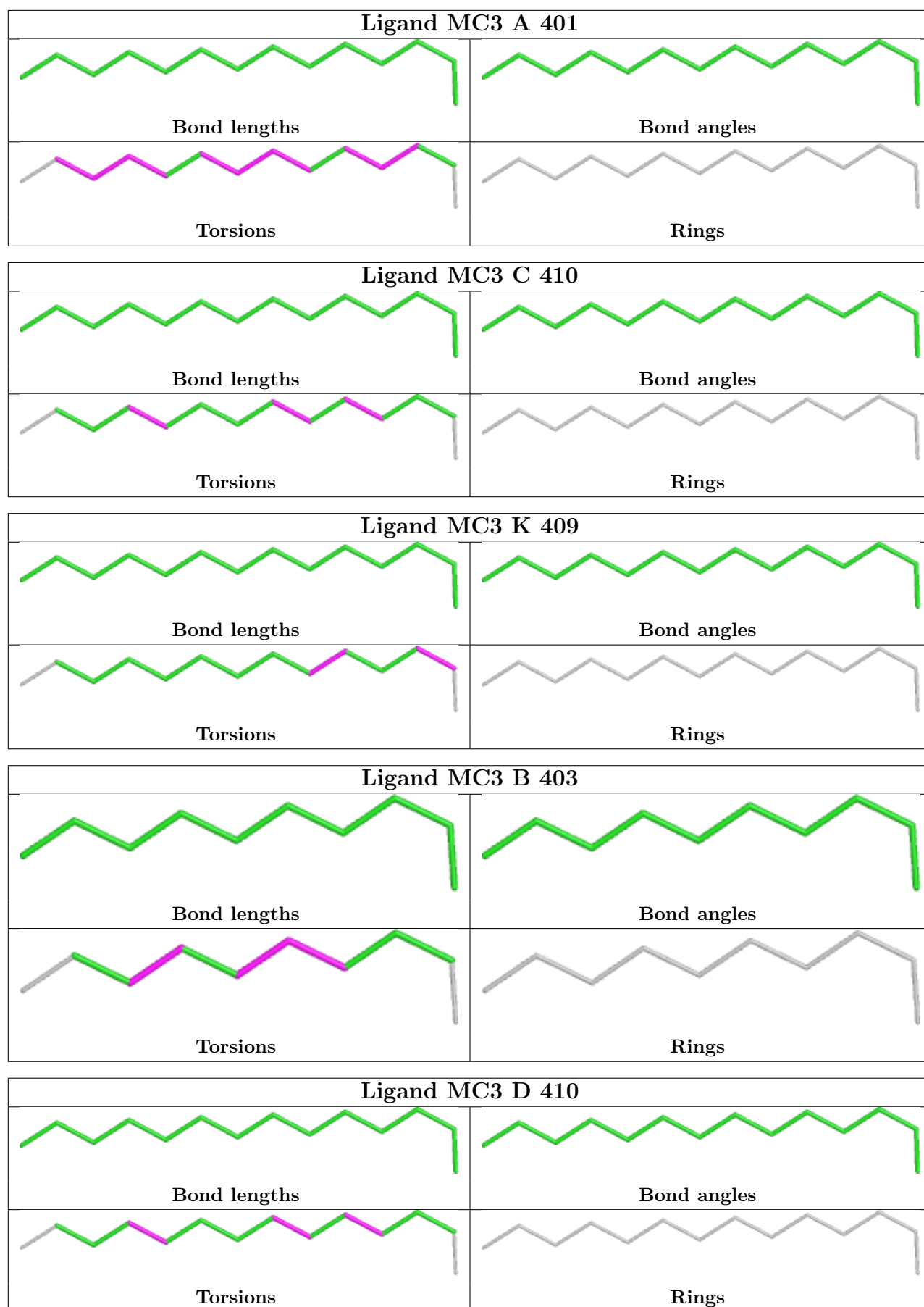
Ligand MC3 J 407	
 Bond lengths	 Bond angles
 Torsions	 Rings

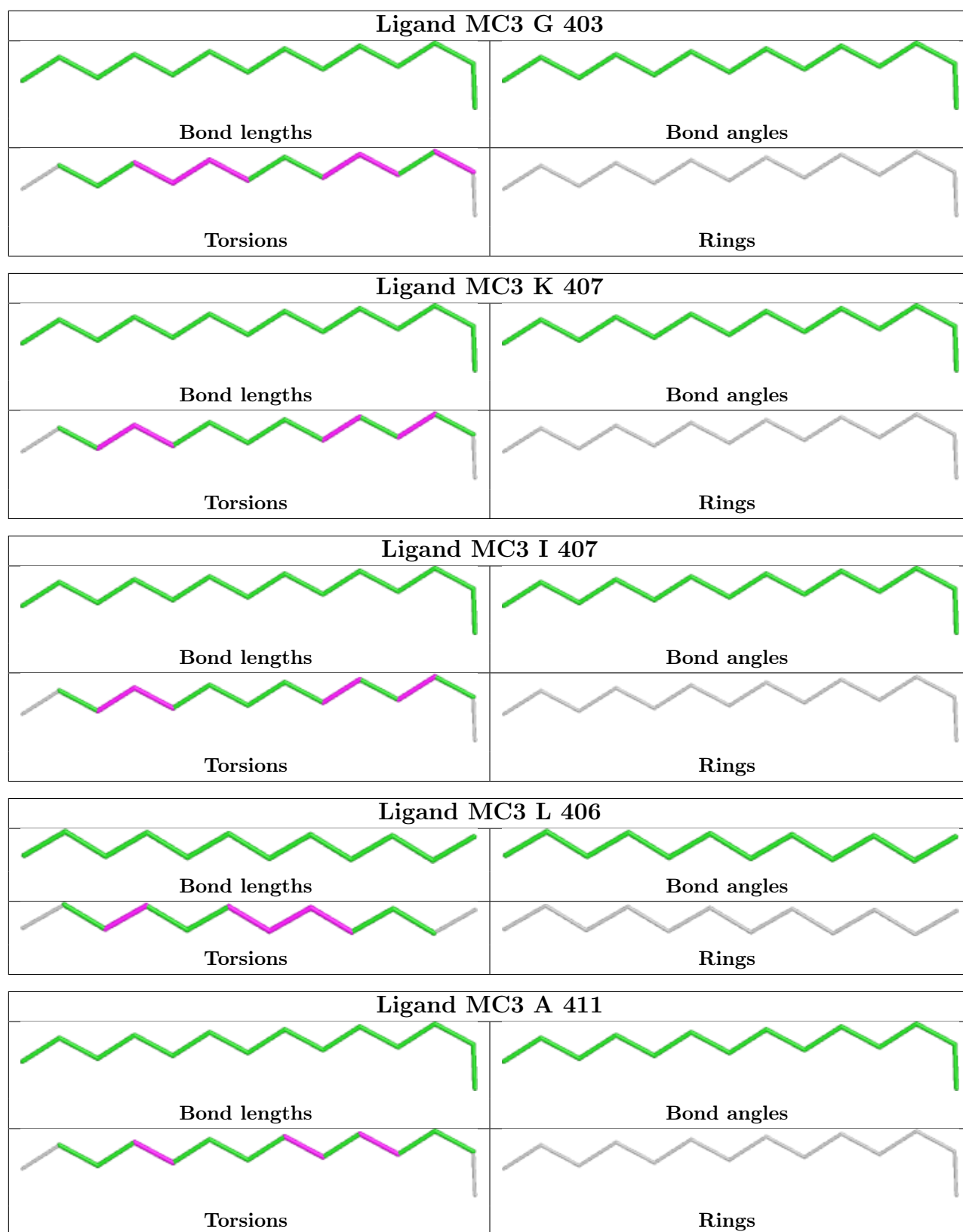


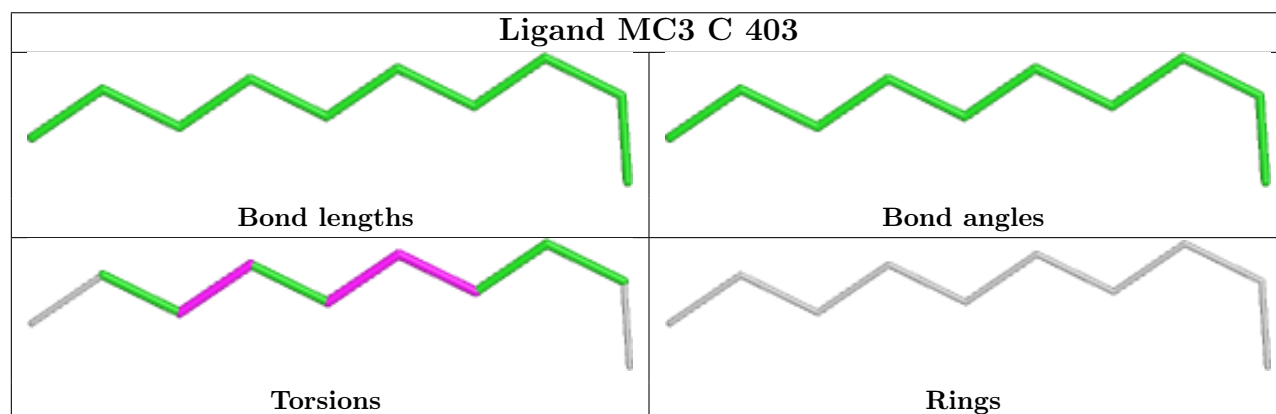
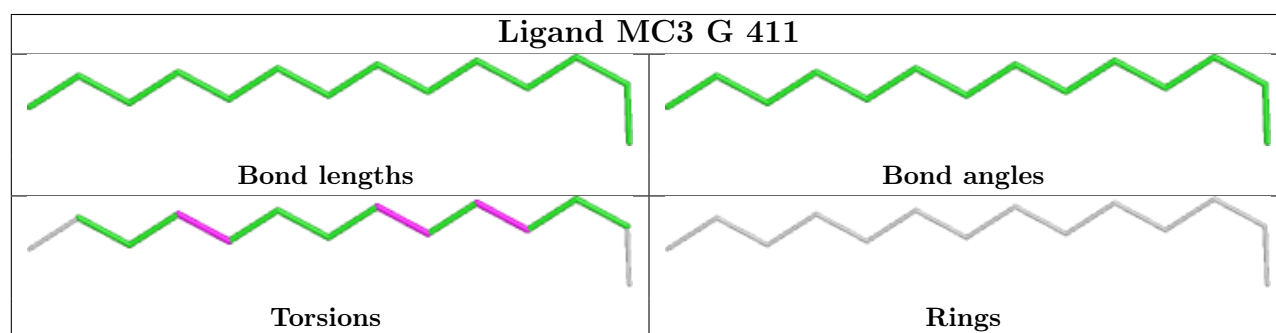
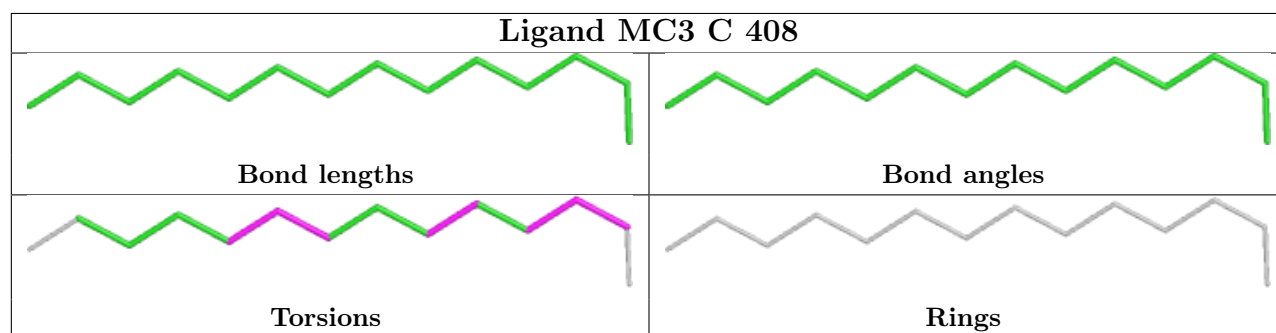
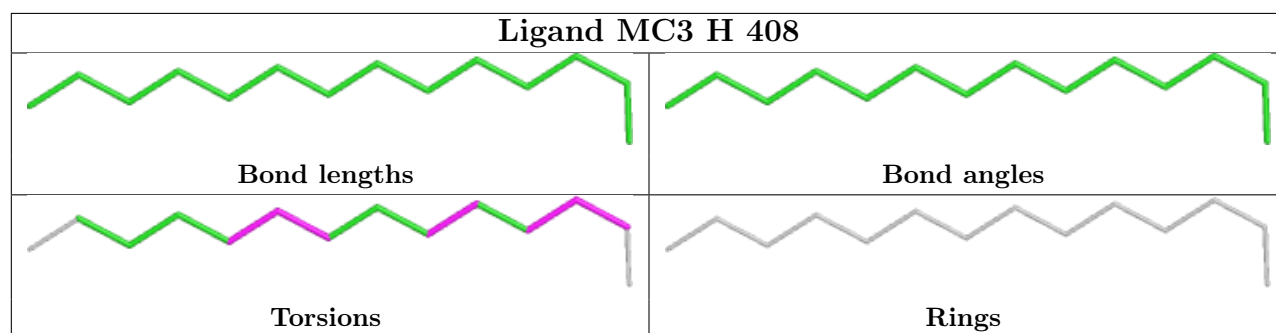
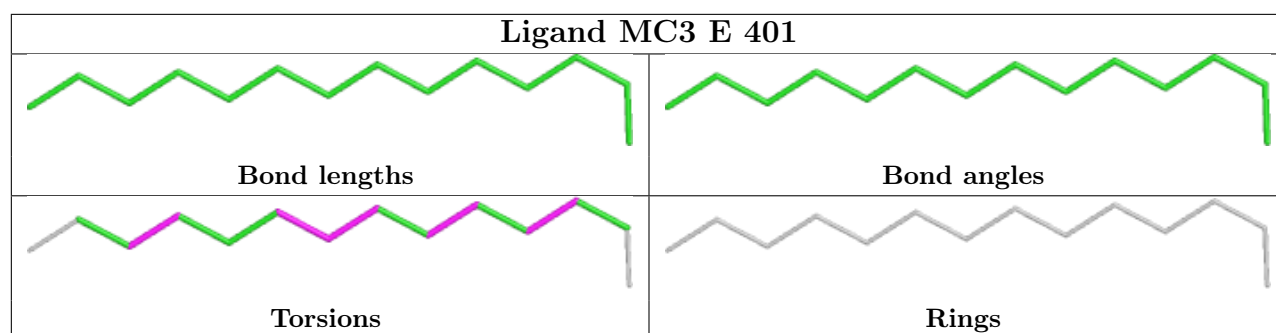


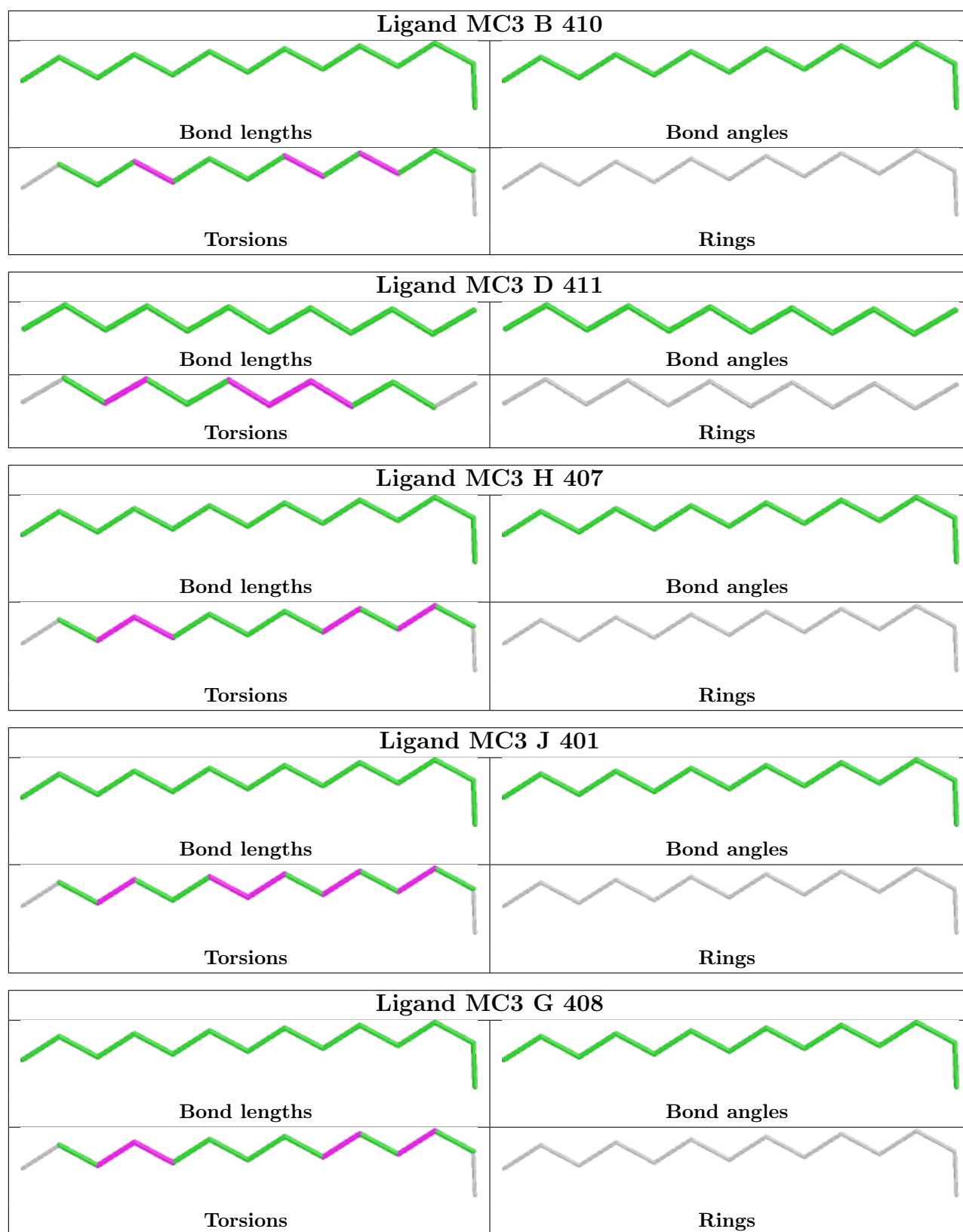


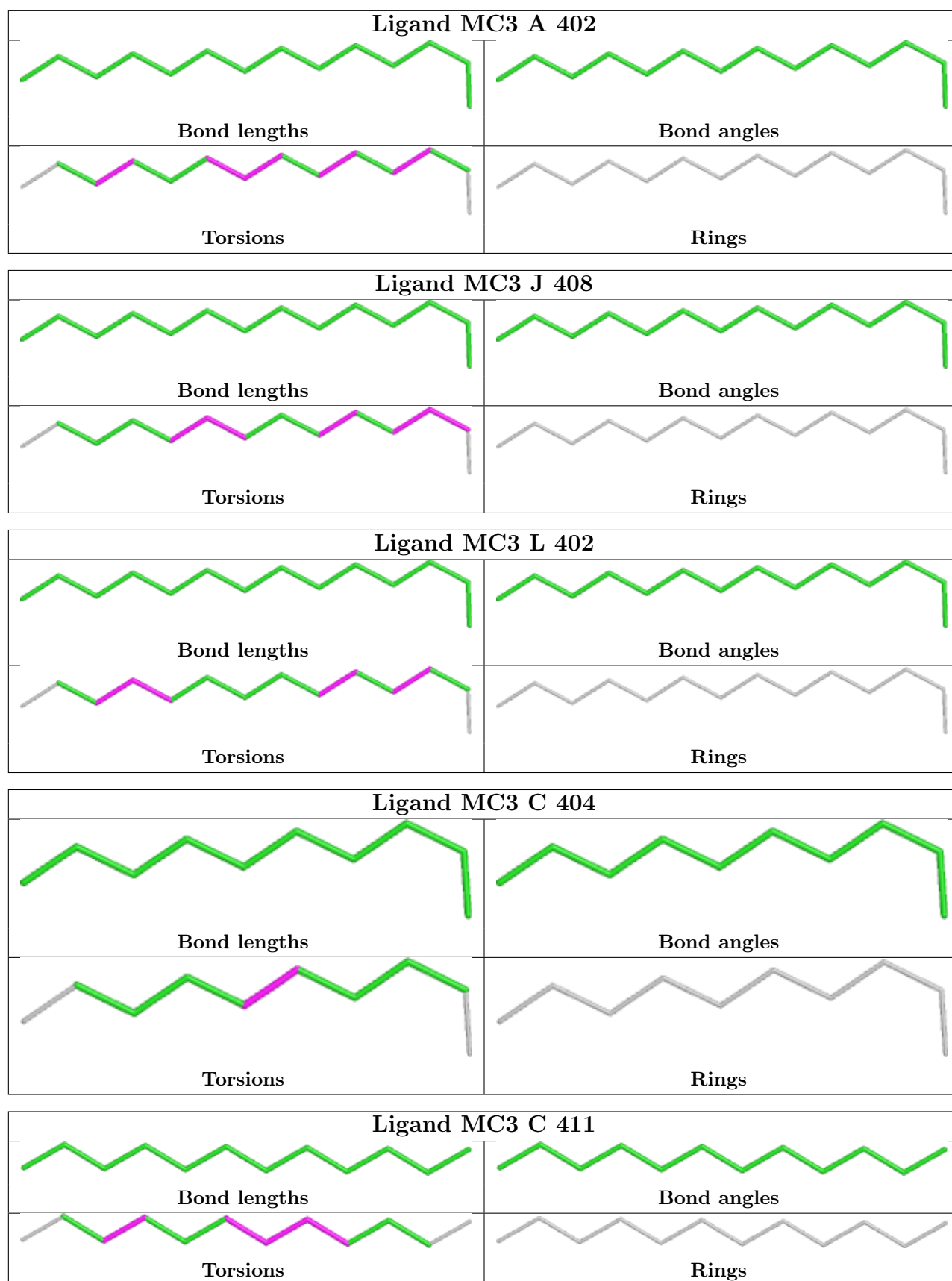


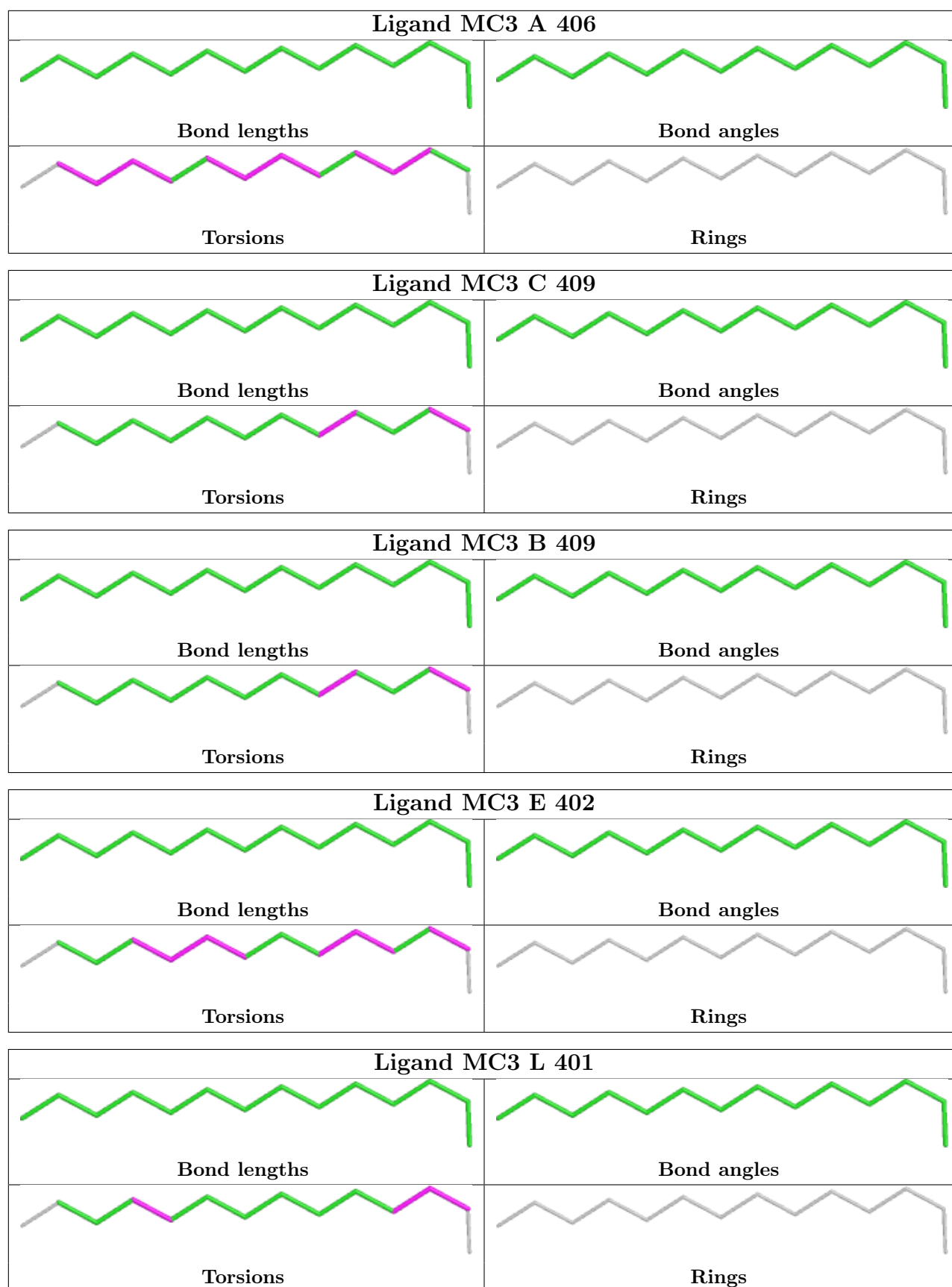


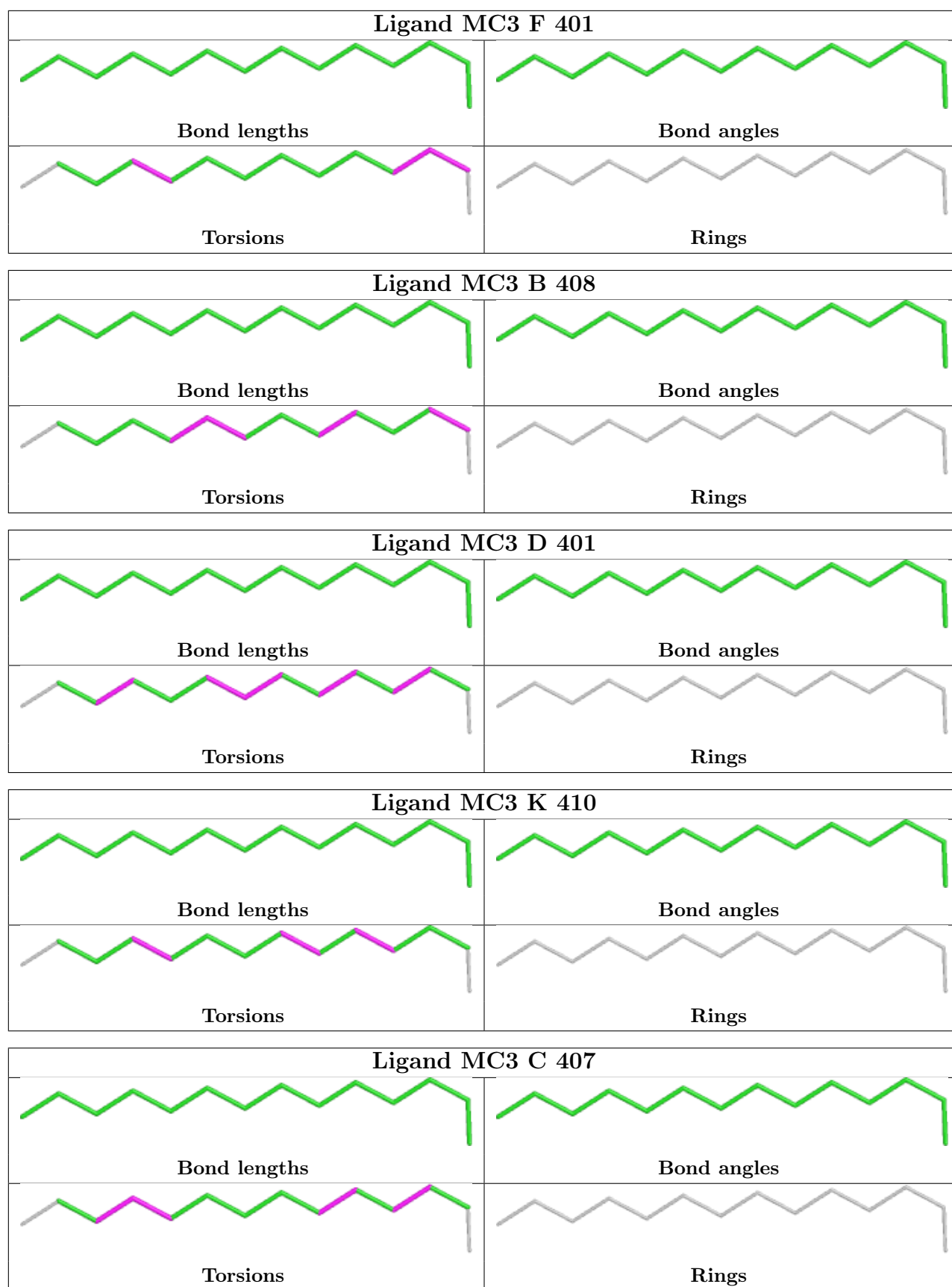


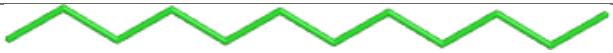
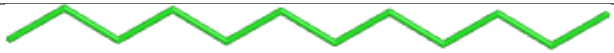










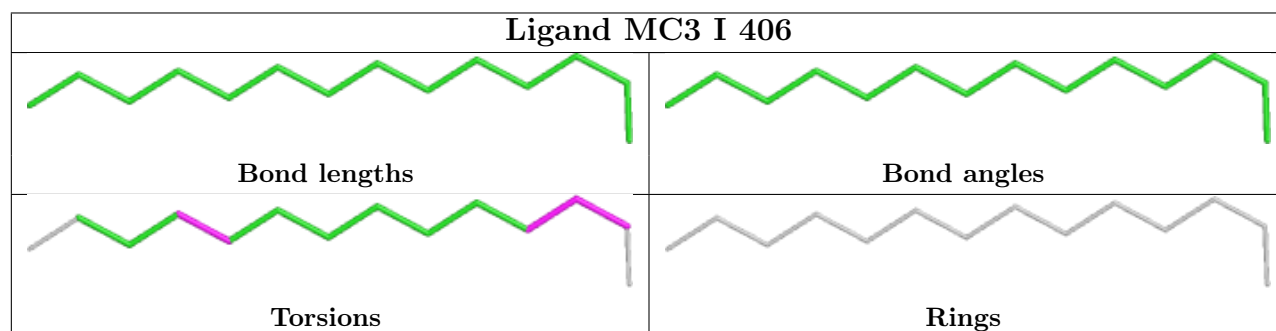
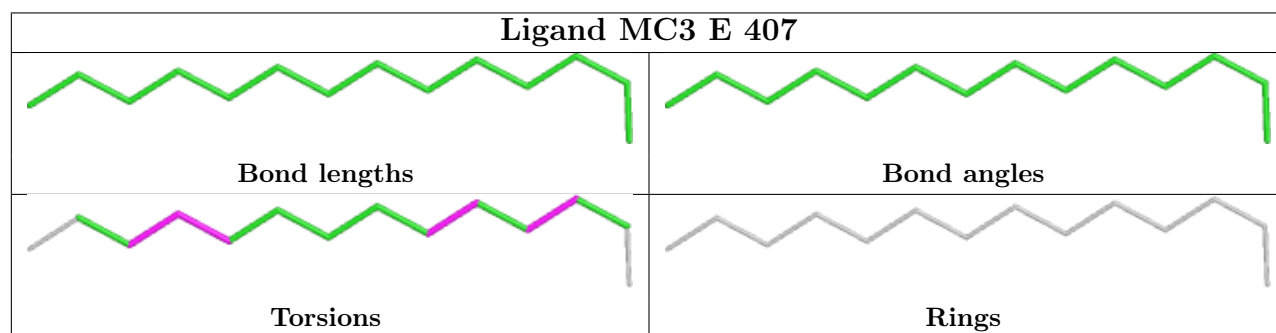
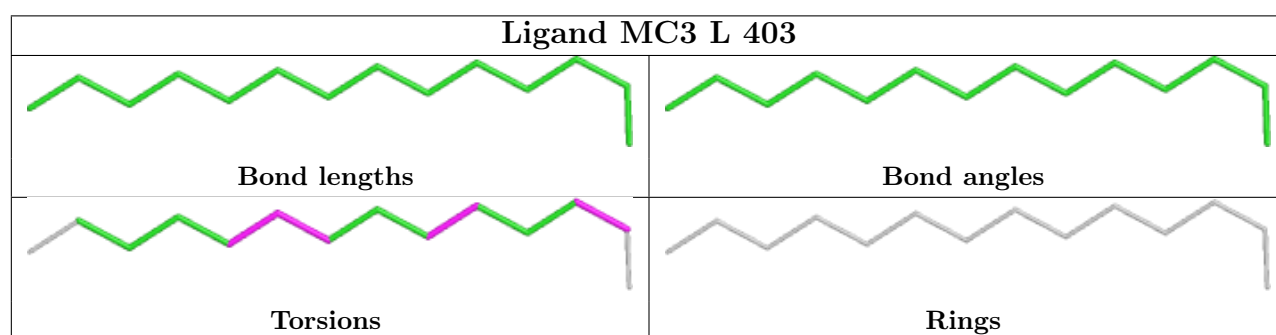
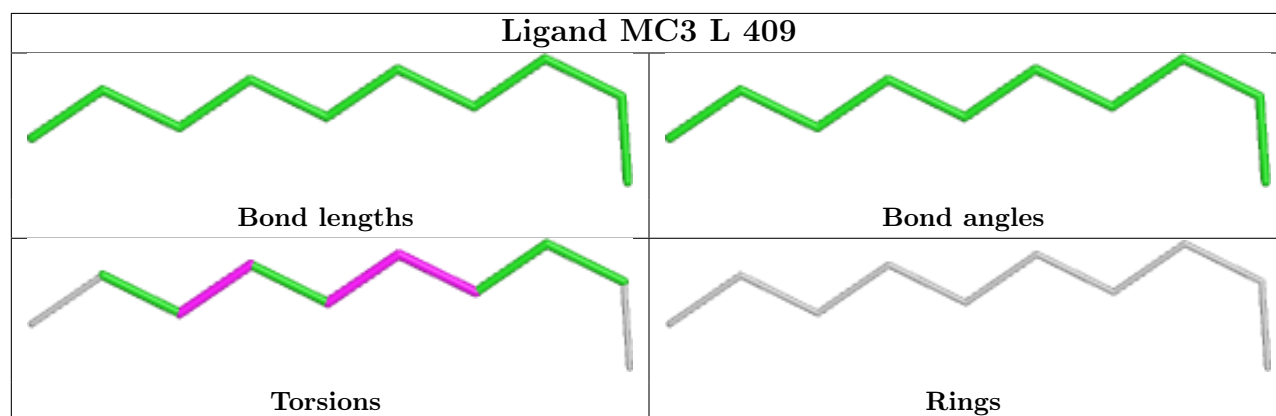
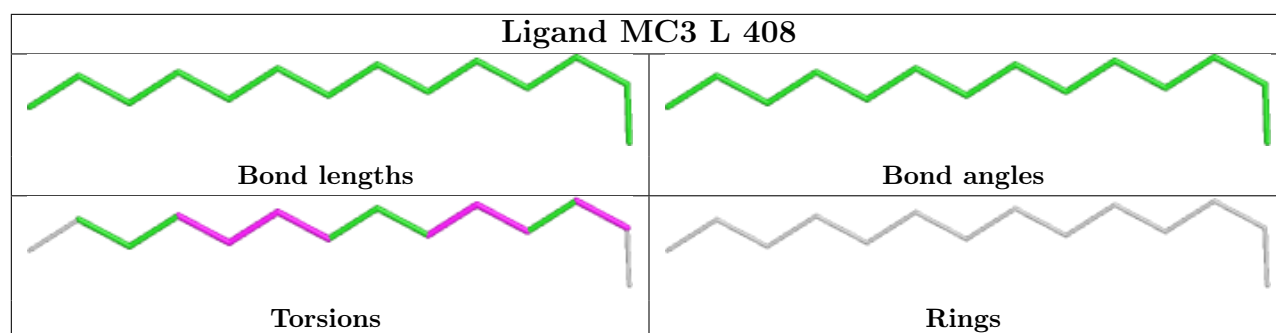
Ligand MC3 I 411	
 Bond lengths	 Bond angles
 Torsions	 Rings

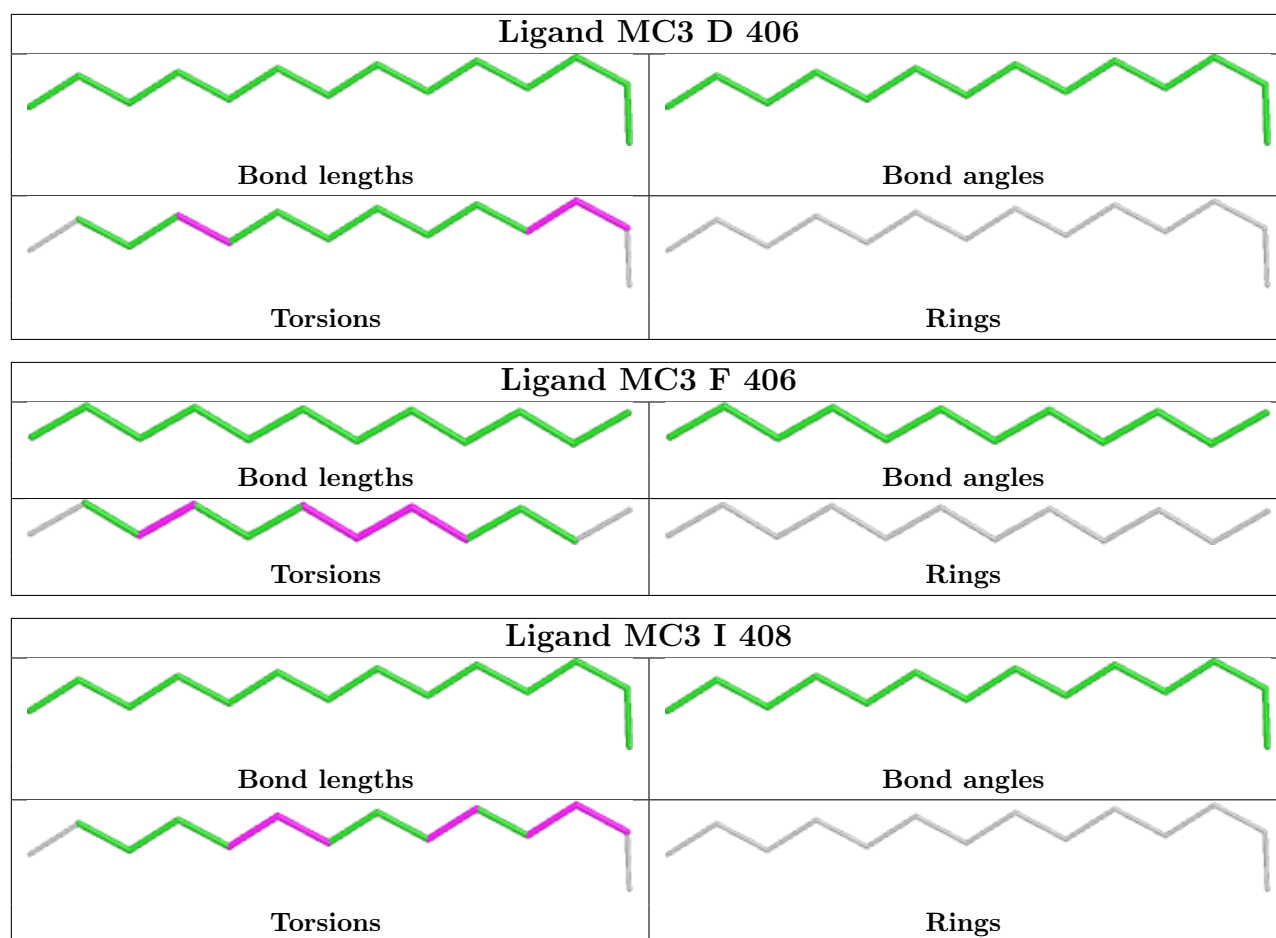
Ligand MC3 A 403	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand MC3 E 408	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand MC3 K 408	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand MC3 B 401	
 Bond lengths	 Bond angles
 Torsions	 Rings





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.

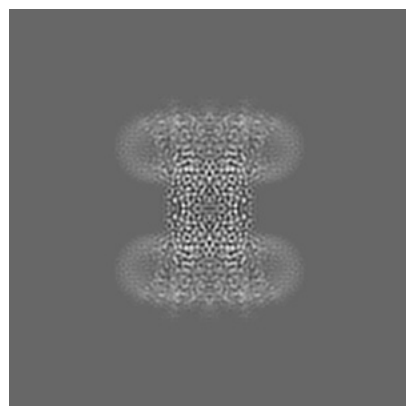
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-33315. These allow visual inspection of the internal detail of the map and identification of artifacts.

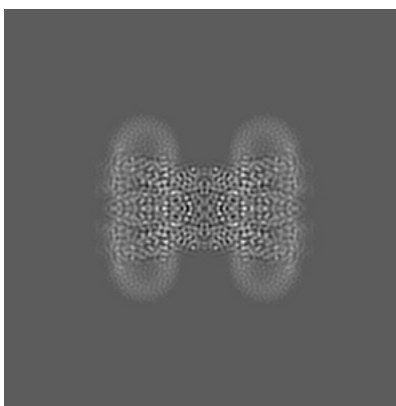
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

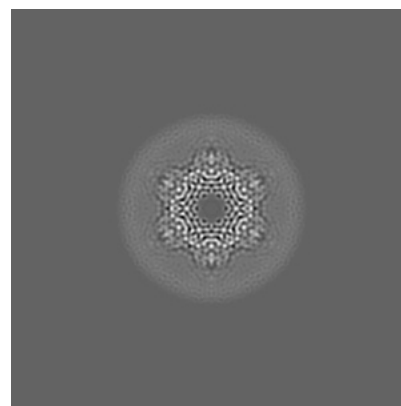
6.1.1 Primary map



X

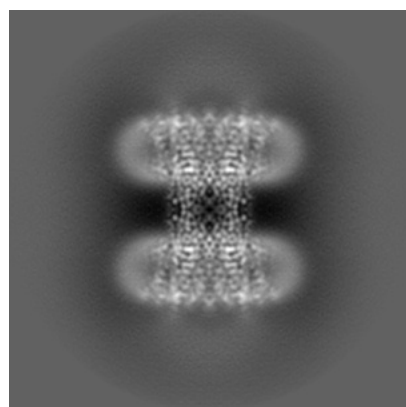


Y

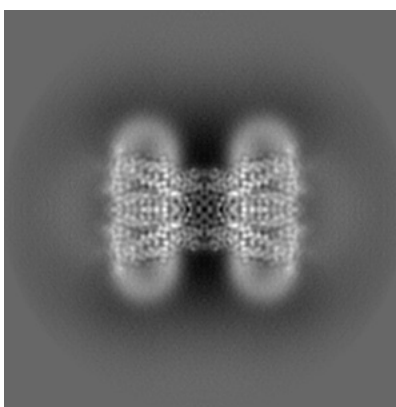


Z

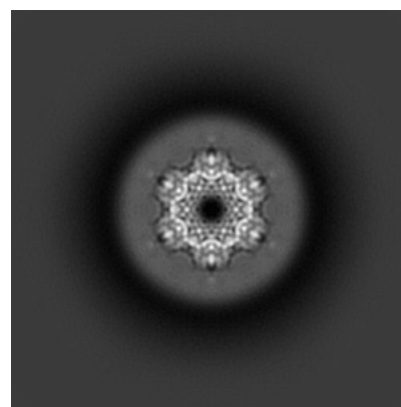
6.1.2 Raw map



X



Y

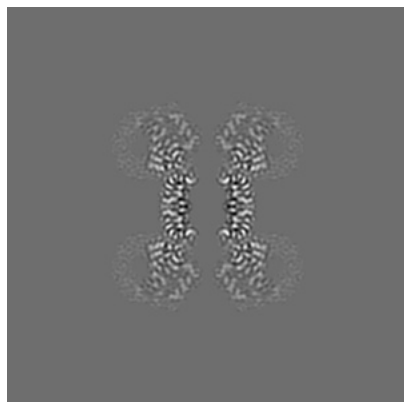


Z

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

6.2 Central slices [i](#)

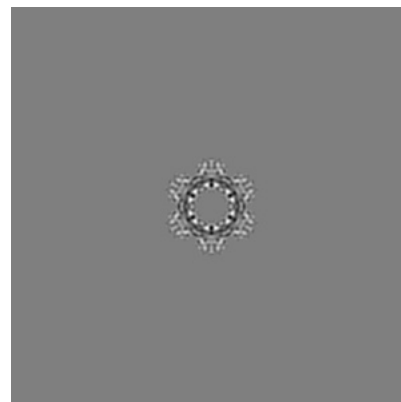
6.2.1 Primary map



X Index: 200

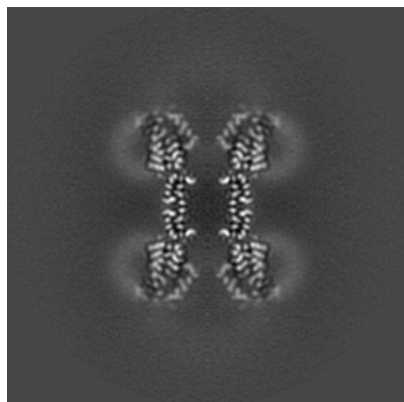


Y Index: 200

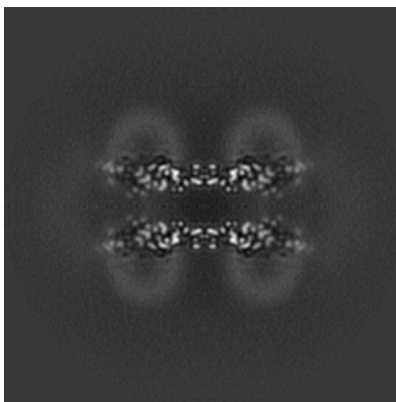


Z Index: 200

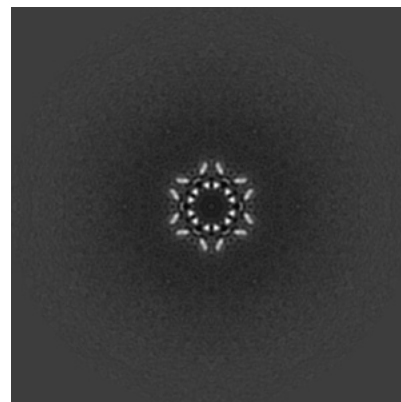
6.2.2 Raw map



X Index: 200



Y Index: 200

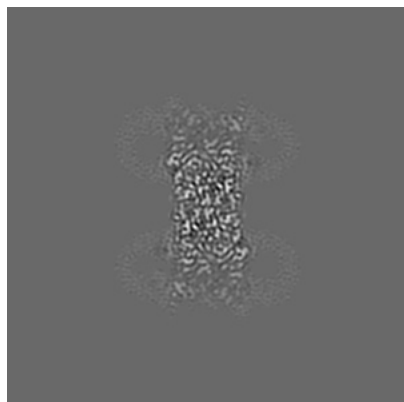


Z Index: 200

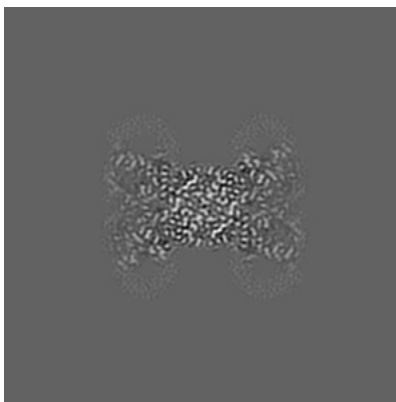
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

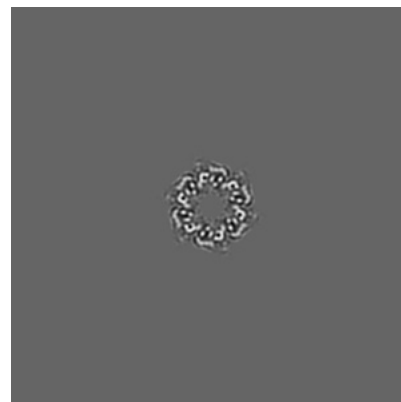
6.3.1 Primary map



X Index: 175

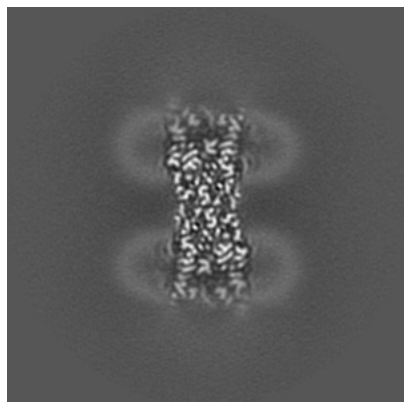


Y Index: 179

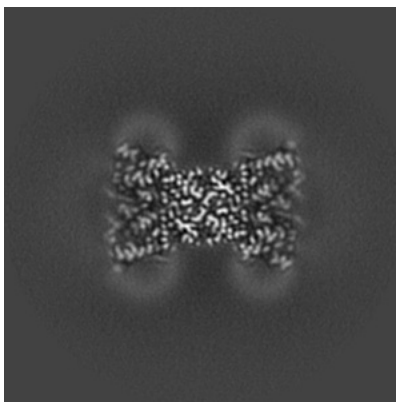


Z Index: 181

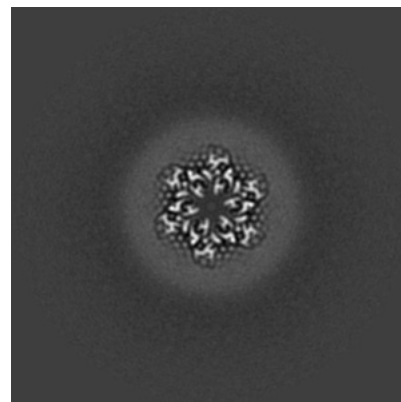
6.3.2 Raw map



X Index: 174



Y Index: 222

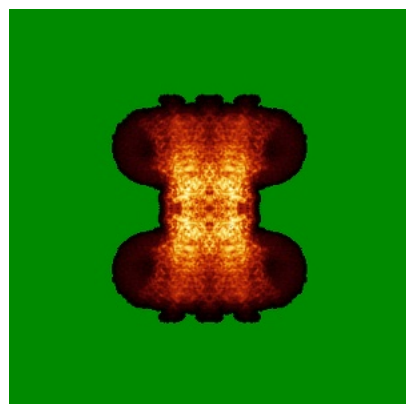


Z Index: 162

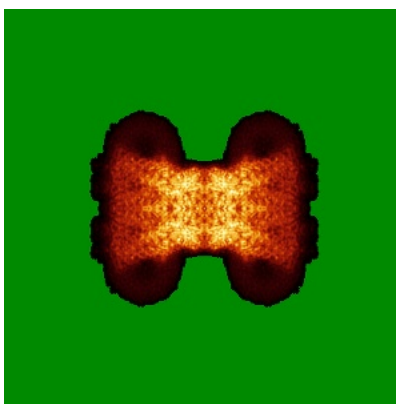
The images above show the largest variance slices of the map in three orthogonal directions.

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

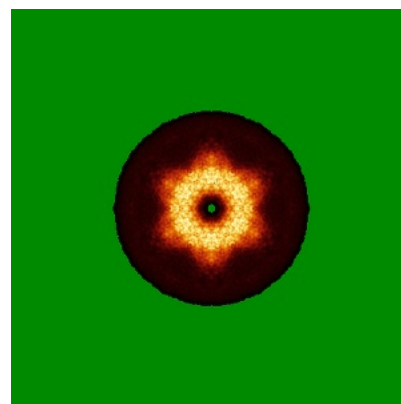
6.4.1 Primary map



X

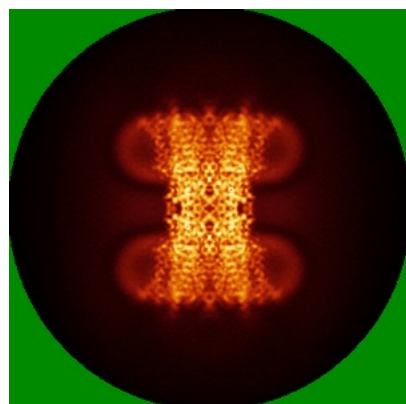


Y

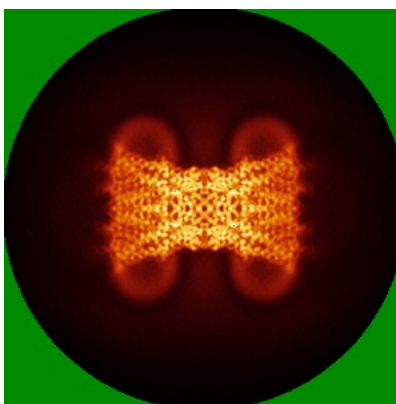


Z

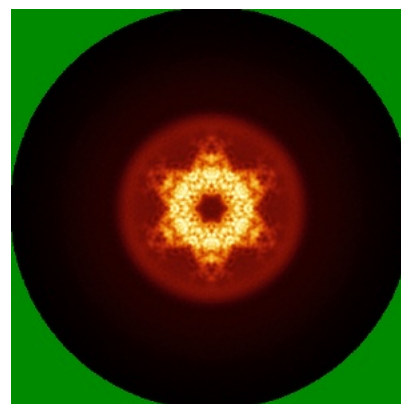
6.4.2 Raw map



X



Y

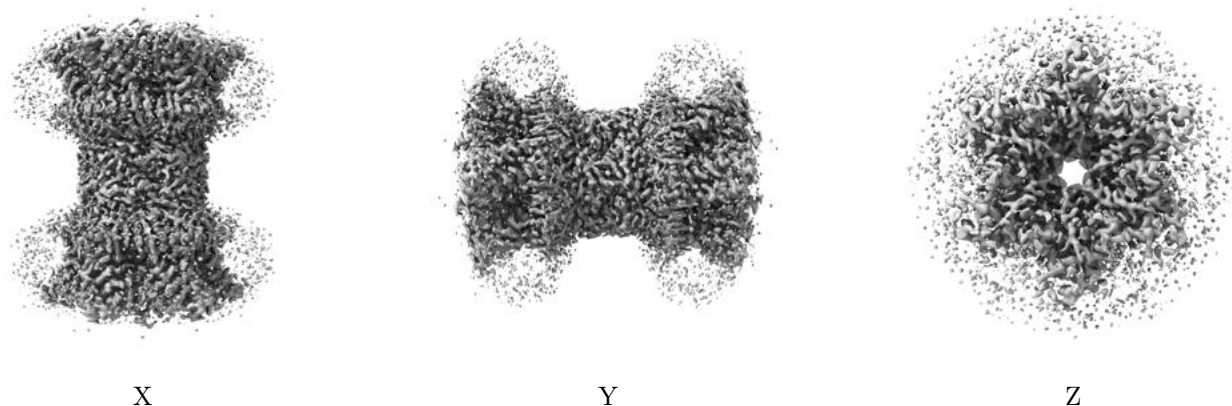


Z

The images above show the map standard deviation projections with false color in three orthogonal directions. Minimum values are shown in green, max in blue, and dark to light orange shades represent small to large values respectively.

6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.01. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

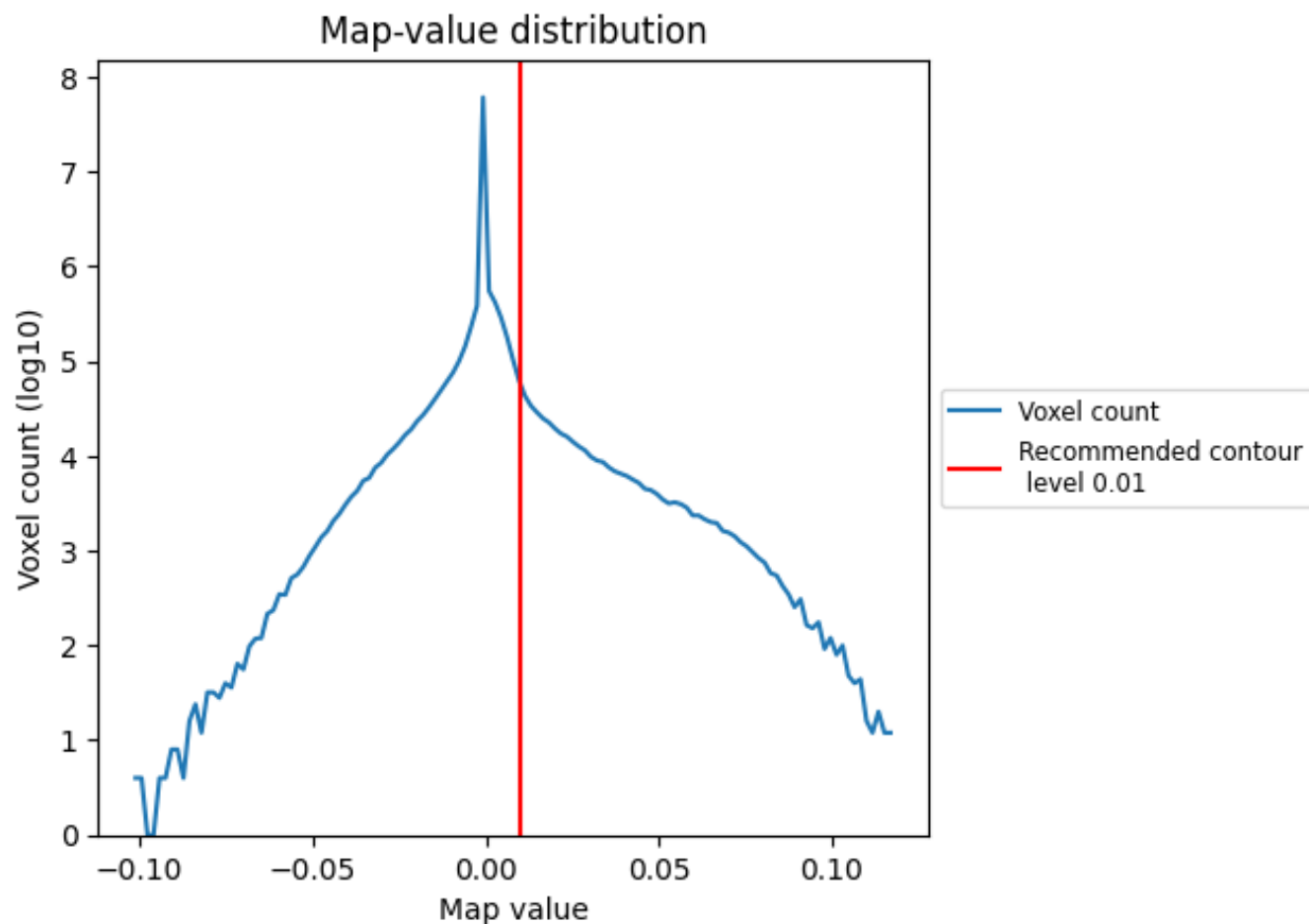
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

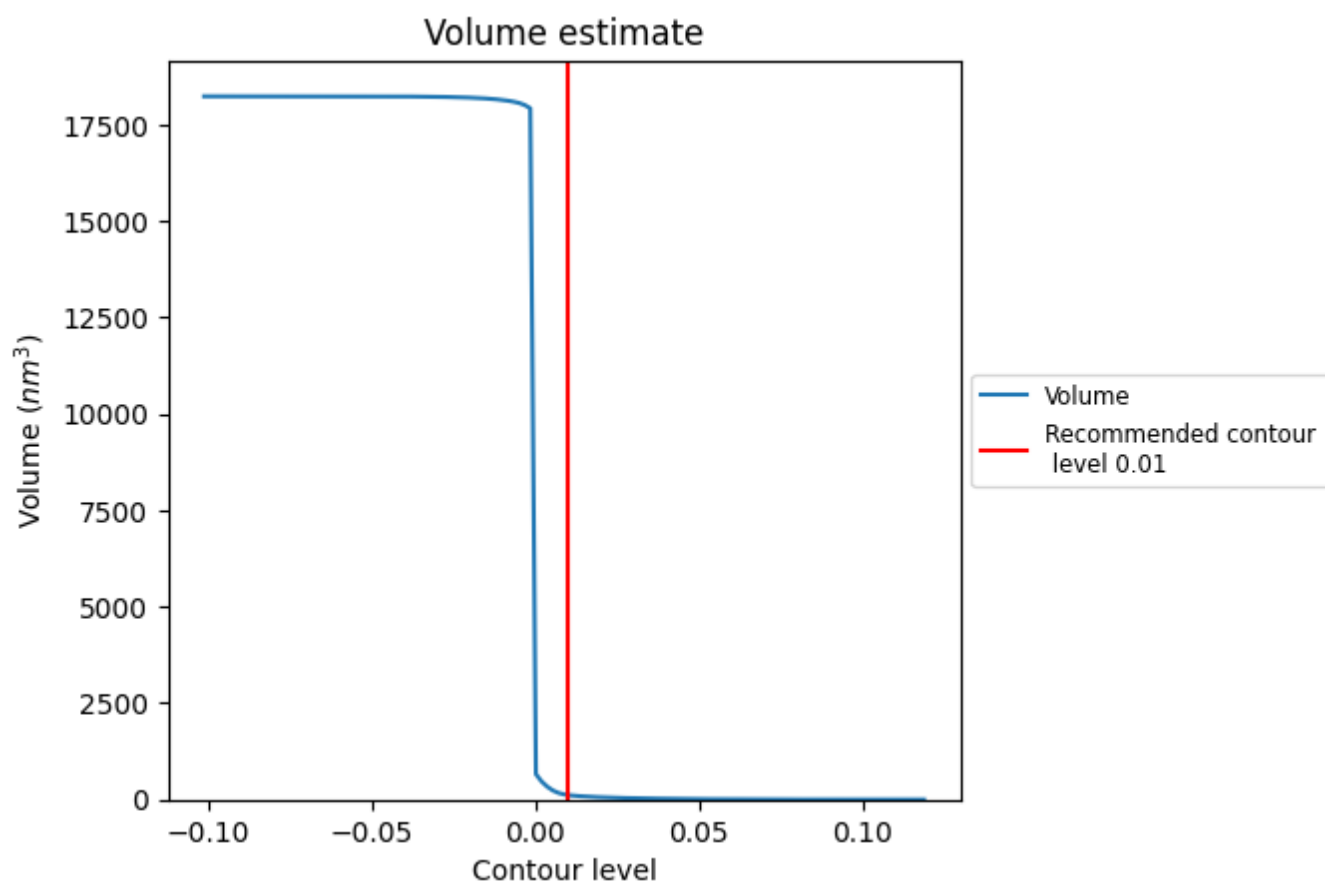
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

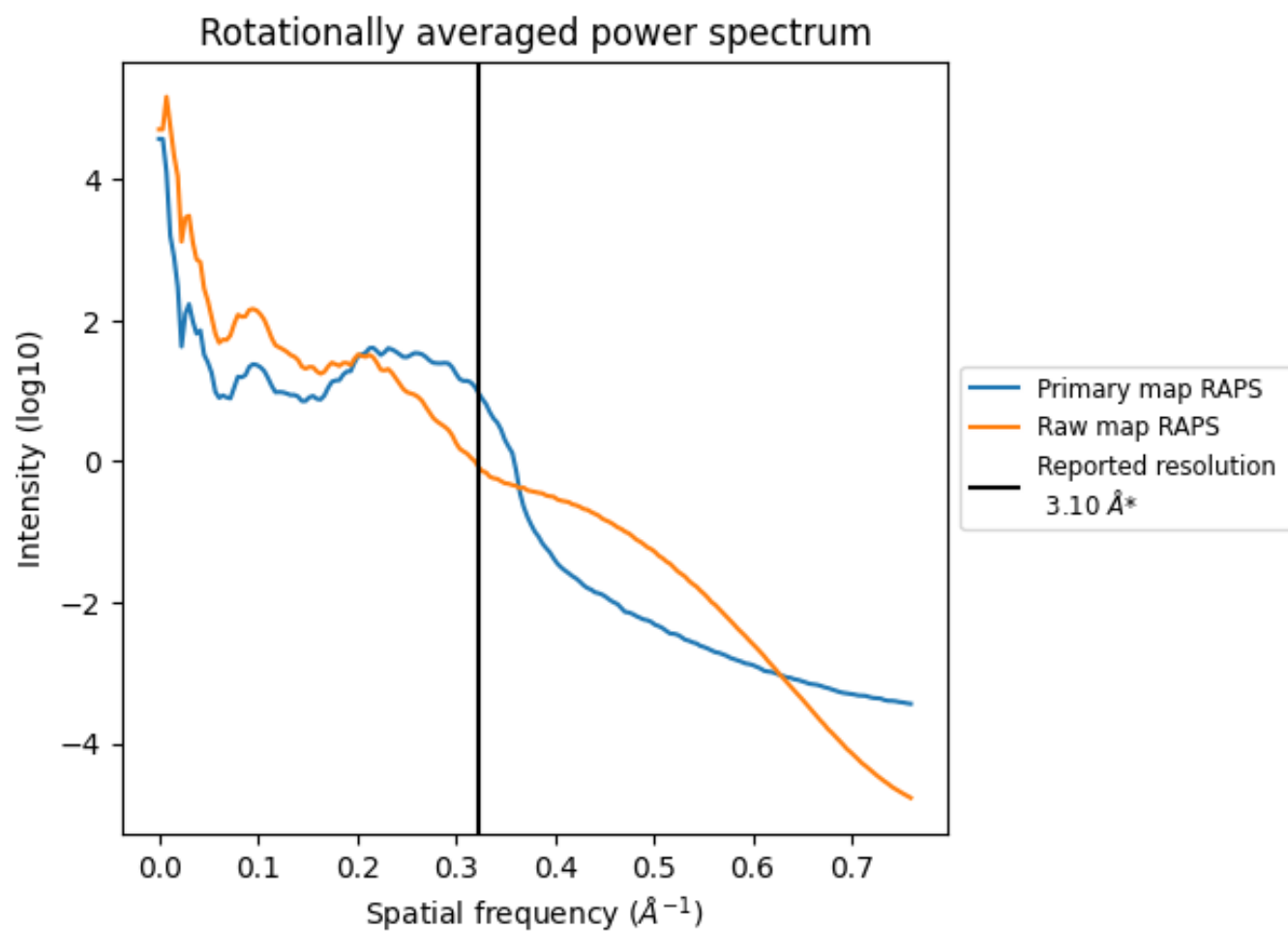
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 117 nm³; this corresponds to an approximate mass of 106 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ

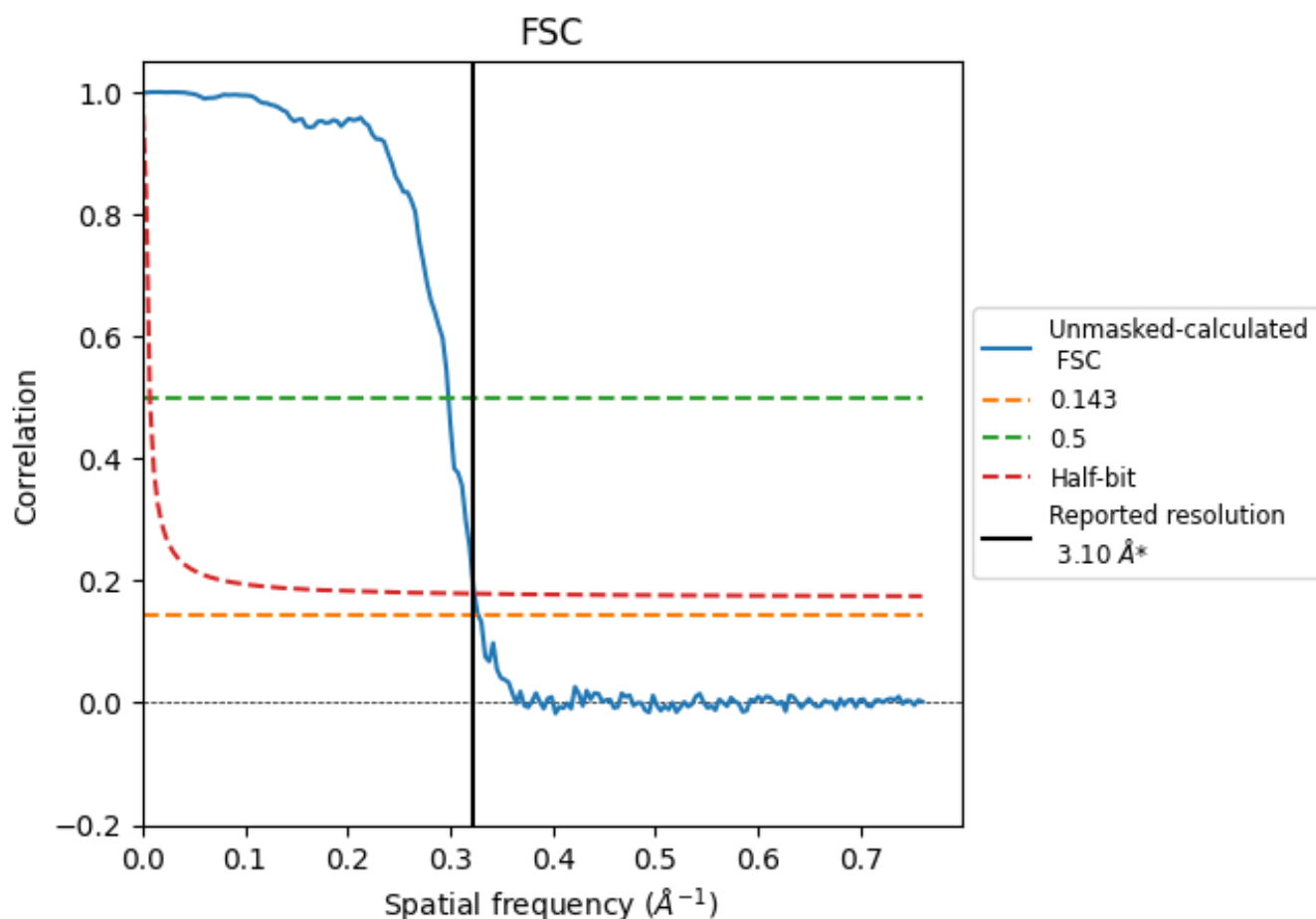


*Reported resolution corresponds to spatial frequency of $0.323 \text{ \AA}^{-1}$

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 0.323 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

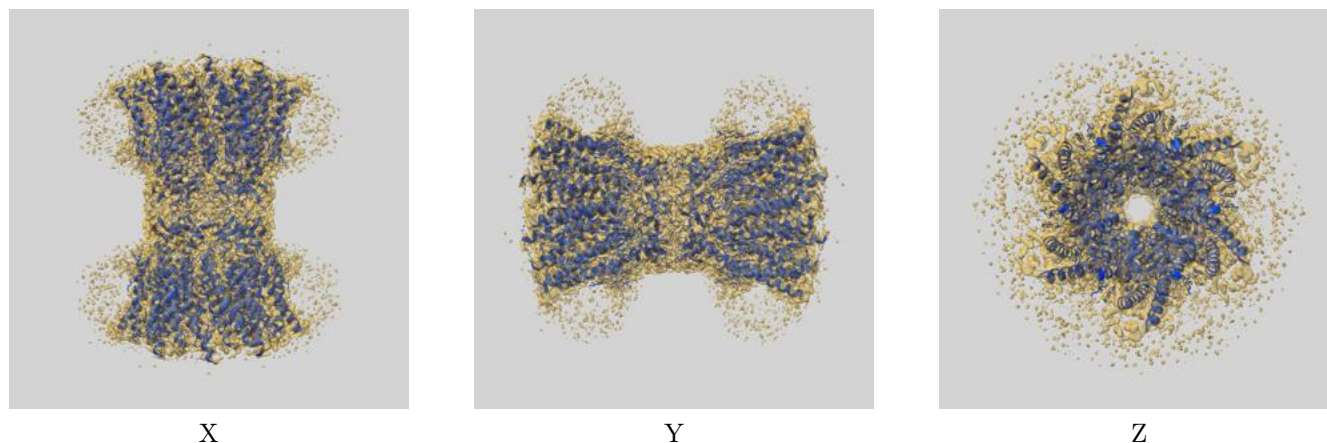
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.10	-	-
Author-provided FSC curve	-	-	-
Unmasked-calculated*	3.05	3.35	3.09

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

9 Map-model fit [i](#)

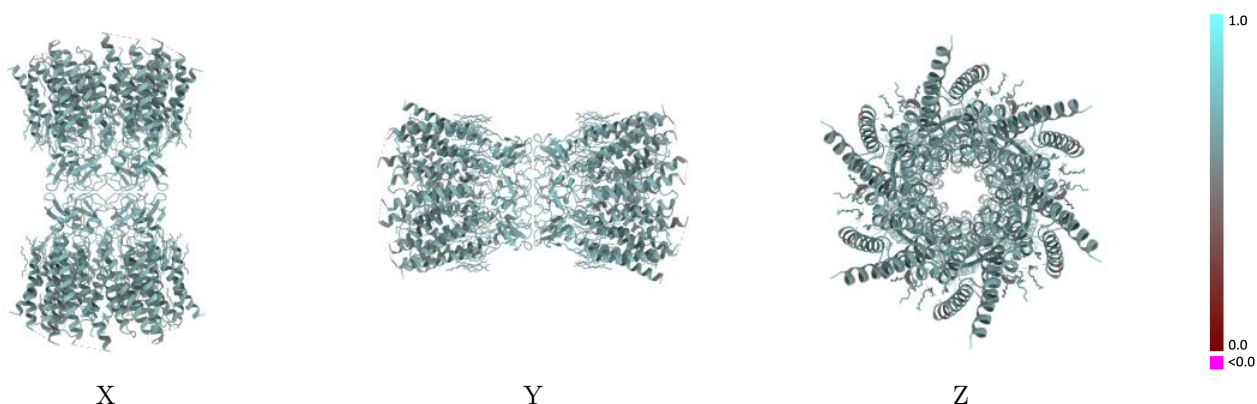
This section contains information regarding the fit between EMDB map EMD-33315 and PDB model 7XNH. Per-residue inclusion information can be found in section [3](#) on page [15](#).

9.1 Map-model overlay [i](#)



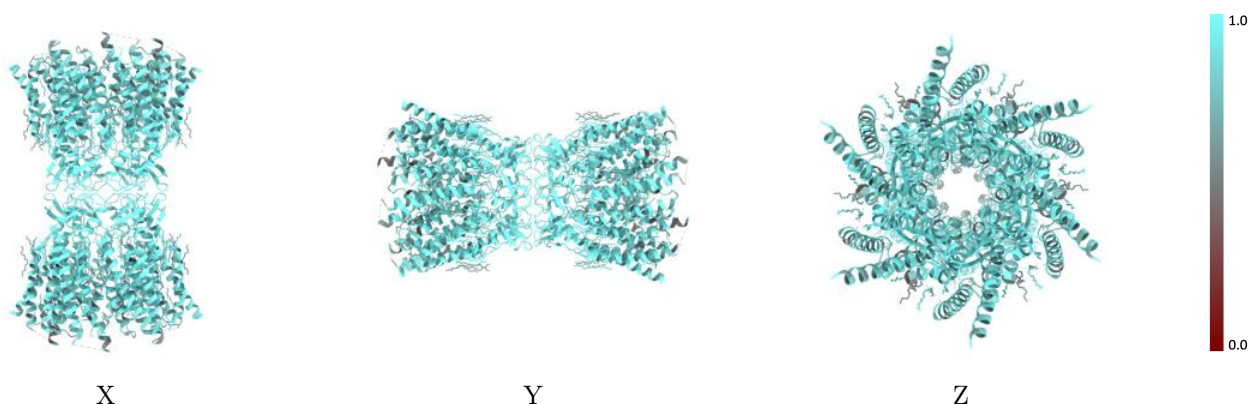
The images above show the 3D surface view of the map at the recommended contour level 0.01 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



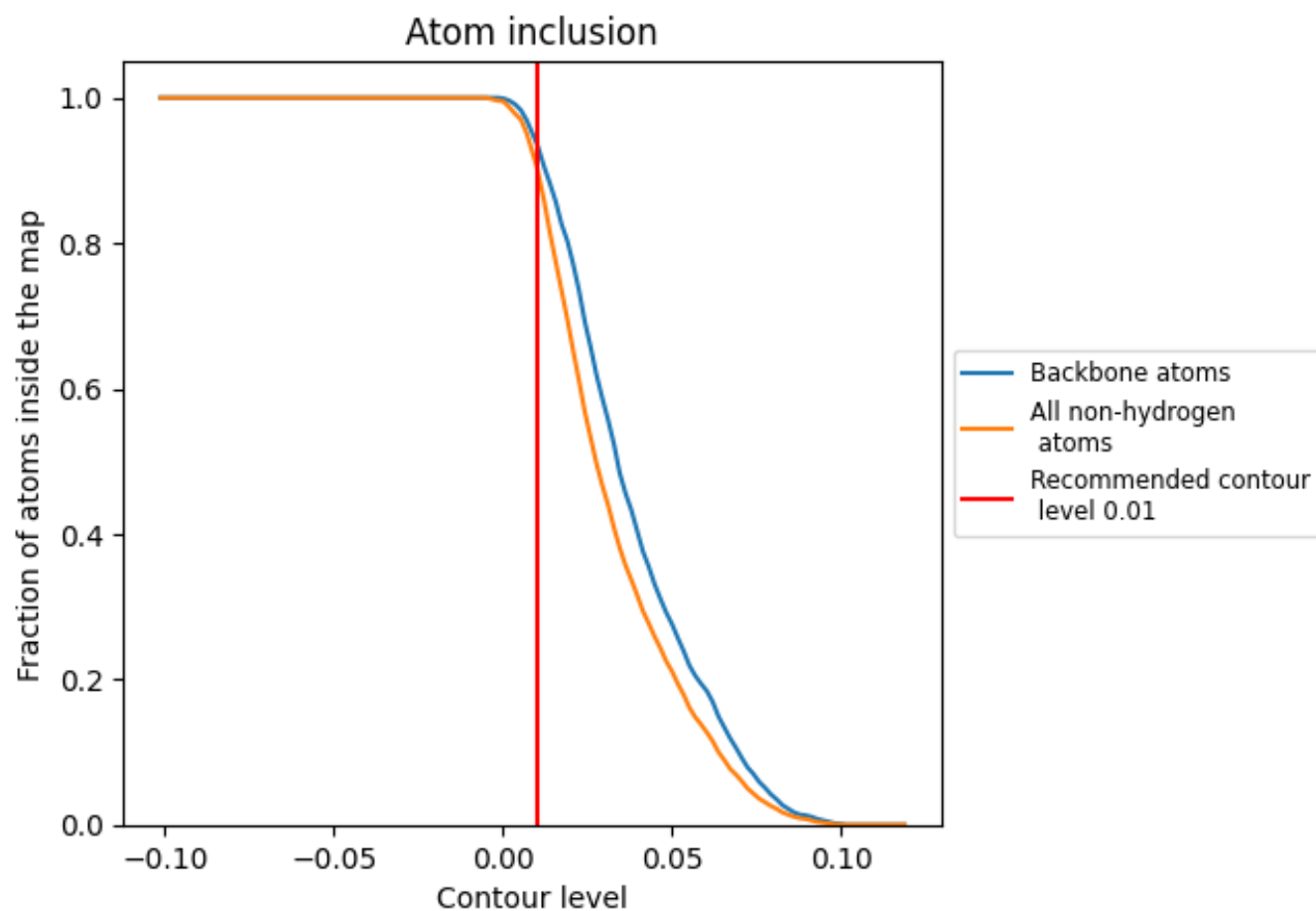
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.01).

9.4 Atom inclusion [i](#)



At the recommended contour level, 94% of all backbone atoms, 90% of all non-hydrogen atoms, are inside the map.

9.5 Map-model fit summary ⓘ

The table lists the average atom inclusion at the recommended contour level (0.01) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.9050	0.6230
A	0.9050	0.6230
B	0.9010	0.6220
C	0.9080	0.6210
D	0.9010	0.6210
E	0.9020	0.6240
F	0.9060	0.6240
G	0.9030	0.6220
H	0.9030	0.6220
I	0.9040	0.6240
J	0.9080	0.6230
K	0.9050	0.6230
L	0.9080	0.6240

1.0

0.0

<0.0