



wwPDB EM Validation Summary Report ⓘ

Mar 10, 2026 – 02:27 AM UTC

PDB ID : 6YEZ / pdb_00006yez
EMDB ID : EMD-10798
Title : Plant PSI-ferredoxin-plastocyanin supercomplex
Authors : Caspy, I.; Nelson, N.; Shkolnisky, Y.; Klaiman, D.; Sheinker, A.
Deposited on : 2020-03-25
Resolution : 2.70 Å(reported)
Based on initial model : 5L8R

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

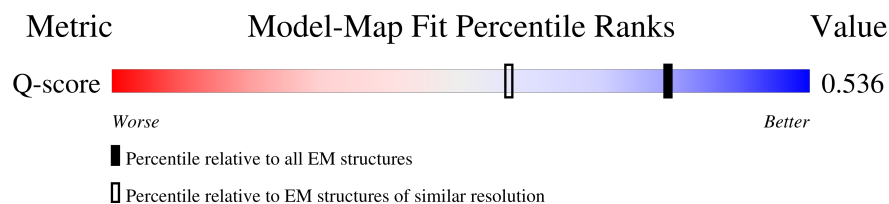
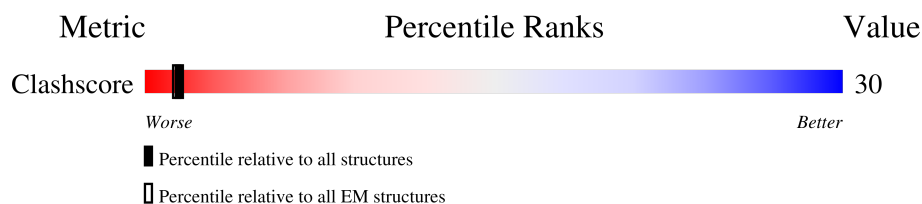
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

ELECTRON MICROSCOPY

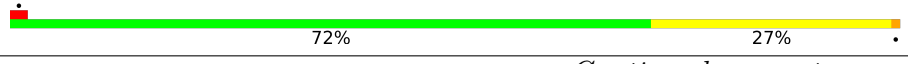
The reported resolution of this entry is 2.70 Å.

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.



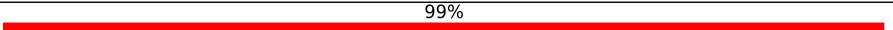
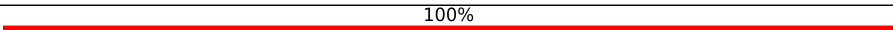
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Q-score	-	25397	10327 (2.20 - 3.20)

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	743	 67% 32% .
2	B	733	 72% 28% .
3	C	80	 66% 32% .
4	D	143	 73% 27% .
5	E	66	 5% 88% 12%
6	F	154	 72% 27% .

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Mol	Chain	Length	Quality of chain
7	G	97	
8	H	93	
9	I	31	
10	J	42	
11	K	81	
12	L	159	
13	1	193	
14	2	208	
15	3	221	
16	4	198	
17	N	97	
18	P	99	

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
19	CL0	A	1011	X	-	-	-
20	CLA	1	601	X	-	-	-
20	CLA	1	602	X	-	-	-
20	CLA	1	603	X	-	X	-
20	CLA	1	604	X	-	-	-
20	CLA	1	605	X	-	-	-
20	CLA	1	606	X	-	-	-
20	CLA	1	607	X	-	-	-
20	CLA	1	608	X	-	-	-
20	CLA	1	611	X	-	-	-
20	CLA	1	613	X	-	-	-
20	CLA	1	614	X	-	-	-
20	CLA	2	601	X	-	-	-
20	CLA	2	602	X	-	-	-
20	CLA	2	603	X	-	-	-
20	CLA	2	604	X	-	X	-
20	CLA	2	605	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
20	CLA	2	606	X	-	-	-
20	CLA	2	607	X	-	-	-
20	CLA	2	608	X	-	-	-
20	CLA	2	612	X	-	-	-
20	CLA	3	601	X	-	-	-
20	CLA	3	602	X	-	-	-
20	CLA	3	603	X	-	-	-
20	CLA	3	605	X	-	-	-
20	CLA	3	606	X	-	-	-
20	CLA	3	608	X	-	-	-
20	CLA	3	610	X	-	-	-
20	CLA	3	612	X	-	-	-
20	CLA	3	613	X	-	-	-
20	CLA	3	614	X	-	-	-
20	CLA	3	617	X	-	-	-
20	CLA	4	601	X	-	X	-
20	CLA	4	602	X	-	-	-
20	CLA	4	603	X	-	-	-
20	CLA	4	604	X	-	X	-
20	CLA	4	606	X	-	-	-
20	CLA	4	607	X	-	-	-
20	CLA	4	608	X	-	-	-
20	CLA	4	609	X	-	-	-
20	CLA	4	612	X	-	-	-
20	CLA	4	617	X	-	-	-
20	CLA	A	1101	X	-	-	-
20	CLA	A	1102	X	-	-	-
20	CLA	A	1103	X	-	-	-
20	CLA	A	1104	X	-	-	-
20	CLA	A	1105	X	-	-	-
20	CLA	A	1106	X	-	-	-
20	CLA	A	1107	X	-	-	-
20	CLA	A	1108	X	-	-	-
20	CLA	A	1109	X	-	-	-
20	CLA	A	1110	X	-	-	-
20	CLA	A	1111	X	-	-	-
20	CLA	A	1112	X	-	-	-
20	CLA	A	1113	X	-	-	-
20	CLA	A	1114	X	-	-	-
20	CLA	A	1115	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
20	CLA	A	1116	X	-	-	-
20	CLA	A	1117	X	-	-	-
20	CLA	A	1118	X	-	-	-
20	CLA	A	1119	X	-	-	-
20	CLA	A	1120	X	-	-	-
20	CLA	A	1121	X	-	-	-
20	CLA	A	1122	X	-	-	-
20	CLA	A	1123	X	-	-	-
20	CLA	A	1124	X	-	-	-
20	CLA	A	1125	X	-	-	-
20	CLA	A	1126	X	-	-	-
20	CLA	A	1127	X	-	-	-
20	CLA	A	1128	X	-	-	-
20	CLA	A	1129	X	-	-	-
20	CLA	A	1130	X	-	-	-
20	CLA	A	1131	X	-	-	-
20	CLA	A	1132	X	-	X	-
20	CLA	A	1133	X	-	-	-
20	CLA	A	1134	X	-	-	-
20	CLA	A	1135	X	-	-	-
20	CLA	A	1136	X	-	-	-
20	CLA	A	1137	X	-	-	-
20	CLA	A	1138	X	-	-	-
20	CLA	A	1139	X	-	-	-
20	CLA	A	1140	X	-	-	-
20	CLA	A	1141	X	-	-	-
20	CLA	B	1021	X	-	-	-
20	CLA	B	1022	X	-	-	-
20	CLA	B	1023	X	-	-	-
20	CLA	B	1201	X	-	-	-
20	CLA	B	1202	X	-	-	-
20	CLA	B	1203	X	-	-	-
20	CLA	B	1204	X	-	-	-
20	CLA	B	1205	X	-	-	-
20	CLA	B	1206	X	-	-	-
20	CLA	B	1207	X	-	-	-
20	CLA	B	1208	X	-	-	-
20	CLA	B	1209	X	-	-	-
20	CLA	B	1210	X	-	-	-
20	CLA	B	1211	X	-	-	-
20	CLA	B	1212	X	-	-	-
20	CLA	B	1213	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
20	CLA	B	1214	X	-	-	-
20	CLA	B	1215	X	-	-	-
20	CLA	B	1216	X	-	-	-
20	CLA	B	1217	X	-	-	-
20	CLA	B	1218	X	-	-	-
20	CLA	B	1219	X	-	-	-
20	CLA	B	1220	X	-	-	-
20	CLA	B	1221	X	-	-	-
20	CLA	B	1222	X	-	-	-
20	CLA	B	1223	X	-	-	-
20	CLA	B	1224	X	-	-	-
20	CLA	B	1225	X	-	-	-
20	CLA	B	1226	X	-	-	-
20	CLA	B	1227	X	-	-	-
20	CLA	B	1228	X	-	-	-
20	CLA	B	1229	X	-	-	-
20	CLA	B	1230	X	-	-	-
20	CLA	B	1231	X	-	-	-
20	CLA	B	1232	X	-	-	-
20	CLA	B	1234	X	-	-	-
20	CLA	B	1235	X	-	-	-
20	CLA	B	1236	X	-	-	-
20	CLA	B	1237	X	-	-	-
20	CLA	B	1238	X	-	-	-
20	CLA	B	1239	X	-	-	-
20	CLA	B	1240	X	-	-	-
20	CLA	F	1301	X	-	-	-
20	CLA	F	1302	X	-	-	-
20	CLA	G	1601	X	-	-	-
20	CLA	G	1602	X	-	-	-
20	CLA	G	1603	X	-	-	-
20	CLA	H	1701	X	-	-	-
20	CLA	J	1901	X	-	-	-
20	CLA	K	1401	X	-	-	-
20	CLA	K	1402	X	-	-	-
20	CLA	K	1403	X	-	-	-
20	CLA	K	1404	X	-	-	-
20	CLA	L	1501	X	-	-	-
20	CLA	L	1502	X	-	-	-
20	CLA	L	1503	X	-	-	-
22	BCR	2	503	-	-	X	-
22	BCR	L	4019	-	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
25	SF4	C	3003	-	-	X	-
29	LUT	1	502	-	-	X	-
29	LUT	2	501	X	-	-	-
29	LUT	3	502	X	-	-	-
30	CHL	1	609	X	-	-	-
30	CHL	1	610	X	-	-	-
30	CHL	1	612	X	-	-	-
30	CHL	2	609	X	-	X	-
30	CHL	2	610	X	-	-	-
30	CHL	2	611	X	-	-	-
30	CHL	2	613	X	-	-	-
30	CHL	2	615	X	-	-	-
30	CHL	3	604	X	-	-	-
30	CHL	3	607	X	-	-	-
30	CHL	3	611	X	-	-	-
30	CHL	4	610	X	-	-	-
30	CHL	4	611	X	-	-	-
30	CHL	4	613	X	-	-	-
30	CHL	4	615	X	-	-	-
31	XAT	2	502	X	-	-	-
31	XAT	4	502	X	-	-	-

2 Entry composition

There are 36 unique types of molecules in this entry. The entry contains 39217 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 Photosystem I P700 chlorophyll a apoprotein A1.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	743	Total	C	N	O	S	0	0
			5858	3839	998	1003	18		

- Molecule 2 is a protein called Photosystem I P700 chlorophyll a apoprotein A2.

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	733	Total	C	N	O	S	0	0
			5857	3848	998	997	14		

- Molecule 3 is a protein called Photosystem I iron-sulfur center.

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	80	Total	C	N	O	S	0	0
			612	379	107	115	11		

- Molecule 4 is a protein called PsaD.

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	143	Total	C	N	O	S	0	0
			1132	731	194	204	3		

- Molecule 5 is a protein called PsaE.

Mol	Chain	Residues	Atoms				AltConf	Trace
5	E	66	Total	C	N	O	0	0
			528	336	93	99		

- Molecule 6 is a protein called PsaF.

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	154	Total	C	N	O	S	0	0
			1206	782	207	215	2		

- Molecule 7 is a protein called PsaG.

Mol	Chain	Residues	Atoms				AltConf	Trace
7	G	97	Total	C	N	O	0	0
			757	492	125	140		

- Molecule 8 is a protein called PsaH.

Mol	Chain	Residues	Atoms				AltConf	Trace
8	H	93	Total	C	N	O	0	0
			712	466	112	134		

- Molecule 9 is a protein called Photosystem I reaction center subunit VIII.

Mol	Chain	Residues	Atoms					AltConf	Trace
9	I	31	Total	C	N	O	S	0	0
			240	165	38	36	1		

- Molecule 10 is a protein called PsaJ.

Mol	Chain	Residues	Atoms					AltConf	Trace
10	J	42	Total	C	N	O	S	0	0
			338	231	51	55	1		

- Molecule 11 is a protein called Photosystem I reaction center subunit X psaK.

Mol	Chain	Residues	Atoms					AltConf	Trace
11	K	81	Total	C	N	O	S	0	0
			569	362	99	105	3		

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
K	86	ALA	VAL	conflict	UNP E1C9L3

- Molecule 12 is a protein called PsaL.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	L	159	Total	C	N	O	S	0	0
			1197	788	191	217	1		

- Molecule 13 is a protein called Lhca1.

Mol	Chain	Residues	Atoms					AltConf	Trace
13	1	193	Total	C	N	O	S	0	0
			1508	982	252	269	5		

- Molecule 14 is a protein called Chlorophyll a-b binding protein, chloroplastic.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	2	208	Total	C	N	O	S	0	0
			1620	1059	265	292	4		

- Molecule 15 is a protein called Chlorophyll a-b binding protein 3, chloroplastic.

Mol	Chain	Residues	Atoms					AltConf	Trace
15	3	221	Total	C	N	O	S	0	0
			1706	1118	278	305	5		

- Molecule 16 is a protein called Chlorophyll a-b binding protein P4, chloroplastic.

Mol	Chain	Residues	Atoms					AltConf	Trace
16	4	198	Total	C	N	O	S	0	0
			1559	1022	253	281	3		

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
4	89	LYS	ARG	conflict	UNP Q9SQL2
4	128	ASP	ALA	conflict	UNP Q9SQL2
4	149	PHE	SER	conflict	UNP Q9SQL2

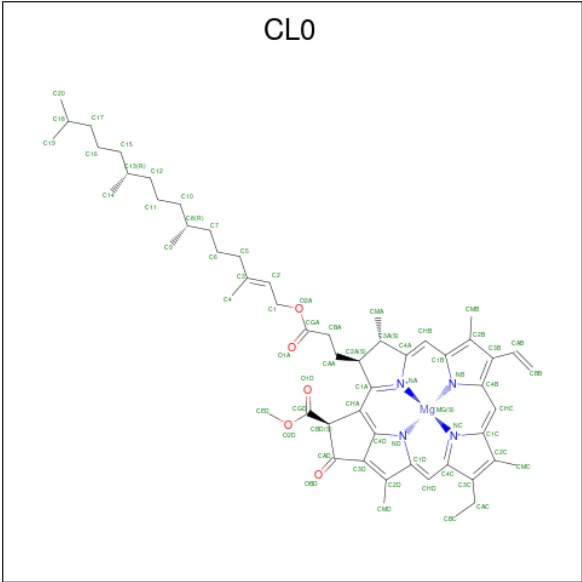
- Molecule 17 is a protein called Ferredoxin-1, chloroplastic.

Mol	Chain	Residues	Atoms					AltConf	Trace
17	N	97	Total	C	N	O	S	0	0
			724	448	111	160	5		

- Molecule 18 is a protein called Plastocyanin, chloroplastic.

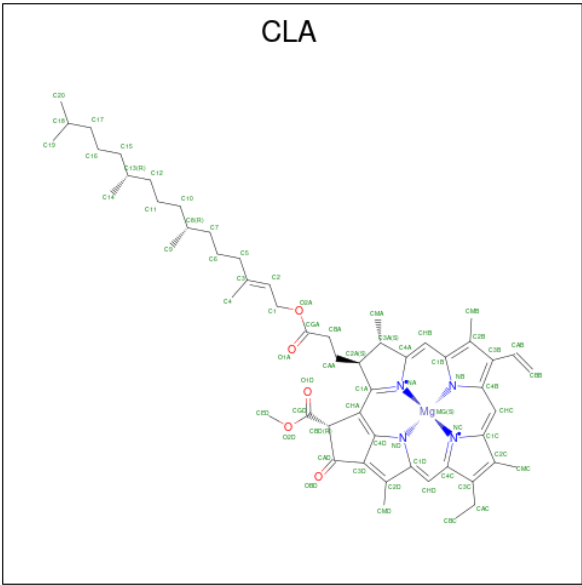
Mol	Chain	Residues	Atoms					AltConf	Trace
18	P	99	Total	C	N	O	S	0	0
			728	460	115	150	3		

- Molecule 19 is CHLOROPHYLL A ISOMER (CCD ID: CL0) (formula: C₅₅H₇₂MgN₄O₅).



Mol	Chain	Residues	Atoms					AltConf
19	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	

- Molecule 20 is CHLOROPHYLL A (CCD ID: CLA) (formula: C₅₅H₇₂MgN₄O₅).



Mol	Chain	Residues	Atoms					AltConf
20	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
20	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
20	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	

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Mol	Chain	Residues	Atoms					AltConf
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 45	C 35	Mg 1	N 4	O 5	0
20	A	1	Total 46	C 36	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 56	C 46	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	A	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 55	C 45	Mg 1	N 4	O 5	0

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Mol	Chain	Residues	Atoms					AltConf
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	A	1	Total 51	C 41	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0

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Mol	Chain	Residues	Atoms					AltConf
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	B	1	Total 46	C 36	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	B	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 46	C 36	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0

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Mol	Chain	Residues	Atoms					AltConf
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 58	C 48	Mg 1	N 4	O 5	0
20	B	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	B	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	B	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	F	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	F	1	Total 65	C 55	Mg 1	N 4	O 5	0

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Mol	Chain	Residues	Atoms					AltConf
20	G	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	G	1	Total 46	C 36	Mg 1	N 4	O 5	0
20	G	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	H	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	J	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	K	1	Total 45	C 35	Mg 1	N 4	O 5	0
20	K	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	K	1	Total 48	C 38	Mg 1	N 4	O 5	0
20	K	1	Total 46	C 36	Mg 1	N 4	O 5	0
20	L	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	L	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	L	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	1	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	1	1	Total 46	C 36	Mg 1	N 4	O 5	0
20	1	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	1	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	1	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	1	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	1	1	Total 46	C 36	Mg 1	N 4	O 5	0
20	1	1	Total 46	C 36	Mg 1	N 4	O 5	0
20	1	1	Total 65	C 55	Mg 1	N 4	O 5	0

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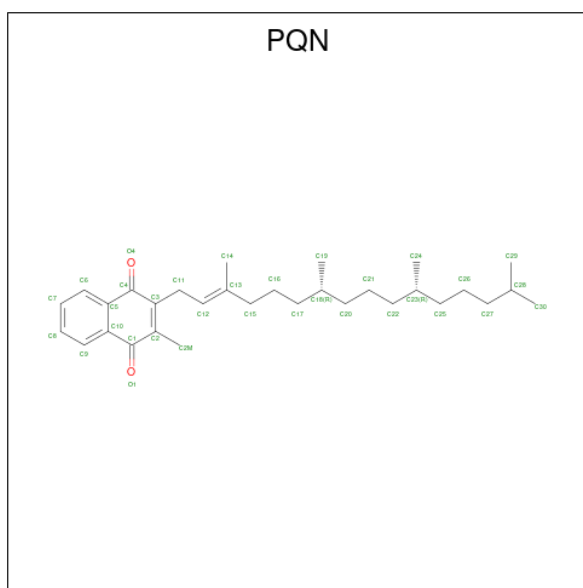
Mol	Chain	Residues	Atoms					AltConf
20	1	1	Total 45	C 35	Mg 1	N 4	O 5	0
20	1	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	2	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	2	1	Total 52	C 42	Mg 1	N 4	O 5	0
20	2	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	2	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	2	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	2	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	2	1	Total 60	C 50	Mg 1	N 4	O 5	0
20	2	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	2	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	3	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	3	1	Total 52	C 42	Mg 1	N 4	O 5	0
20	3	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	3	1	Total 55	C 45	Mg 1	N 4	O 5	0
20	3	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	3	1	Total 48	C 38	Mg 1	N 4	O 5	0
20	3	1	Total 65	C 55	Mg 1	N 4	O 5	0
20	3	1	Total 50	C 40	Mg 1	N 4	O 5	0
20	3	1	Total 46	C 36	Mg 1	N 4	O 5	0
20	3	1	Total 42	C 34	Mg 1	N 4	O 3	0

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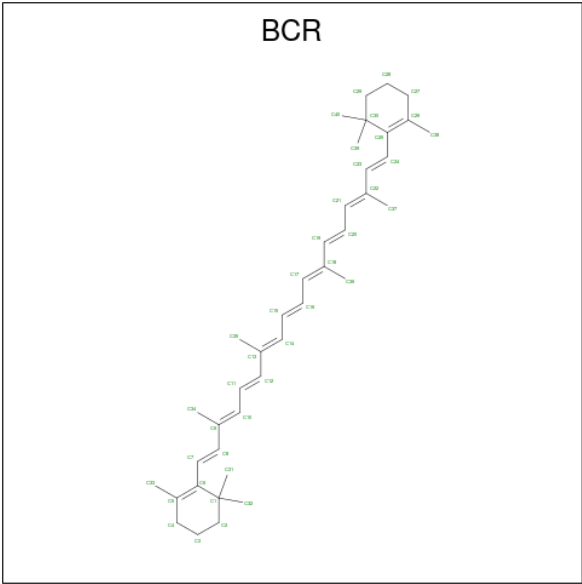
Mol	Chain	Residues	Atoms					AltConf
20	3	1	Total	C	Mg	N	O	0
			60	50	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			50	40	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			60	50	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			50	40	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			60	50	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			60	50	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			50	40	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			60	50	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			46	36	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
20	4	1	Total	C	Mg	N	O	0
			65	55	1	4	5	

- Molecule 21 is PHYLLOQUINONE (CCD ID: PQN) (formula: $C_{31}H_{46}O_2$).



Mol	Chain	Residues	Atoms			AltConf
21	A	1	Total	C	O	0
			33	31	2	
21	B	1	Total	C	O	0
			33	31	2	

- Molecule 22 is BETA-CAROTENE (CCD ID: BCR) (formula: C₄₀H₅₆).



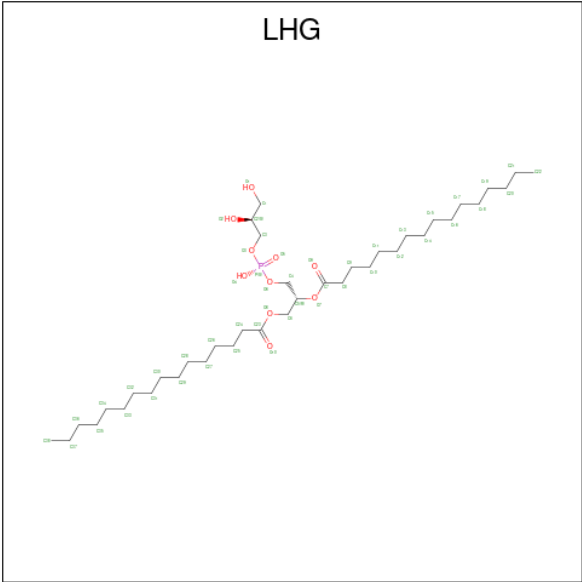
Mol	Chain	Residues	Atoms		AltConf
22	A	1	Total	C	0
			40	40	
22	A	1	Total	C	0
			40	40	
22	A	1	Total	C	0
			40	40	
22	A	1	Total	C	0
			40	40	
22	A	1	Total	C	0
			40	40	
22	A	1	Total	C	0
			40	40	
22	B	1	Total	C	0
			40	40	
22	B	1	Total	C	0
			40	40	
22	B	1	Total	C	0
			40	40	
22	B	1	Total	C	0
			40	40	

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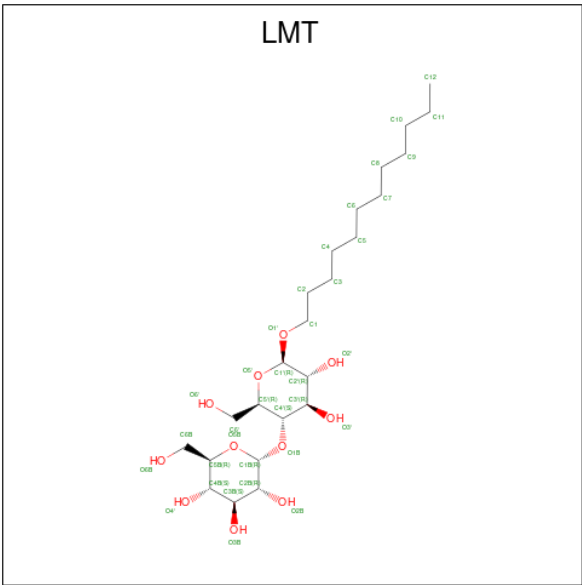
Mol	Chain	Residues	Atoms	AltConf
22	B	1	Total C 40 40	0
22	F	1	Total C 40 40	0
22	F	1	Total C 40 40	0
22	G	1	Total C 40 40	0
22	H	1	Total C 40 40	0
22	I	1	Total C 40 40	0
22	I	1	Total C 40 40	0
22	J	1	Total C 40 40	0
22	K	1	Total C 40 40	0
22	K	1	Total C 40 40	0
22	L	1	Total C 40 40	0
22	L	1	Total C 40 40	0
22	1	1	Total C 40 40	0
22	1	1	Total C 40 40	0
22	2	1	Total C 40 40	0
22	3	1	Total C 40 40	0
22	3	1	Total C 40 40	0

- Molecule 23 is 1,2-DIPALMITOYL-PHOSPHATIDYL-GLYCEROLE (CCD ID: LHG) (formula: $C_{38}H_{75}O_{10}P$).



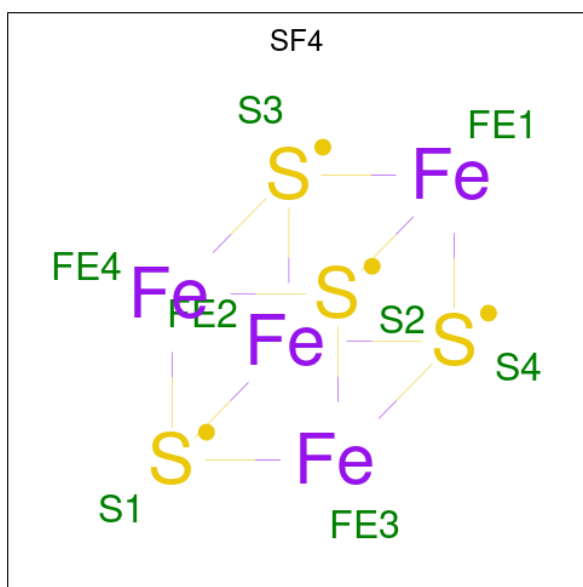
Mol	Chain	Residues	Atoms				AltConf
23	A	1	Total	C	O	P	0
			40	29	10	1	
23	A	1	Total	C	O	P	0
			49	38	10	1	
23	B	1	Total	C	O	P	0
			21	10	10	1	
23	B	1	Total	C	O	P	0
			49	38	10	1	
23	1	1	Total	C	O	P	0
			49	38	10	1	
23	2	1	Total	C	O	P	0
			35	24	10	1	
23	3	1	Total	C	O	P	0
			17	8	8	1	
23	4	1	Total	C	O	P	0
			35	24	10	1	

- Molecule 24 is DODECYL-BETA-D-MALTOSIDE (CCD ID: LMT) (formula: C₂₄H₄₆O₁₁).



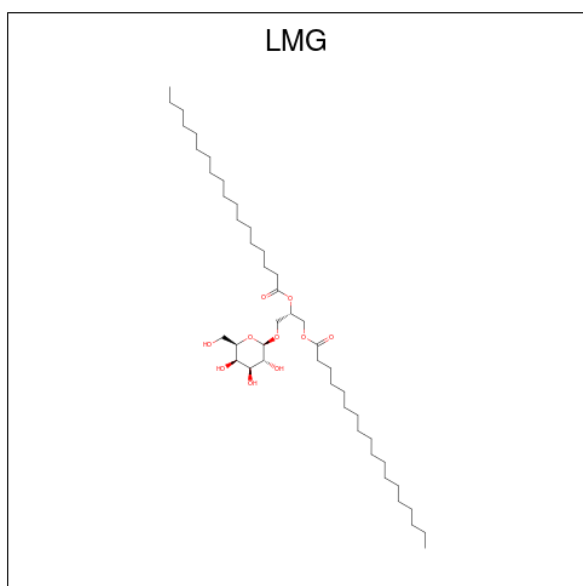
Mol	Chain	Residues	Atoms			AltConf
24	A	1	Total	C	O	0
			35	24	11	
24	B	1	Total	C	O	0
			32	21	11	
24	B	1	Total	C	O	0
			31	20	11	
24	G	1	Total	C	O	0
			31	20	11	
24	G	1	Total	C	O	0
			35	24	11	
24	2	1	Total	C	O	0
			35	24	11	

- Molecule 25 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄).



Mol	Chain	Residues	Atoms			AltConf
25	A	1	Total	Fe	S	0
			8	4	4	
25	C	1	Total	Fe	S	0
			8	4	4	
25	C	1	Total	Fe	S	0
			8	4	4	

- Molecule 26 is 1,2-DISTEAROYL-MONOGALACTOSYL-DIGLYCERIDE (CCD ID: LMG) (formula: $C_{45}H_{86}O_{10}$).



Mol	Chain	Residues	Atoms			AltConf
26	A	1	Total	C	O	0
			50	40	10	
26	B	1	Total	C	O	0
			35	25	10	
26	B	1	Total	C	O	0
			33	23	10	
26	B	1	Total	C	O	0
			34	24	10	
26	F	1	Total	C	O	0
			47	37	10	
26	F	1	Total	C	O	0
			36	26	10	
26	F	1	Total	C	O	0
			34	24	10	
26	F	1	Total	C	O	0
			13	7	6	
26	F	1	Total	C	O	0
			30	20	10	
26	G	1	Total	C	O	0
			25	15	10	
26	G	1	Total	C	O	0
			49	39	10	
26	G	1	Total	C	O	0
			50	40	10	
26	1	1	Total	C	O	0
			46	36	10	
26	2	1	Total	C	O	0
			25	15	10	
26	2	1	Total	C	O	0
			36	26	10	
26	2	1	Total	C	O	0
			30	20	10	
26	2	1	Total	C	O	0
			30	20	10	
26	2	1	Total	C	O	0
			13	7	6	
26	3	1	Total	C	O	0
			30	20	10	

- Molecule 27 is CALCIUM ION (CCD ID: CA) (formula: Ca).

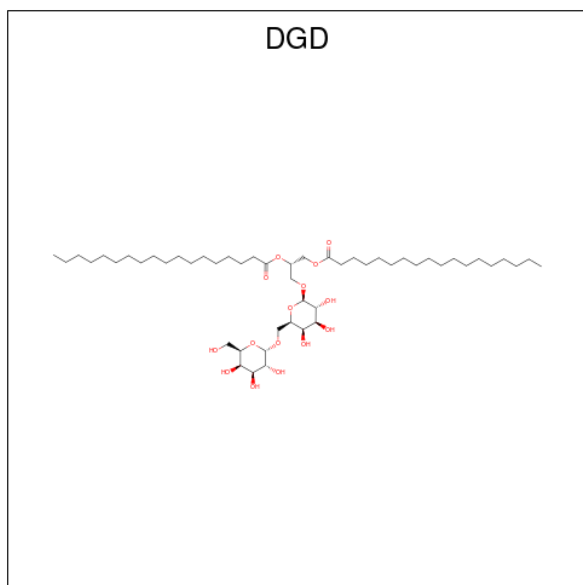
Mol	Chain	Residues	Atoms		AltConf
27	A	1	Total	Ca	0
			1	1	

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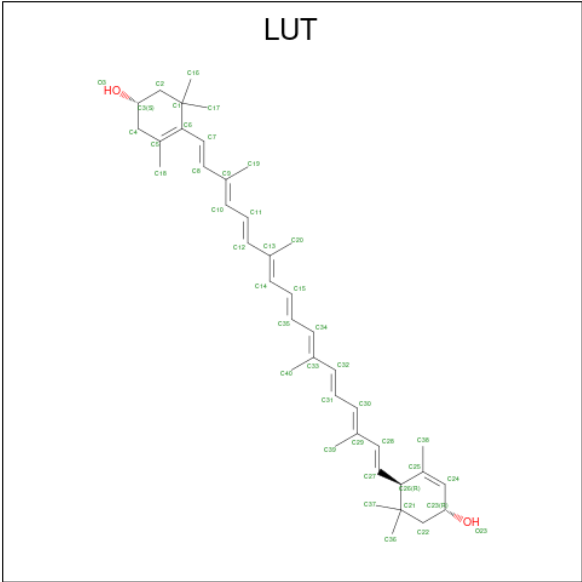
Mol	Chain	Residues	Atoms		AltConf
27	B	1	Total	Ca	0
			1	1	

- Molecule 28 is DIGALACTOSYL DIACYL GLYCEROL (DGDG) (CCD ID: DGD) (formula: $C_{51}H_{96}O_{15}$).



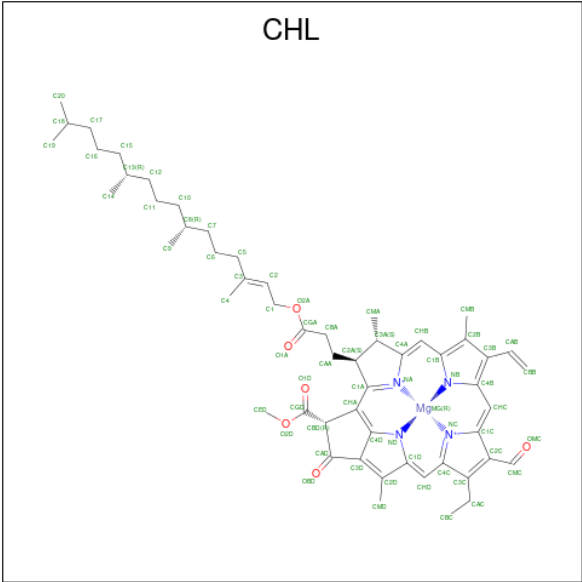
Mol	Chain	Residues	Atoms			AltConf
28	B	1	Total	C	O	0
			61	46	15	
28	F	1	Total	C	O	0
			57	42	15	
28	G	1	Total	C	O	0
			47	32	15	
28	J	1	Total	C	O	0
			58	43	15	
28	1	1	Total	C	O	0
			41	26	15	
28	3	1	Total	C	O	0
			51	36	15	
28	4	1	Total	C	O	0
			51	36	15	

- Molecule 29 is (3R,3'R,6S)-4,5-DIDEHYDRO-5,6-DIHYDRO-BETA,BETA-CAROTENE-3,3'-DIOL (CCD ID: LUT) (formula: $C_{40}H_{56}O_2$).



Mol	Chain	Residues	Atoms			AltConf
29	J	1	Total	C	O	0
			42	40	2	
29	1	1	Total	C	O	0
			42	40	2	
29	1	1	Total	C	O	0
			42	40	2	
29	2	1	Total	C	O	0
			42	40	2	
29	3	1	Total	C	O	0
			42	40	2	
29	3	1	Total	C	O	0
			42	40	2	
29	4	1	Total	C	O	0
			42	40	2	

- Molecule 30 is CHLOROPHYLL B (CCD ID: CHL) (formula: C₅₅H₇₀MgN₄O₆).



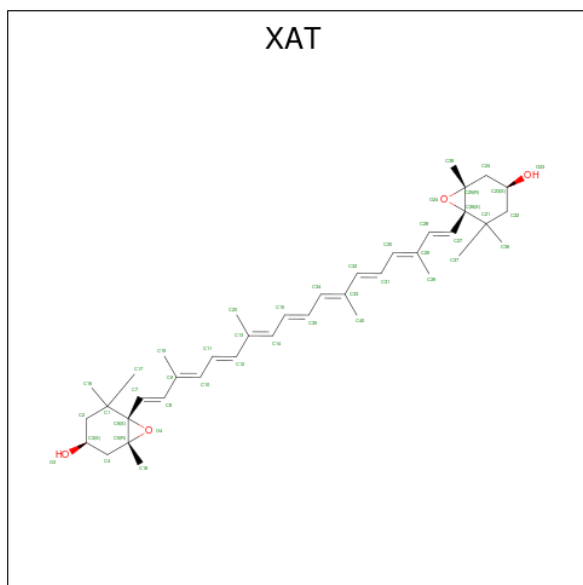
Mol	Chain	Residues	Atoms					AltConf
30	1	1	Total 47	C 36	Mg 1	N 4	O 6	0
30	1	1	Total 61	C 50	Mg 1	N 4	O 6	0
30	1	1	Total 56	C 45	Mg 1	N 4	O 6	0
30	2	1	Total 56	C 45	Mg 1	N 4	O 6	0
30	2	1	Total 48	C 37	Mg 1	N 4	O 6	0
30	2	1	Total 46	C 35	Mg 1	N 4	O 6	0
30	2	1	Total 56	C 45	Mg 1	N 4	O 6	0
30	2	1	Total 66	C 55	Mg 1	N 4	O 6	0
30	3	1	Total 66	C 55	Mg 1	N 4	O 6	0
30	3	1	Total 51	C 40	Mg 1	N 4	O 6	0
30	3	1	Total 47	C 36	Mg 1	N 4	O 6	0
30	4	1	Total 47	C 36	Mg 1	N 4	O 6	0
30	4	1	Total 51	C 40	Mg 1	N 4	O 6	0
30	4	1	Total 61	C 50	Mg 1	N 4	O 6	0

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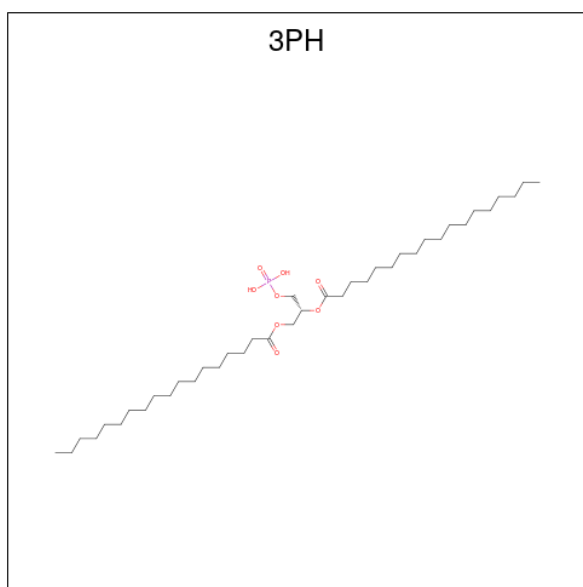
Mol	Chain	Residues	Atoms					AltConf
30	4	1	Total	C	Mg	N	O	0
			43	34	1	4	4	

- Molecule 31 is (3S,5R,6S,3'S,5'R,6'S)-5,6,5',6'-DIEPOXY-5,6,5',6'-TETRAHYDRO-BETA, BETA-CAROTENE-3,3'-DIOL (CCD ID: XAT) (formula: C₄₀H₅₆O₄).



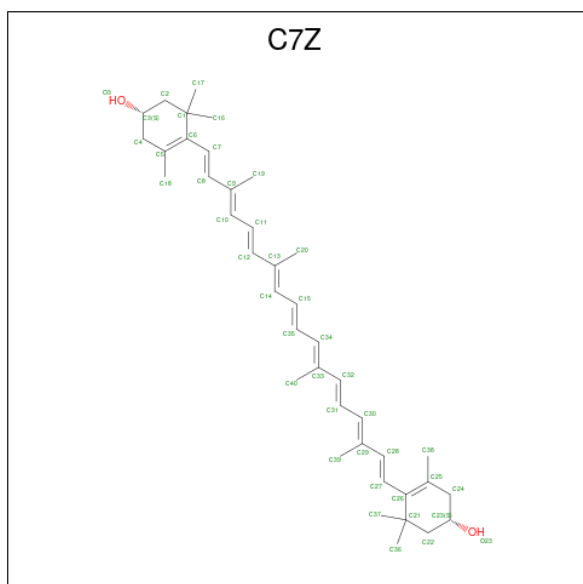
Mol	Chain	Residues	Atoms			AltConf
31	2	1	Total	C	O	0
			44	40	4	
31	4	1	Total	C	O	0
			44	40	4	

- Molecule 32 is 1,2-DIACYL-GLYCEROL-3-SN-PHOSPHATE (CCD ID: 3PH) (formula: C₃₉H₇₇O₈P).



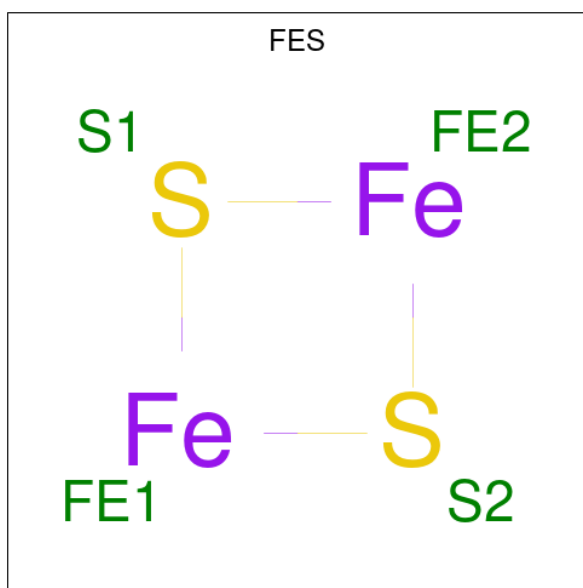
Mol	Chain	Residues	Atoms				AltConf
32	2	1	Total	C	O	P	0
			33	24	8	1	

- Molecule 33 is (1 {S})-3,5,5-trimethyl-4-[(1 {E},3 {E},5 {E},7 {E},9 {E},11 {E},13 {E},15 {E},17 {E})-3,7,12,16-tetramethyl-18-[(4 {S})-2,6,6-trimethyl-4-oxidanyl-cyclohexen-1-yl]octadeca-1,3,5,7,9,11,13,15,17-nonaenyl]cyclohex-3-en-1-ol (CCD ID: C7Z) (formula: $C_{40}H_{56}O_2$).



Mol	Chain	Residues	Atoms				AltConf
33	4	1	Total	C	O		0
			42	40	2		

- Molecule 34 is FE2/S2 (INORGANIC) CLUSTER (CCD ID: FES) (formula: Fe_2S_2).



Mol	Chain	Residues	Atoms			AltConf
34	N	1	Total	Fe	S	0
			4	2	2	

- Molecule 35 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		AltConf
35	P	1	Total	Cu	0
			1	1	

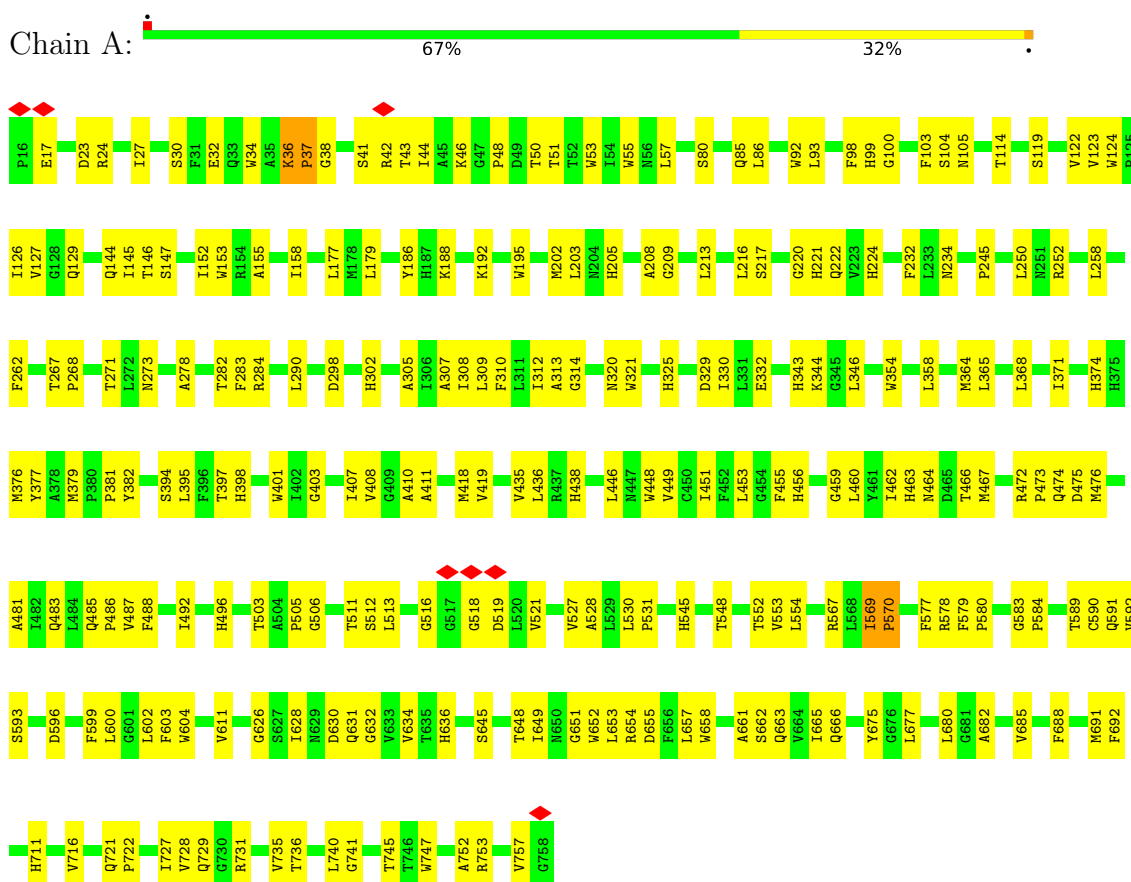
- Molecule 36 is water.

Mol	Chain	Residues	Atoms		AltConf
36	B	2	Total	O	0
			2	2	

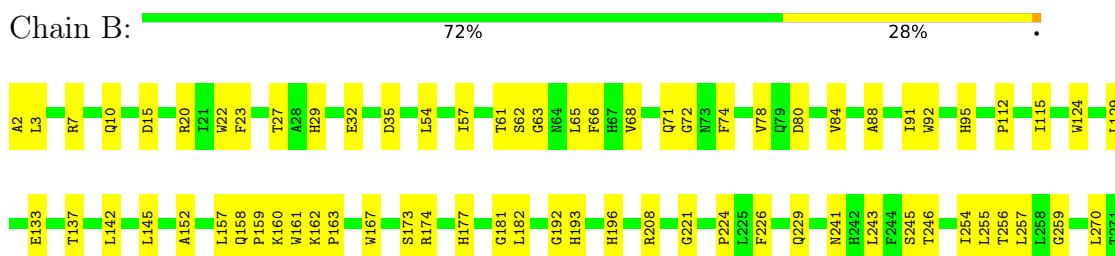
3 Residue-property plots

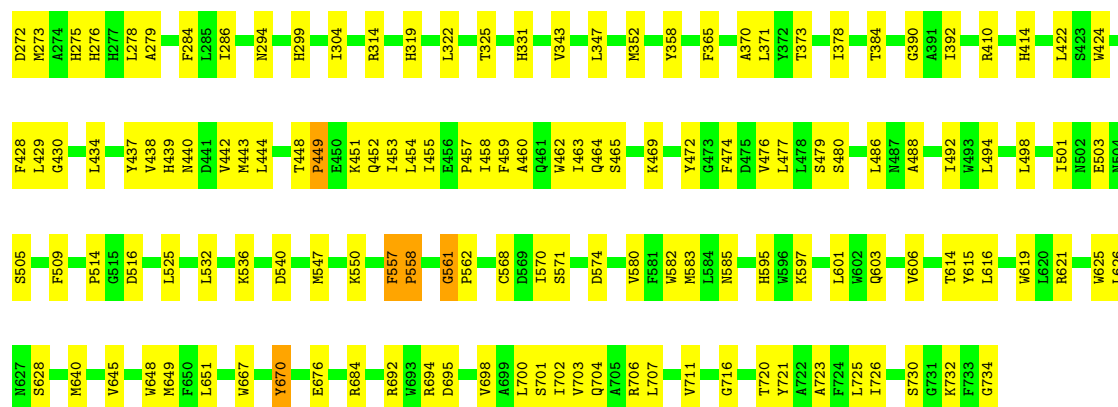
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion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

- Molecule 1: Photosystem I P700 chlorophyll a apoprotein A1



- Molecule 2: Photosystem I P700 chlorophyll a apoprotein A2





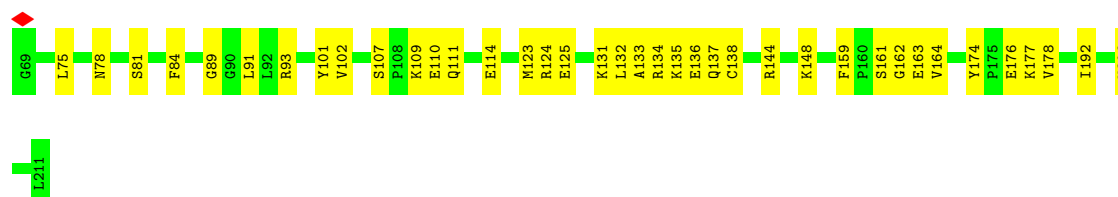
• Molecule 3: Photosystem I iron-sulfur center

Chain C: 66% 32% .



• Molecule 4: PsaD

Chain D: 73% 27%



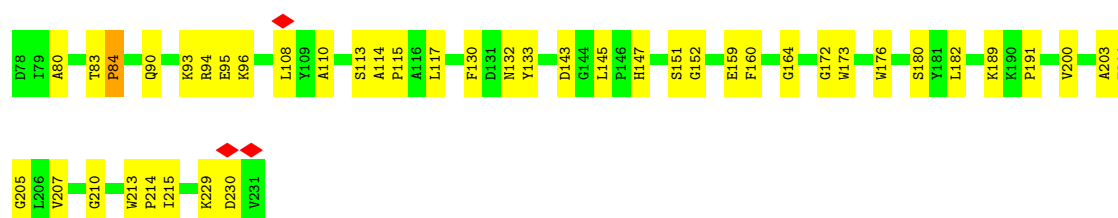
• Molecule 5: PsaE

Chain E: 5% 88% 12%

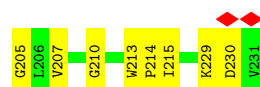


• Molecule 6: PsaF

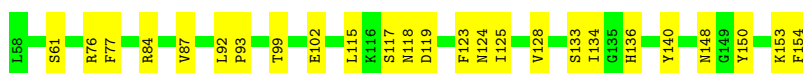
Chain F: 72% 27%



• Molecule 7: PsaG

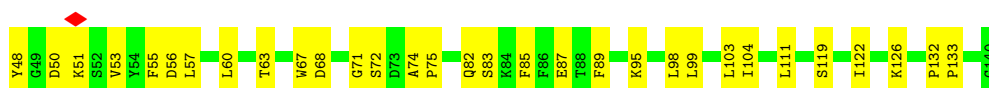


Chain G:  74% 26%



• Molecule 8: PsaH

Chain H:  67% 33%



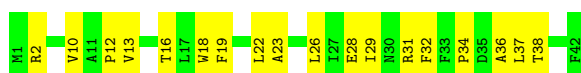
• Molecule 9: Photosystem I reaction center subunit VIII

Chain I:  61% 39%



• Molecule 10: PsaJ

Chain J:  57% 43%



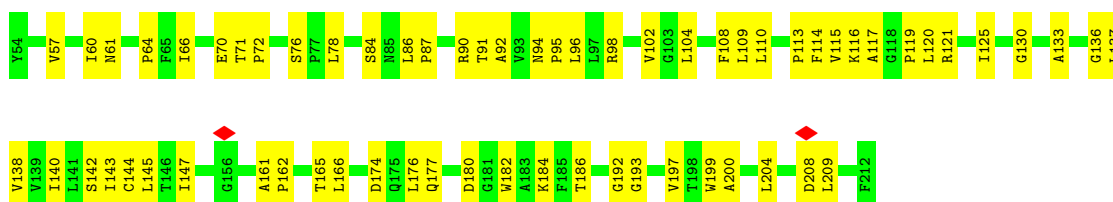
• Molecule 11: Photosystem I reaction center subunit X psaK

Chain K:  32% 63% 37%



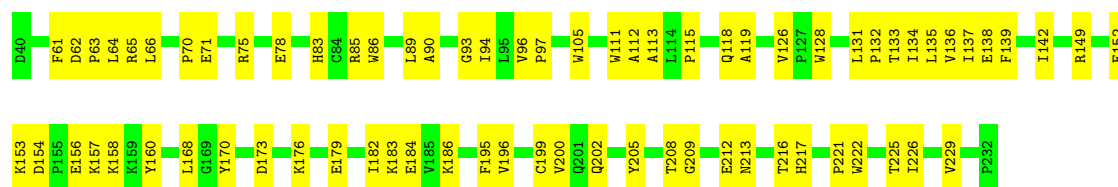
• Molecule 12: PsaL

Chain L:  60% 40%



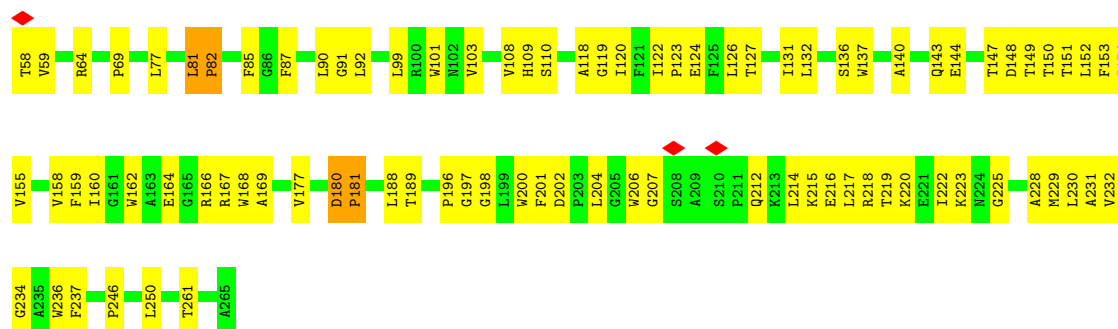
• Molecule 13: Lhca1

Chain 1:  63% 37%



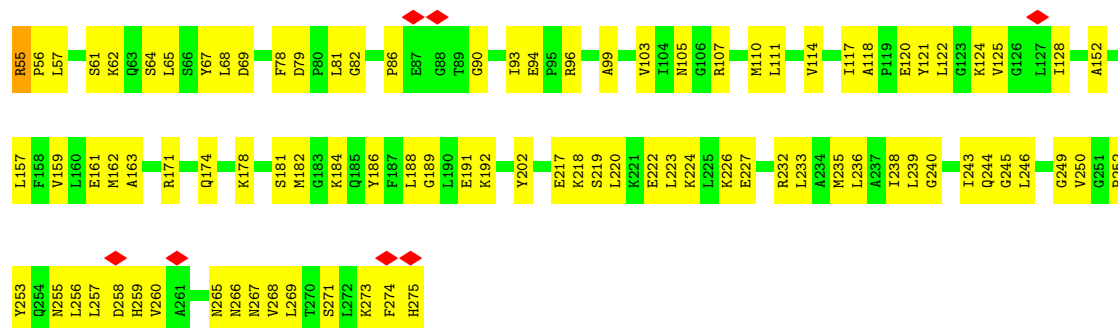
• Molecule 14: Chlorophyll a-b binding protein, chloroplastic

Chain 2:  58% 40%



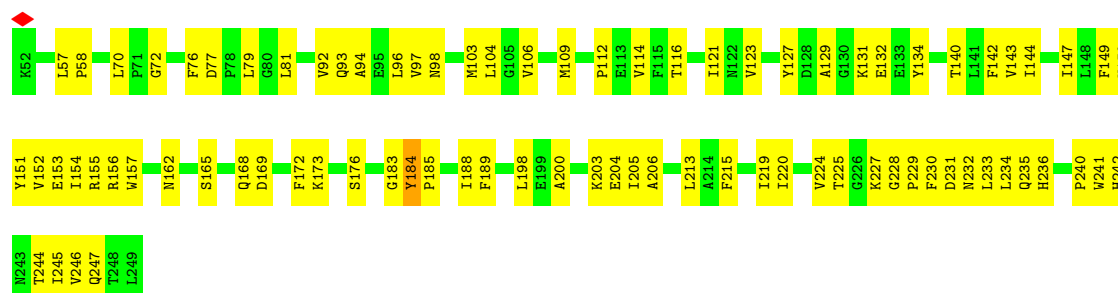
• Molecule 15: Chlorophyll a-b binding protein 3, chloroplastic

Chain 3:  58% 41%



• Molecule 16: Chlorophyll a-b binding protein P4, chloroplastic

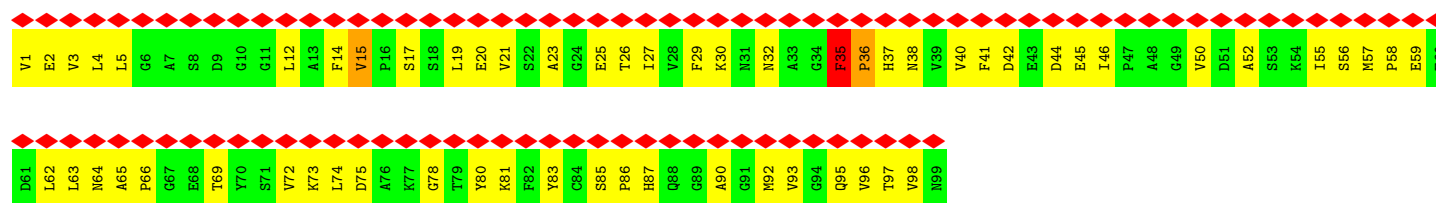
Chain 4:  58% 41%



- Molecule 17: Ferredoxin-1, chloroplastic



- Molecule 18: Plastocyanin, chloroplastic



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	102216	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 ($e^-/\text{\AA}^2$)	47.05	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	2000	Depositor
Magnification	105000	Depositor
Image detector	GATAN K3 BIOQUANTUM (6k x 4k)	Depositor
Maximum map value	0.184	Depositor
Minimum map value	-0.047	Depositor
Average map value	0.001	Depositor
Map value standard deviation	0.006	Depositor
Recommended contour level	0.022	Depositor
Map size ($\text{\AA}$)	246.6, 246.6, 246.6	wwPDB
Map dimensions	300, 300, 300	wwPDB
Map angles ($^\circ$)	90.0, 90.0, 90.0	wwPDB
Pixel spacing ($\text{\AA}$)	0.822, 0.822, 0.822	Depositor

5 Model quality i

5.1 Standard geometry i

Bond lengths and bond angles in the following residue types are not validated in this section: SF4, CLA, FES, LMT, BCR, LMG, C7Z, XAT, DGD, CL0, 3PH, LUT, CHL, LHG, CU, PQN, CA

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	0.61	4/6057 (0.1%)	0.71	7/8264 (0.1%)
2	B	0.60	5/6069 (0.1%)	0.63	10/8286 (0.1%)
3	C	1.00	2/625 (0.3%)	1.21	4/846 (0.5%)
4	D	0.39	0/1163	0.55	0/1572
5	E	0.30	0/540	0.43	0/734
6	F	0.83	2/1234 (0.2%)	0.99	4/1670 (0.2%)
7	G	0.30	0/776	0.42	0/1054
8	H	0.61	0/733	0.78	0/995
9	I	0.32	0/246	0.57	0/335
10	J	0.42	0/349	0.52	0/476
11	K	0.65	0/576	0.87	0/779
12	L	0.80	2/1232 (0.2%)	1.18	7/1684 (0.4%)
13	1	0.36	0/1558	0.52	0/2125
14	2	0.86	4/1679 (0.2%)	1.11	6/2302 (0.3%)
15	3	0.60	1/1760 (0.1%)	0.85	5/2390 (0.2%)
16	4	0.55	1/1608 (0.1%)	0.76	4/2191 (0.2%)
17	N	1.19	2/736 (0.3%)	1.71	14/1000 (1.4%)
18	P	1.19	3/743 (0.4%)	1.20	3/1009 (0.3%)
All	All	0.67	26/27684 (0.1%)	0.83	64/37712 (0.2%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
18	P	0	2

The worst 5 of 26 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	561	GLY	C-N	20.63	1.50	1.33
1	A	37	PRO	N-CA	18.42	1.70	1.47
14	2	82	PRO	N-CA	18.38	1.70	1.47
1	A	570	PRO	N-CA	18.14	1.70	1.47
6	F	84	PRO	N-CA	18.13	1.70	1.47

The worst 5 of 64 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
14	2	180	ASP	CA-C-N	19.71	139.66	119.56
14	2	180	ASP	C-N-CA	19.71	139.66	119.56
2	B	448	THR	CA-C-N	16.77	138.44	119.32
2	B	448	THR	C-N-CA	16.77	138.44	119.32
3	C	21	CYS	CA-C-N	16.74	137.28	119.87

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
18	P	15	VAL	Peptide
18	P	35	PHE	Peptide

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	5858	0	5719	336	0
2	B	5857	0	5653	231	0
3	C	612	0	591	31	0
4	D	1132	0	1141	47	0
5	E	528	0	528	5	0
6	F	1206	0	1231	48	0
7	G	757	0	743	24	0
8	H	712	0	701	38	0
9	I	240	0	264	30	0
10	J	338	0	345	35	0
11	K	569	0	596	38	0
12	L	1197	0	1197	87	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
13	1	1508	0	1489	117	0
14	2	1620	0	1557	135	0
15	3	1706	0	1661	125	0
16	4	1559	0	1527	144	0
17	N	724	0	672	69	0
18	P	728	0	699	92	0
19	A	65	0	72	11	0
20	1	608	0	563	96	0
20	2	522	0	499	95	0
20	3	578	0	495	60	0
20	4	631	0	595	169	0
20	A	2643	0	2749	345	0
20	B	2610	0	2747	241	0
20	F	130	0	144	21	0
20	G	166	0	153	21	0
20	H	60	0	58	18	0
20	J	50	0	38	7	0
20	K	199	0	158	16	0
20	L	160	0	134	49	0
21	A	33	0	46	8	0
21	B	33	0	46	3	0
22	1	80	0	105	5	0
22	2	40	0	53	23	0
22	3	80	0	105	10	0
22	A	240	0	316	25	0
22	B	200	0	265	19	0
22	F	80	0	105	12	0
22	G	40	0	53	4	0
22	H	40	0	53	18	0
22	I	80	0	105	17	0
22	J	40	0	53	5	0
22	K	80	0	106	15	0
22	L	80	0	106	46	0
23	1	49	0	74	11	0
23	2	35	0	40	7	0
23	3	17	0	12	0	0
23	4	35	0	40	11	0
23	A	89	0	127	3	0
23	B	70	0	86	6	0
24	2	35	0	46	7	0
24	A	35	0	45	1	0
24	B	63	0	68	1	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
24	G	66	0	79	0	0
25	A	8	0	0	0	0
25	C	16	0	0	3	0
26	1	46	0	65	8	0
26	2	134	0	133	17	0
26	3	30	0	30	0	0
26	A	50	0	73	3	0
26	B	102	0	114	16	0
26	F	160	0	188	8	0
26	G	124	0	161	15	0
27	A	1	0	0	0	0
27	B	1	0	0	0	0
28	1	41	0	40	1	0
28	3	51	0	60	3	0
28	4	51	0	60	4	0
28	B	61	0	83	7	0
28	F	57	0	75	18	0
28	G	47	0	52	3	0
28	J	58	0	77	11	0
29	1	84	0	110	39	0
29	2	42	0	55	12	0
29	3	84	0	110	7	0
29	4	42	0	55	19	0
29	J	42	0	55	9	0
30	1	164	0	134	28	0
30	2	272	0	223	52	0
30	3	164	0	136	8	0
30	4	202	0	152	11	0
31	2	44	0	56	13	0
31	4	44	0	56	14	0
32	2	33	0	39	5	0
33	4	42	0	0	4	0
34	N	4	0	0	0	0
35	P	1	0	0	0	0
36	B	2	0	0	0	0
All	All	39217	0	39245	2323	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 30.

The worst 5 of 2323 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:558:PRO:N	2:B:558:PRO:CA	1.69	1.44
14:2:181:PRO:CA	14:2:181:PRO:N	1.70	1.43
6:F:84:PRO:N	6:F:84:PRO:CA	1.70	1.41
1:A:570:PRO:N	1:A:570:PRO:CA	1.70	1.41
3:C:22:PRO:CA	3:C:22:PRO:N	1.69	1.38

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

There are no protein backbone outliers to report in this entry.

5.3.2 Protein sidechains [i](#)

There are no protein residues with a non-rotameric sidechain to report in this entry.

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

Of 244 ligands modelled in this entry, 3 are monoatomic - leaving 241 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the

expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
20	CLA	A	1111	-	69,73,73	1.41	6 (8%)	82,113,113	1.96	19 (23%)
20	CLA	A	1128	-	69,73,73	1.41	6 (8%)	82,113,113	1.93	19 (23%)
20	CLA	A	1132	-	69,73,73	1.35	6 (8%)	82,113,113	1.93	18 (21%)
22	BCR	G	4011	-	41,41,41	1.63	4 (9%)	56,56,56	4.71	17 (30%)
20	CLA	3	614	-	46,50,73	1.67	7 (15%)	53,85,113	2.16	13 (24%)
31	XAT	4	502	-	41,47,47	0.76	1 (2%)	54,74,74	2.09	14 (25%)
20	CLA	A	1103	-	69,73,73	1.39	6 (8%)	82,113,113	1.99	20 (24%)
20	CLA	A	1106	1	69,73,73	1.40	6 (8%)	82,113,113	1.91	20 (24%)
26	LMG	B	5003	-	35,35,55	0.79	1 (2%)	43,43,63	1.15	4 (9%)
30	CHL	1	610	13	41,55,74	1.80	9 (21%)	35,91,114	2.10	9 (25%)
22	BCR	F	4016	-	41,41,41	1.68	5 (12%)	56,56,56	4.46	16 (28%)
26	LMG	F	5006	-	13,13,55	0.55	0	18,18,63	0.77	0
20	CLA	4	617	-	69,73,73	1.34	6 (8%)	82,113,113	1.98	18 (21%)
28	DGD	3	803	-	52,52,67	0.93	3 (5%)	66,66,81	1.15	4 (6%)
20	CLA	B	1220	-	59,63,73	1.51	6 (10%)	70,101,113	1.95	16 (22%)
20	CLA	1	602	13	50,54,73	1.61	6 (12%)	59,90,113	2.02	16 (27%)
30	CHL	4	610	-	41,55,74	1.90	9 (21%)	35,91,114	1.98	9 (25%)
20	CLA	K	1403	-	52,56,73	1.57	6 (11%)	61,92,113	2.13	17 (27%)
20	CLA	4	604	-	64,68,73	1.40	6 (9%)	76,107,113	1.98	19 (25%)
33	C7Z	4	505	-	43,43,43	3.91	17 (39%)	56,60,60	5.58	32 (57%)
22	BCR	3	506	-	41,41,41	1.69	4 (9%)	56,56,56	4.57	17 (30%)
20	CLA	1	604	-	69,73,73	1.38	6 (8%)	82,113,113	1.89	18 (21%)
26	LMG	B	5007	-	34,34,55	0.51	0	42,42,63	1.08	3 (7%)
29	LUT	4	501	-	42,43,43	2.48	4 (9%)	51,60,60	2.13	15 (29%)
20	CLA	A	1136	-	69,73,73	1.38	6 (8%)	82,113,113	1.84	17 (20%)
22	BCR	A	4002	-	41,41,41	1.66	5 (12%)	56,56,56	4.41	17 (30%)
20	CLA	A	1112	-	69,73,73	1.40	6 (8%)	82,113,113	1.87	17 (20%)
30	CHL	2	609	-	60,74,74	1.27	9 (15%)	58,114,114	1.63	11 (18%)
22	BCR	1	503	-	41,41,41	1.65	4 (9%)	56,56,56	4.36	14 (25%)
20	CLA	A	1129	-	69,73,73	1.39	6 (8%)	82,113,113	1.89	17 (20%)
26	LMG	F	5003	-	36,36,55	0.78	1 (2%)	44,44,63	1.10	4 (9%)
26	LMG	2	803	-	36,36,55	0.69	1 (2%)	44,44,63	1.07	3 (6%)
20	CLA	B	1207	-	69,73,73	1.35	6 (8%)	82,113,113	1.93	19 (23%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
20	CLA	3	612	-	54,58,73	1.52	6 (11%)	64,95,113	2.12	17 (26%)
30	CHL	4	613	-	55,69,74	1.37	9 (16%)	52,108,114	1.79	12 (23%)
20	CLA	1	606	-	54,58,73	1.57	6 (11%)	64,95,113	2.09	19 (29%)
30	CHL	4	615	-	37,51,74	1.45	8 (21%)	30,86,114	2.37	11 (36%)
20	CLA	A	1113	-	49,53,73	1.61	6 (12%)	58,89,113	2.04	14 (24%)
20	CLA	B	1229	-	69,73,73	1.39	6 (8%)	82,113,113	1.97	21 (25%)
26	LMG	2	802	-	25,25,55	0.54	0	33,33,63	1.16	3 (9%)
20	CLA	B	1228	-	64,68,73	1.43	6 (9%)	76,107,113	1.91	13 (17%)
20	CLA	B	1022	-	69,73,73	1.45	6 (8%)	82,113,113	1.88	20 (24%)
20	CLA	1	608	-	50,54,73	1.61	6 (12%)	59,90,113	2.14	16 (27%)
20	CLA	3	601	-	59,63,73	1.49	6 (10%)	70,101,113	2.09	20 (28%)
26	LMG	2	805	-	30,30,55	0.53	0	38,38,63	1.10	2 (5%)
20	CLA	4	601	16	64,68,73	1.40	6 (9%)	76,107,113	1.94	18 (23%)
20	CLA	1	601	13	69,73,73	1.39	6 (8%)	82,113,113	1.95	21 (25%)
20	CLA	2	607	23	64,68,73	1.41	6 (9%)	76,107,113	2.04	20 (26%)
19	CL0	A	1011	-	58,73,73	2.94	16 (27%)	60,113,113	2.89	21 (35%)
20	CLA	B	1218	-	69,73,73	1.37	6 (8%)	82,113,113	1.95	17 (20%)
20	CLA	B	1231	-	64,68,73	1.47	6 (9%)	76,107,113	1.97	15 (19%)
29	LUT	1	502	-	42,43,43	2.50	2 (4%)	51,60,60	2.06	11 (21%)
20	CLA	B	1023	-	69,73,73	1.42	7 (10%)	82,113,113	1.84	15 (18%)
20	CLA	B	1226	-	69,73,73	1.41	6 (8%)	82,113,113	1.99	17 (20%)
20	CLA	B	1235	-	69,73,73	1.38	6 (8%)	82,113,113	1.90	19 (23%)
20	CLA	B	1237	-	69,73,73	1.36	6 (8%)	82,113,113	2.00	21 (25%)
20	CLA	A	1013	-	69,73,73	1.43	6 (8%)	82,113,113	1.83	20 (24%)
20	CLA	G	1602	-	50,54,73	1.65	6 (12%)	59,90,113	2.04	17 (28%)
20	CLA	4	608	-	50,54,73	1.58	7 (14%)	59,90,113	2.11	17 (28%)
24	LMT	G	5005	-	32,32,36	1.20	4 (12%)	43,43,47	1.03	2 (4%)
20	CLA	K	1401	-	49,53,73	1.60	6 (12%)	58,89,113	2.09	15 (25%)
22	BCR	B	4005	-	41,41,41	1.64	4 (9%)	56,56,56	4.48	17 (30%)
20	CLA	G	1603	-	69,73,73	1.37	6 (8%)	82,113,113	2.06	20 (24%)
30	CHL	3	604	-	60,74,74	1.48	7 (11%)	58,114,114	1.67	10 (17%)
20	CLA	3	602	-	56,60,73	1.52	7 (12%)	65,97,113	2.11	19 (29%)
24	LMT	A	5004	-	36,36,36	1.19	7 (19%)	47,47,47	1.09	3 (6%)
26	LMG	A	5006	-	50,50,55	1.17	5 (10%)	58,58,63	1.20	3 (5%)
20	CLA	B	1210	-	69,73,73	1.43	6 (8%)	82,113,113	1.99	21 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
22	BCR	A	4007	-	41,41,41	1.67	6 (14%)	56,56,56	4.54	16 (28%)
24	LMT	B	5008	-	32,32,36	1.28	6 (18%)	43,43,47	1.04	3 (6%)
22	BCR	3	503	-	41,41,41	1.61	4 (9%)	56,56,56	4.57	18 (32%)
22	BCR	L	4020	-	41,41,41	1.62	4 (9%)	56,56,56	4.61	15 (26%)
20	CLA	B	1202	-	69,73,73	1.36	6 (8%)	82,113,113	1.92	21 (25%)
20	CLA	4	606	-	54,58,73	1.54	6 (11%)	64,95,113	2.04	19 (29%)
20	CLA	L	1503	-	54,58,73	1.53	7 (12%)	64,95,113	2.08	20 (31%)
20	CLA	K	1404	11	50,54,73	1.59	7 (14%)	59,90,113	2.02	14 (23%)
26	LMG	G	5001	-	49,49,55	1.12	4 (8%)	57,57,63	1.18	6 (10%)
20	CLA	1	603	-	59,63,73	1.49	6 (10%)	70,101,113	2.10	20 (28%)
20	CLA	3	610	15	69,73,73	1.37	6 (8%)	82,113,113	1.88	21 (25%)
26	LMG	F	5004	-	34,34,55	0.46	0	42,42,63	1.13	3 (7%)
28	DGD	1	803	-	42,42,67	0.87	1 (2%)	56,56,81	1.06	3 (5%)
26	LMG	F	5001	-	30,30,55	0.53	0	38,38,63	1.18	2 (5%)
28	DGD	B	5005	-	62,62,67	1.21	6 (9%)	76,76,81	1.05	5 (6%)
20	CLA	1	614	13	64,68,73	1.43	6 (9%)	76,107,113	1.99	18 (23%)
20	CLA	B	1222	36	69,73,73	1.41	6 (8%)	82,113,113	1.94	21 (25%)
20	CLA	A	1104	-	69,73,73	1.42	6 (8%)	82,113,113	1.93	19 (23%)
20	CLA	B	1205	-	69,73,73	1.39	6 (8%)	82,113,113	1.98	19 (23%)
20	CLA	A	1123	-	69,73,73	1.38	6 (8%)	82,113,113	1.89	19 (23%)
26	LMG	1	802	-	46,46,55	1.05	3 (6%)	54,54,63	1.08	2 (3%)
22	BCR	B	4010	-	41,41,41	1.66	5 (12%)	56,56,56	4.43	16 (28%)
20	CLA	B	1227	-	69,73,73	1.39	6 (8%)	82,113,113	1.91	20 (24%)
20	CLA	3	605	-	59,63,73	1.51	6 (10%)	70,101,113	2.09	21 (30%)
20	CLA	B	1232	-	59,63,73	1.52	6 (10%)	70,101,113	2.04	22 (31%)
30	CHL	1	612	-	55,69,74	1.77	9 (16%)	52,108,114	1.80	12 (23%)
20	CLA	2	612	-	59,63,73	1.48	6 (10%)	70,101,113	2.00	15 (21%)
20	CLA	4	605	-	64,68,73	1.48	6 (9%)	76,107,113	1.89	20 (26%)
22	BCR	B	4009	-	41,41,41	1.62	4 (9%)	56,56,56	4.36	12 (21%)
28	DGD	J	5001	-	59,59,67	1.16	6 (10%)	73,73,81	1.03	2 (2%)
20	CLA	A	1116	-	60,64,73	1.51	6 (10%)	71,102,113	2.00	17 (23%)
20	CLA	A	1139	-	69,73,73	1.41	6 (8%)	82,113,113	1.91	16 (19%)
20	CLA	3	613	-	50,54,73	1.59	6 (12%)	59,90,113	2.07	15 (25%)
22	BCR	F	4014	-	41,41,41	1.64	5 (12%)	56,56,56	4.43	16 (28%)
20	CLA	B	1236	-	54,58,73	1.59	6 (11%)	64,95,113	2.10	19 (29%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
20	CLA	B	1224	-	69,73,73	1.39	6 (8%)	82,113,113	1.93	17 (20%)
26	LMG	2	806	-	13,13,55	0.57	0	18,18,63	0.55	0
28	DGD	G	5003	-	48,48,67	0.86	2 (4%)	62,62,81	1.02	2 (3%)
20	CLA	B	1215	-	69,73,73	1.40	6 (8%)	82,113,113	1.98	20 (24%)
29	LUT	1	501	-	42,43,43	2.47	2 (4%)	51,60,60	1.91	16 (31%)
21	PQN	A	2001	-	34,34,34	0.39	0	43,45,45	1.27	4 (9%)
20	CLA	A	1117	-	69,73,73	1.41	6 (8%)	82,113,113	1.82	15 (18%)
23	LHG	4	801	-	34,34,48	0.42	0	37,40,54	1.16	3 (8%)
20	CLA	2	602	-	56,60,73	1.50	7 (12%)	65,97,113	2.04	18 (27%)
20	CLA	B	1213	-	64,68,73	1.45	6 (9%)	76,107,113	1.91	17 (22%)
23	LHG	B	5002	-	48,48,48	0.40	0	51,54,54	1.12	3 (5%)
23	LHG	1	801	-	48,48,48	0.45	1 (2%)	51,54,54	1.21	4 (7%)
20	CLA	B	1221	-	69,73,73	1.42	6 (8%)	82,113,113	1.94	20 (24%)
20	CLA	B	1214	-	69,73,73	1.42	6 (8%)	82,113,113	1.94	18 (21%)
26	LMG	F	5002	-	47,47,55	1.09	4 (8%)	55,55,63	1.22	4 (7%)
29	LUT	3	501	-	42,43,43	2.44	2 (4%)	51,60,60	2.08	19 (37%)
20	CLA	J	1901	10	54,58,73	1.54	7 (12%)	64,95,113	2.12	19 (29%)
20	CLA	B	1239	-	69,73,73	1.42	6 (8%)	82,113,113	1.89	21 (25%)
20	CLA	4	607	-	64,68,73	1.41	8 (12%)	76,107,113	1.97	19 (25%)
22	BCR	I	4018	-	41,41,41	1.70	5 (12%)	56,56,56	4.02	19 (33%)
20	CLA	L	1501	12	54,58,73	1.52	7 (12%)	64,95,113	2.11	18 (28%)
20	CLA	A	1107	1	69,73,73	1.39	6 (8%)	82,113,113	1.96	19 (23%)
20	CLA	B	1217	-	50,54,73	1.63	6 (12%)	59,90,113	2.07	15 (25%)
26	LMG	2	804	-	30,30,55	0.55	0	38,38,63	1.15	4 (10%)
20	CLA	G	1601	-	59,63,73	1.55	6 (10%)	70,101,113	1.96	17 (24%)
20	CLA	B	1230	-	62,66,73	1.46	6 (9%)	73,104,113	2.04	20 (27%)
20	CLA	A	1133	-	69,73,73	1.40	7 (10%)	82,113,113	1.81	16 (19%)
23	LHG	A	5002	-	48,48,48	0.38	0	51,54,54	1.14	5 (9%)
26	LMG	B	5004	-	33,33,55	0.55	0	41,41,63	1.33	6 (14%)
30	CHL	3	611	-	41,55,74	1.80	9 (21%)	35,91,114	2.08	11 (31%)
20	CLA	B	1216	-	69,73,73	1.37	6 (8%)	82,113,113	1.83	17 (20%)
25	SF4	C	3003	3	0,12,12	-	-	-	-	-
30	CHL	4	611	-	45,59,74	1.58	9 (20%)	40,96,114	2.06	11 (27%)
22	BCR	A	4008	-	41,41,41	1.69	6 (14%)	56,56,56	4.21	19 (33%)
20	CLA	B	1208	-	64,68,73	1.42	6 (9%)	76,107,113	1.88	15 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
20	CLA	F	1302	6	69,73,73	1.38	6 (8%)	82,113,113	1.85	15 (18%)
22	BCR	L	4019	-	41,41,41	1.62	4 (9%)	56,56,56	4.53	15 (26%)
22	BCR	I	4020	-	41,41,41	1.62	4 (9%)	56,56,56	4.51	14 (25%)
20	CLA	2	604	14	69,73,73	1.35	7 (10%)	82,113,113	1.90	19 (23%)
28	DGD	F	5005	-	58,58,67	1.13	4 (6%)	72,72,81	1.13	6 (8%)
22	BCR	A	4011	-	41,41,41	1.68	5 (12%)	56,56,56	4.46	19 (33%)
21	PQN	B	2002	-	34,34,34	0.43	0	43,45,45	1.30	5 (11%)
20	CLA	4	612	-	69,73,73	1.35	7 (10%)	82,113,113	1.91	18 (21%)
20	CLA	A	1115	-	69,73,73	1.37	6 (8%)	82,113,113	1.93	16 (19%)
20	CLA	A	1110	-	59,63,73	1.52	6 (10%)	70,101,113	2.04	21 (30%)
22	BCR	1	504	-	41,41,41	1.66	4 (9%)	56,56,56	4.81	15 (26%)
22	BCR	B	4004	-	41,41,41	1.64	4 (9%)	56,56,56	4.97	23 (41%)
22	BCR	K	4001	-	41,41,41	1.63	4 (9%)	56,56,56	4.53	15 (26%)
34	FES	N	101	17	0,4,4	-	-	-	-	-
20	CLA	A	1012	-	69,73,73	1.43	6 (8%)	82,113,113	1.95	17 (20%)
20	CLA	B	1225	-	69,73,73	1.41	6 (8%)	82,113,113	1.88	18 (21%)
20	CLA	1	613	-	49,53,73	1.67	6 (12%)	58,89,113	1.97	13 (22%)
20	CLA	2	606	-	54,58,73	1.53	6 (11%)	64,95,113	2.12	20 (31%)
24	LMT	G	5004	-	36,36,36	1.18	6 (16%)	47,47,47	1.06	3 (6%)
20	CLA	A	1105	-	64,68,73	1.44	6 (9%)	76,107,113	2.00	16 (21%)
20	CLA	A	1121	-	64,68,73	1.45	6 (9%)	76,107,113	1.99	18 (23%)
30	CHL	2	611	-	42,56,74	1.60	9 (21%)	36,92,114	2.01	9 (25%)
29	LUT	2	501	-	42,43,43	2.41	2 (4%)	51,60,60	2.34	17 (33%)
20	CLA	3	617	-	64,68,73	1.39	6 (9%)	76,107,113	1.95	19 (25%)
20	CLA	A	1125	-	69,73,73	1.40	6 (8%)	82,113,113	1.97	20 (24%)
20	CLA	L	1502	-	64,68,73	1.39	6 (9%)	76,107,113	1.98	20 (26%)
22	BCR	B	4006	-	41,41,41	1.65	4 (9%)	56,56,56	4.38	21 (37%)
23	LHG	2	801	20	34,34,48	0.45	0	37,40,54	1.15	4 (10%)
20	CLA	A	1108	-	54,58,73	1.60	6 (11%)	64,95,113	2.05	18 (28%)
22	BCR	A	4003	-	41,41,41	1.65	4 (9%)	56,56,56	4.39	18 (32%)
20	CLA	B	1238	36	69,73,73	1.34	7 (10%)	82,113,113	1.92	18 (21%)
32	3PH	2	807	-	32,32,47	1.03	4 (12%)	35,37,52	1.23	2 (5%)
20	CLA	2	603	14	69,73,73	1.37	8 (11%)	82,113,113	1.91	19 (23%)
22	BCR	J	4012	-	41,41,41	1.65	4 (9%)	56,56,56	4.38	18 (32%)
22	BCR	K	4002	-	41,41,41	1.61	4 (9%)	56,56,56	4.63	18 (32%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
20	CLA	A	1127	-	69,73,73	1.42	7 (10%)	82,113,113	1.86	16 (19%)
25	SF4	C	3002	3	0,12,12	-	-	-		
22	BCR	H	4021	-	41,41,41	1.62	4 (9%)	56,56,56	4.57	15 (26%)
31	XAT	2	502	-	41,47,47	0.79	1 (2%)	54,74,74	2.14	13 (24%)
20	CLA	3	606	-	54,58,73	1.58	7 (12%)	64,95,113	2.01	13 (20%)
20	CLA	A	1141	23	64,68,73	1.41	6 (9%)	76,107,113	1.93	19 (25%)
20	CLA	B	1240	23	69,73,73	1.42	6 (8%)	82,113,113	1.99	21 (25%)
20	CLA	B	1021	-	69,73,73	1.43	7 (10%)	82,113,113	1.92	18 (21%)
20	CLA	F	1301	-	69,73,73	1.38	6 (8%)	82,113,113	1.88	17 (20%)
20	CLA	4	609	16	54,58,73	1.52	6 (11%)	64,95,113	2.10	19 (29%)
20	CLA	K	1402	-	64,68,73	1.40	7 (10%)	76,107,113	2.00	21 (27%)
20	CLA	1	611	-	69,73,73	1.38	6 (8%)	82,113,113	1.90	17 (20%)
20	CLA	A	1122	-	69,73,73	1.38	6 (8%)	82,113,113	1.88	17 (20%)
20	CLA	B	1234	-	59,63,73	1.54	6 (10%)	70,101,113	1.97	14 (20%)
30	CHL	2	610	-	50,64,74	1.36	9 (18%)	46,102,114	1.86	11 (23%)
20	CLA	B	1211	-	69,73,73	1.39	6 (8%)	82,113,113	1.95	19 (23%)
20	CLA	2	601	-	64,68,73	1.45	6 (9%)	76,107,113	1.89	17 (22%)
20	CLA	A	1109	-	69,73,73	1.39	6 (8%)	82,113,113	1.91	20 (24%)
23	LHG	B	5001	20	20,20,48	0.60	0	23,26,54	1.53	2 (8%)
20	CLA	B	1223	-	69,73,73	1.40	6 (8%)	82,113,113	1.94	17 (20%)
20	CLA	A	1124	-	59,63,73	1.51	6 (10%)	70,101,113	2.01	18 (25%)
28	DGD	4	802	-	52,52,67	0.93	2 (3%)	66,66,81	1.04	3 (4%)
20	CLA	B	1209	-	50,54,73	1.59	6 (12%)	59,90,113	2.02	12 (20%)
24	LMT	2	808	-	36,36,36	0.42	0	47,47,47	0.74	1 (2%)
20	CLA	A	1130	-	59,63,73	1.49	6 (10%)	70,101,113	1.95	17 (24%)
20	CLA	3	608	-	52,56,73	1.55	7 (13%)	61,92,113	2.09	19 (31%)
20	CLA	A	1119	-	69,73,73	1.42	6 (8%)	82,113,113	1.89	19 (23%)
30	CHL	1	609	13	50,64,74	1.47	9 (18%)	46,102,114	1.84	11 (23%)
30	CHL	2	613	-	40,54,74	1.43	9 (22%)	34,90,114	2.09	8 (23%)
20	CLA	1	605	-	69,73,73	1.38	6 (8%)	82,113,113	2.04	18 (21%)
30	CHL	2	615	-	50,64,74	1.25	9 (18%)	46,102,114	1.93	13 (28%)
20	CLA	A	1118	-	54,58,73	1.57	6 (11%)	64,95,113	2.20	20 (31%)
20	CLA	A	1137	-	69,73,73	1.38	6 (8%)	82,113,113	1.91	17 (20%)
20	CLA	B	1212	-	59,63,73	1.50	6 (10%)	70,101,113	2.04	17 (24%)
22	BCR	2	503	-	41,41,41	1.73	5 (12%)	56,56,56	5.47	26 (46%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
20	CLA	A	1126	-	69,73,73	1.41	6 (8%)	82,113,113	1.90	17 (20%)
30	CHL	3	607	23	45,59,74	1.68	10 (22%)	40,96,114	1.96	12 (30%)
25	SF4	A	3001	2,1	0,12,12	-	-	-		
23	LHG	A	5001	20	39,39,48	0.46	0	42,45,54	1.27	4 (9%)
20	CLA	B	1206	2	69,73,73	1.42	6 (8%)	82,113,113	1.88	17 (20%)
20	CLA	A	1140	-	69,73,73	1.39	6 (8%)	82,113,113	1.87	21 (25%)
26	LMG	3	802	-	30,30,55	0.57	0	38,38,63	1.15	3 (7%)
20	CLA	4	603	-	69,73,73	1.34	7 (10%)	82,113,113	1.81	18 (21%)
29	LUT	3	502	-	42,43,43	2.41	1 (2%)	51,60,60	2.07	10 (19%)
24	LMT	B	5006	-	33,33,36	1.21	6 (18%)	44,44,47	1.01	3 (6%)
22	BCR	A	4017	-	41,41,41	1.64	5 (12%)	56,56,56	4.92	19 (33%)
20	CLA	B	1201	-	69,73,73	1.38	6 (8%)	82,113,113	1.87	16 (19%)
20	CLA	A	1101	-	69,73,73	1.35	6 (8%)	82,113,113	1.94	19 (23%)
20	CLA	A	1120	-	64,68,73	1.43	6 (9%)	76,107,113	1.93	16 (21%)
20	CLA	A	1131	-	69,73,73	1.41	6 (8%)	82,113,113	1.88	17 (20%)
23	LHG	3	801	30	16,16,48	0.85	1 (6%)	17,20,54	0.70	1 (5%)
29	LUT	J	4013	-	42,43,43	2.33	2 (4%)	51,60,60	2.39	15 (29%)
20	CLA	A	1102	-	69,73,73	1.37	6 (8%)	82,113,113	1.93	20 (24%)
20	CLA	A	1134	1	59,63,73	1.47	6 (10%)	70,101,113	1.99	17 (24%)
20	CLA	A	1135	-	55,59,73	1.57	6 (10%)	64,96,113	2.08	16 (25%)
20	CLA	B	1204	-	69,73,73	1.39	6 (8%)	82,113,113	1.92	19 (23%)
20	CLA	B	1219	-	69,73,73	1.39	6 (8%)	82,113,113	1.90	16 (19%)
20	CLA	H	1701	8	64,68,73	1.39	6 (9%)	76,107,113	1.94	18 (23%)
20	CLA	1	607	-	50,54,73	1.63	6 (12%)	59,90,113	2.03	15 (25%)
20	CLA	2	608	-	54,58,73	1.59	6 (11%)	64,95,113	2.19	18 (28%)
26	LMG	G	5002	-	50,50,55	1.19	5 (10%)	58,58,63	1.23	5 (8%)
20	CLA	2	605	-	69,73,73	1.38	6 (8%)	82,113,113	1.95	18 (21%)
26	LMG	G	5006	-	25,25,55	0.57	0	33,33,63	1.25	4 (12%)
20	CLA	A	1138	-	69,73,73	1.40	6 (8%)	82,113,113	1.85	15 (18%)
20	CLA	3	603	15	59,63,73	1.47	6 (10%)	70,101,113	2.08	21 (30%)
20	CLA	B	1203	-	69,73,73	1.39	6 (8%)	82,113,113	1.87	20 (24%)
20	CLA	4	602	-	54,58,73	1.53	7 (12%)	64,95,113	2.02	17 (26%)
20	CLA	A	1114	-	50,54,73	1.61	6 (12%)	59,90,113	2.09	16 (27%)

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
20	CLA	A	1111	-	1/1/15/20	18/39/115/115	-
20	CLA	A	1128	-	1/1/15/20	16/39/115/115	-
20	CLA	A	1132	-	1/1/15/20	13/39/115/115	-
22	BCR	G	4011	-	-	8/29/63/63	0/2/2/2
20	CLA	3	614	-	1/1/10/20	3/12/88/115	-
31	XAT	4	502	-	2/2/12/26	2/31/93/93	0/4/4/4
20	CLA	A	1103	-	1/1/15/20	18/39/115/115	-
20	CLA	A	1106	1	1/1/15/20	17/39/115/115	-
30	CHL	1	610	13	3/3/16/26	5/17/115/137	-
26	LMG	B	5003	-	-	8/30/50/70	0/1/1/1
22	BCR	F	4016	-	-	14/29/63/63	0/2/2/2
26	LMG	F	5006	-	-	1/4/24/70	0/1/1/1
20	CLA	4	617	-	1/1/15/20	19/39/115/115	-
28	DGD	3	803	-	-	10/40/80/95	0/2/2/2
20	CLA	B	1220	-	1/1/13/20	10/27/103/115	-
20	CLA	1	602	13	1/1/11/20	8/17/93/115	-
30	CHL	4	610	-	3/3/16/26	3/17/115/137	-
20	CLA	K	1403	-	1/1/11/20	8/19/95/115	-
20	CLA	4	604	-	1/1/14/20	15/33/109/115	-
33	C7Z	4	505	-	-	16/29/67/67	0/2/2/2
22	BCR	3	506	-	-	14/29/63/63	0/2/2/2
20	CLA	1	604	-	1/1/15/20	14/39/115/115	-
26	LMG	B	5007	-	-	11/29/49/70	0/1/1/1
29	LUT	4	501	-	-	4/29/67/67	0/2/2/2
20	CLA	A	1136	-	1/1/15/20	14/39/115/115	-
30	CHL	2	609	-	4/4/20/26	9/39/137/137	-
20	CLA	A	1112	-	1/1/15/20	18/39/115/115	-
22	BCR	A	4002	-	-	11/29/63/63	0/2/2/2
22	BCR	1	503	-	-	16/29/63/63	0/2/2/2
20	CLA	A	1129	-	1/1/15/20	15/39/115/115	-
26	LMG	F	5003	-	-	6/31/51/70	0/1/1/1
26	LMG	2	803	-	-	12/31/51/70	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
20	CLA	B	1207	-	1/1/15/20	22/39/115/115	-
20	CLA	3	612	-	1/1/12/20	9/21/97/115	-
30	CHL	4	613	-	4/4/19/26	8/33/131/137	-
20	CLA	1	606	-	1/1/12/20	5/21/97/115	-
30	CHL	4	615	-	3/3/15/26	2/12/110/137	-
20	CLA	A	1113	-	1/1/11/20	7/15/91/115	-
20	CLA	B	1229	-	1/1/15/20	12/39/115/115	-
26	LMG	2	802	-	-	8/20/40/70	0/1/1/1
20	CLA	B	1228	-	1/1/14/20	11/33/109/115	-
20	CLA	B	1022	-	1/1/15/20	6/39/115/115	-
20	CLA	1	608	-	1/1/11/20	9/17/93/115	-
20	CLA	3	601	-	1/1/13/20	12/27/103/115	-
26	LMG	2	805	-	-	7/25/45/70	0/1/1/1
20	CLA	4	601	16	1/1/14/20	14/33/109/115	-
20	CLA	1	601	13	1/1/15/20	18/39/115/115	-
20	CLA	2	607	23	1/1/14/20	15/33/109/115	-
19	CL0	A	1011	-	3/3/20/25	8/37/135/135	-
20	CLA	B	1218	-	1/1/15/20	8/39/115/115	-
20	CLA	B	1231	-	1/1/14/20	10/33/109/115	-
29	LUT	1	502	-	-	5/29/67/67	0/2/2/2
20	CLA	B	1023	-	1/1/15/20	10/39/115/115	-
20	CLA	B	1226	-	1/1/15/20	21/39/115/115	-
20	CLA	B	1235	-	1/1/15/20	15/39/115/115	-
20	CLA	B	1237	-	1/1/15/20	18/39/115/115	-
20	CLA	A	1013	-	1/1/15/20	16/39/115/115	-
20	CLA	G	1602	-	1/1/11/20	6/17/93/115	-
20	CLA	4	608	-	1/1/11/20	5/17/93/115	-
24	LMT	G	5005	-	-	6/17/57/61	0/2/2/2
20	CLA	K	1401	-	1/1/11/20	7/15/91/115	-
22	BCR	B	4005	-	-	10/29/63/63	0/2/2/2
20	CLA	G	1603	-	1/1/15/20	17/39/115/115	-
30	CHL	3	604	-	4/4/20/26	9/39/137/137	-
20	CLA	3	602	-	1/1/12/20	11/24/100/115	-
24	LMT	A	5004	-	-	9/21/61/61	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
26	LMG	A	5006	-	-	11/45/65/70	0/1/1/1
20	CLA	B	1210	-	1/1/15/20	14/39/115/115	-
22	BCR	A	4007	-	-	9/29/63/63	0/2/2/2
24	LMT	B	5008	-	-	8/17/57/61	0/2/2/2
22	BCR	3	503	-	-	15/29/63/63	0/2/2/2
22	BCR	L	4020	-	-	9/29/63/63	0/2/2/2
20	CLA	B	1202	-	1/1/15/20	14/39/115/115	-
20	CLA	4	606	-	1/1/12/20	6/21/97/115	-
20	CLA	L	1503	-	1/1/12/20	10/21/97/115	-
20	CLA	K	1404	11	1/1/11/20	7/17/93/115	-
26	LMG	G	5001	-	-	21/44/64/70	0/1/1/1
20	CLA	1	603	-	1/1/13/20	7/27/103/115	-
20	CLA	3	610	15	1/1/15/20	21/39/115/115	-
26	LMG	F	5004	-	-	8/29/49/70	0/1/1/1
28	DGD	1	803	-	-	15/30/70/95	0/2/2/2
26	LMG	F	5001	-	-	6/25/45/70	0/1/1/1
28	DGD	B	5005	-	-	20/50/90/95	0/2/2/2
20	CLA	1	614	13	1/1/14/20	17/33/109/115	-
20	CLA	B	1222	36	1/1/15/20	21/39/115/115	-
20	CLA	A	1104	-	1/1/15/20	15/39/115/115	-
20	CLA	B	1205	-	1/1/15/20	17/39/115/115	-
20	CLA	A	1123	-	1/1/15/20	14/39/115/115	-
26	LMG	1	802	-	-	8/41/61/70	0/1/1/1
30	CHL	1	612	-	4/4/19/26	7/33/131/137	-
20	CLA	B	1227	-	1/1/15/20	14/39/115/115	-
20	CLA	3	605	-	1/1/13/20	10/27/103/115	-
20	CLA	B	1232	-	1/1/13/20	15/27/103/115	-
22	BCR	B	4010	-	-	14/29/63/63	0/2/2/2
20	CLA	2	612	-	1/1/13/20	10/27/103/115	-
20	CLA	4	605	-	-	14/33/109/115	-
22	BCR	B	4009	-	-	9/29/63/63	0/2/2/2
28	DGD	J	5001	-	-	14/47/87/95	0/2/2/2
20	CLA	A	1116	-	1/1/13/20	11/29/105/115	-
20	CLA	A	1139	-	1/1/15/20	12/39/115/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
20	CLA	3	613	-	1/1/11/20	7/17/93/115	-
22	BCR	F	4014	-	-	9/29/63/63	0/2/2/2
20	CLA	B	1236	-	1/1/12/20	8/21/97/115	-
20	CLA	B	1224	-	1/1/15/20	17/39/115/115	-
26	LMG	2	806	-	-	1/4/24/70	0/1/1/1
28	DGD	G	5003	-	-	12/36/76/95	0/2/2/2
20	CLA	B	1215	-	1/1/15/20	16/39/115/115	-
29	LUT	1	501	-	-	3/29/67/67	0/2/2/2
21	PQN	A	2001	-	-	3/23/43/43	0/2/2/2
20	CLA	A	1117	-	1/1/15/20	17/39/115/115	-
23	LHG	4	801	-	-	19/39/39/53	-
20	CLA	2	602	-	1/1/12/20	10/24/100/115	-
20	CLA	B	1213	-	1/1/14/20	7/33/109/115	-
23	LHG	B	5002	-	-	29/53/53/53	-
23	LHG	1	801	-	-	28/53/53/53	-
20	CLA	B	1221	-	1/1/15/20	21/39/115/115	-
20	CLA	B	1214	-	1/1/15/20	16/39/115/115	-
26	LMG	F	5002	-	-	10/42/62/70	0/1/1/1
29	LUT	3	501	-	-	5/29/67/67	0/2/2/2
20	CLA	J	1901	10	1/1/12/20	7/21/97/115	-
20	CLA	B	1239	-	1/1/15/20	18/39/115/115	-
20	CLA	4	607	-	1/1/14/20	13/33/109/115	-
22	BCR	I	4018	-	-	13/29/63/63	0/2/2/2
20	CLA	L	1501	12	1/1/12/20	10/21/97/115	-
20	CLA	A	1107	1	1/1/15/20	16/39/115/115	-
20	CLA	B	1217	-	1/1/11/20	9/17/93/115	-
26	LMG	2	804	-	-	7/25/45/70	0/1/1/1
20	CLA	G	1601	-	1/1/13/20	9/27/103/115	-
20	CLA	B	1230	-	1/1/13/20	14/31/107/115	-
20	CLA	A	1133	-	1/1/15/20	16/39/115/115	-
23	LHG	A	5002	-	-	35/53/53/53	-
26	LMG	B	5004	-	-	13/28/48/70	0/1/1/1
30	CHL	3	611	-	3/3/16/26	0/17/115/137	-
20	CLA	B	1216	-	1/1/15/20	11/39/115/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
25	SF4	C	3003	3	-	-	0/6/5/5
30	CHL	4	611	-	3/3/17/26	2/21/119/137	-
22	BCR	A	4008	-	-	12/29/63/63	0/2/2/2
20	CLA	B	1208	-	1/1/14/20	13/33/109/115	-
20	CLA	F	1302	6	1/1/15/20	17/39/115/115	-
22	BCR	L	4019	-	-	14/29/63/63	0/2/2/2
22	BCR	I	4020	-	-	15/29/63/63	0/2/2/2
20	CLA	2	604	14	1/1/15/20	17/39/115/115	-
28	DGD	F	5005	-	-	19/46/86/95	0/2/2/2
22	BCR	A	4011	-	-	15/29/63/63	0/2/2/2
21	PQN	B	2002	-	-	6/23/43/43	0/2/2/2
20	CLA	4	612	-	1/1/15/20	17/39/115/115	-
20	CLA	A	1115	-	1/1/15/20	19/39/115/115	-
20	CLA	A	1110	-	1/1/13/20	6/27/103/115	-
22	BCR	1	504	-	-	14/29/63/63	0/2/2/2
22	BCR	B	4004	-	-	12/29/63/63	0/2/2/2
22	BCR	K	4001	-	-	12/29/63/63	0/2/2/2
34	FES	N	101	17	-	-	0/1/1/1
20	CLA	A	1012	-	1/1/15/20	21/39/115/115	-
20	CLA	B	1225	-	1/1/15/20	14/39/115/115	-
20	CLA	1	613	-	1/1/11/20	6/15/91/115	-
20	CLA	2	606	-	1/1/12/20	9/21/97/115	-
24	LMT	G	5004	-	-	10/21/61/61	0/2/2/2
20	CLA	A	1105	-	1/1/14/20	18/33/109/115	-
20	CLA	A	1121	-	1/1/14/20	16/33/109/115	-
30	CHL	2	611	-	3/3/16/26	4/18/116/137	-
29	LUT	2	501	-	1/1/12/27	3/29/67/67	0/2/2/2
20	CLA	3	617	-	1/1/14/20	12/33/109/115	-
20	CLA	A	1125	-	1/1/15/20	16/39/115/115	-
20	CLA	L	1502	-	1/1/14/20	21/33/109/115	-
22	BCR	B	4006	-	-	11/29/63/63	0/2/2/2
23	LHG	2	801	20	-	19/39/39/53	-
20	CLA	A	1108	-	1/1/12/20	5/21/97/115	-
22	BCR	A	4003	-	-	13/29/63/63	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
20	CLA	B	1238	36	1/1/15/20	16/39/115/115	-
32	3PH	2	807	-	-	16/34/34/49	-
20	CLA	2	603	14	1/1/15/20	13/39/115/115	-
22	BCR	J	4012	-	-	12/29/63/63	0/2/2/2
22	BCR	K	4002	-	-	11/29/63/63	0/2/2/2
20	CLA	A	1127	-	1/1/15/20	16/39/115/115	-
25	SF4	C	3002	3	-	-	0/6/5/5
22	BCR	H	4021	-	-	12/29/63/63	0/2/2/2
31	XAT	2	502	-	2/2/12/26	4/31/93/93	0/4/4/4
20	CLA	3	606	-	1/1/12/20	10/21/97/115	-
20	CLA	A	1141	23	1/1/14/20	14/33/109/115	-
20	CLA	B	1240	23	1/1/15/20	18/39/115/115	-
20	CLA	B	1021	-	1/1/15/20	12/39/115/115	-
20	CLA	F	1301	-	1/1/15/20	17/39/115/115	-
20	CLA	4	609	16	1/1/12/20	9/21/97/115	-
20	CLA	K	1402	-	1/1/14/20	20/33/109/115	-
20	CLA	1	611	-	1/1/15/20	14/39/115/115	-
20	CLA	A	1122	-	1/1/15/20	15/39/115/115	-
20	CLA	B	1234	-	1/1/13/20	6/27/103/115	-
30	CHL	2	610	-	4/4/18/26	6/27/125/137	-
20	CLA	B	1211	-	1/1/15/20	21/39/115/115	-
20	CLA	2	601	-	1/1/14/20	14/33/109/115	-
20	CLA	A	1109	-	1/1/15/20	13/39/115/115	-
23	LHG	B	5001	20	-	14/23/23/53	-
20	CLA	B	1223	-	1/1/15/20	13/39/115/115	-
20	CLA	A	1124	-	1/1/13/20	8/27/103/115	-
28	DGD	4	802	-	-	16/40/80/95	0/2/2/2
20	CLA	B	1209	-	1/1/11/20	6/17/93/115	-
24	LMT	2	808	-	-	2/21/61/61	0/2/2/2
20	CLA	A	1130	-	1/1/13/20	12/27/103/115	-
20	CLA	3	608	-	1/1/11/20	4/19/95/115	-
20	CLA	A	1119	-	1/1/15/20	14/39/115/115	-
30	CHL	1	609	13	4/4/18/26	0/27/125/137	-
30	CHL	2	613	-	3/3/16/26	8/15/113/137	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
20	CLA	1	605	-	1/1/15/20	17/39/115/115	-
30	CHL	2	615	-	4/4/18/26	5/27/125/137	-
20	CLA	A	1118	-	1/1/12/20	8/21/97/115	-
20	CLA	A	1137	-	1/1/15/20	18/39/115/115	-
20	CLA	B	1212	-	1/1/13/20	9/27/103/115	-
22	BCR	2	503	-	-	14/29/63/63	0/2/2/2
20	CLA	A	1126	-	1/1/15/20	20/39/115/115	-
30	CHL	3	607	23	3/3/17/26	5/21/119/137	-
25	SF4	A	3001	2,1	-	-	0/6/5/5
23	LHG	A	5001	20	-	22/44/44/53	-
20	CLA	B	1206	2	1/1/15/20	17/39/115/115	-
20	CLA	A	1140	-	1/1/15/20	13/39/115/115	-
29	LUT	3	502	-	1/1/12/27	9/29/67/67	0/2/2/2
20	CLA	4	603	-	1/1/15/20	16/39/115/115	-
26	LMG	3	802	-	-	11/25/45/70	0/1/1/1
24	LMT	B	5006	-	-	7/18/58/61	0/2/2/2
22	BCR	A	4017	-	-	6/29/63/63	0/2/2/2
20	CLA	B	1201	-	1/1/15/20	17/39/115/115	-
20	CLA	A	1101	-	1/1/15/20	16/39/115/115	-
20	CLA	A	1120	-	1/1/14/20	15/33/109/115	-
20	CLA	A	1131	-	1/1/15/20	13/39/115/115	-
23	LHG	3	801	30	-	12/19/19/53	-
29	LUT	J	4013	-	-	7/29/67/67	0/2/2/2
20	CLA	A	1102	-	1/1/15/20	24/39/115/115	-
20	CLA	A	1134	1	1/1/13/20	13/27/103/115	-
20	CLA	A	1135	-	1/1/12/20	9/23/99/115	-
20	CLA	B	1204	-	1/1/15/20	20/39/115/115	-
20	CLA	B	1219	-	1/1/15/20	18/39/115/115	-
20	CLA	H	1701	8	1/1/14/20	15/33/109/115	-
20	CLA	1	607	-	1/1/11/20	5/17/93/115	-
20	CLA	2	608	-	1/1/12/20	8/21/97/115	-
26	LMG	G	5002	-	-	18/45/65/70	0/1/1/1
20	CLA	2	605	-	1/1/15/20	20/39/115/115	-
26	LMG	G	5006	-	-	11/20/40/70	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
20	CLA	A	1138	-	1/1/15/20	14/39/115/115	-
20	CLA	3	603	15	1/1/13/20	13/27/103/115	-
20	CLA	B	1203	-	1/1/15/20	13/39/115/115	-
20	CLA	4	602	-	1/1/12/20	7/21/97/115	-
20	CLA	A	1114	-	1/1/11/20	7/17/93/115	-

The worst 5 of 1256 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
29	1	502	LUT	C24-C25	15.26	1.51	1.33
29	4	501	LUT	C24-C25	14.96	1.50	1.33
29	1	501	LUT	C24-C25	14.95	1.50	1.33
29	3	501	LUT	C24-C25	14.84	1.50	1.33
29	3	502	LUT	C24-C25	14.70	1.50	1.33

The worst 5 of 3472 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	H	4021	BCR	C10-C11-C12	19.50	179.72	123.20
22	3	506	BCR	C10-C11-C12	19.08	178.49	123.20
22	1	504	BCR	C10-C11-C12	18.96	178.15	123.20
22	L	4019	BCR	C10-C11-C12	18.96	178.13	123.20
22	3	503	BCR	C10-C11-C12	18.95	178.09	123.20

5 of 201 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
19	A	1011	CL0	NA
19	A	1011	CL0	ND
19	A	1011	CL0	NC
20	A	1101	CLA	ND
20	A	1102	CLA	ND

5 of 2850 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
20	A	1101	CLA	C2-C1-O2A-CGA
20	A	1102	CLA	C3A-C2A-CAA-CBA
20	A	1102	CLA	CHA-CBD-CGD-O1D
20	A	1102	CLA	CHA-CBD-CGD-O2D
20	A	1103	CLA	C1A-C2A-CAA-CBA

There are no ring outliers.

226 monomers are involved in 1507 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	A	1111	CLA	13	0
20	A	1128	CLA	3	0
20	A	1132	CLA	28	0
22	G	4011	BCR	4	0
31	4	502	XAT	14	0
20	A	1103	CLA	13	0
20	A	1106	CLA	10	0
26	B	5003	LMG	13	0
30	1	610	CHL	10	0
22	F	4016	BCR	9	0
20	4	617	CLA	10	0
28	3	803	DGD	3	0
20	B	1220	CLA	5	0
20	1	602	CLA	2	0
30	4	610	CHL	3	0
20	4	604	CLA	28	0
33	4	505	C7Z	4	0
22	3	506	BCR	2	0
20	1	604	CLA	13	0
26	B	5007	LMG	1	0
29	4	501	LUT	19	0
20	A	1136	CLA	11	0
22	A	4002	BCR	2	0
20	A	1112	CLA	11	0
30	2	609	CHL	22	0
22	1	503	BCR	1	0
20	A	1129	CLA	1	0
26	F	5003	LMG	3	0
26	2	803	LMG	4	0
20	B	1207	CLA	11	0
20	3	612	CLA	9	0
30	4	613	CHL	5	0
20	1	606	CLA	9	0
30	4	615	CHL	3	0
20	A	1113	CLA	18	0
20	B	1229	CLA	6	0
26	2	802	LMG	10	0
20	B	1228	CLA	4	0
20	B	1022	CLA	7	0
20	1	608	CLA	9	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	3	601	CLA	8	0
26	2	805	LMG	1	0
20	4	601	CLA	41	0
20	1	601	CLA	12	0
20	2	607	CLA	5	0
19	A	1011	CL0	11	0
20	B	1218	CLA	7	0
20	B	1231	CLA	6	0
29	1	502	LUT	33	0
20	B	1023	CLA	3	0
20	B	1226	CLA	5	0
20	B	1235	CLA	7	0
20	B	1237	CLA	20	0
20	A	1013	CLA	12	0
20	G	1602	CLA	2	0
20	4	608	CLA	13	0
20	K	1401	CLA	12	0
22	B	4005	BCR	3	0
20	G	1603	CLA	14	0
30	3	604	CHL	2	0
20	3	602	CLA	2	0
24	A	5004	LMT	1	0
26	A	5006	LMG	3	0
20	B	1210	CLA	8	0
22	A	4007	BCR	5	0
24	B	5008	LMT	1	0
22	3	503	BCR	8	0
22	L	4020	BCR	18	0
20	B	1202	CLA	5	0
20	4	606	CLA	3	0
20	L	1503	CLA	17	0
20	K	1404	CLA	3	0
26	G	5001	LMG	5	0
20	1	603	CLA	21	0
20	3	610	CLA	3	0
28	1	803	DGD	1	0
26	F	5001	LMG	2	0
28	B	5005	DGD	7	0
20	1	614	CLA	5	0
20	B	1222	CLA	7	0
20	A	1104	CLA	5	0
20	B	1205	CLA	7	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	A	1123	CLA	9	0
26	1	802	LMG	8	0
22	B	4010	BCR	3	0
20	B	1227	CLA	6	0
20	3	605	CLA	5	0
20	B	1232	CLA	11	0
30	1	612	CHL	5	0
20	2	612	CLA	12	0
20	4	605	CLA	11	0
22	B	4009	BCR	4	0
28	J	5001	DGD	11	0
20	A	1116	CLA	13	0
20	A	1139	CLA	10	0
20	3	613	CLA	4	0
22	F	4014	BCR	3	0
20	B	1236	CLA	1	0
20	B	1224	CLA	5	0
28	G	5003	DGD	3	0
20	B	1215	CLA	5	0
29	1	501	LUT	6	0
21	A	2001	PQN	8	0
20	A	1117	CLA	12	0
23	4	801	LHG	11	0
20	2	602	CLA	2	0
20	B	1213	CLA	3	0
23	B	5002	LHG	4	0
23	1	801	LHG	11	0
20	B	1221	CLA	3	0
20	B	1214	CLA	11	0
26	F	5002	LMG	3	0
29	3	501	LUT	3	0
20	J	1901	CLA	7	0
20	B	1239	CLA	7	0
20	4	607	CLA	10	0
22	I	4018	BCR	8	0
20	L	1501	CLA	14	0
20	A	1107	CLA	9	0
26	2	804	LMG	2	0
20	G	1601	CLA	5	0
20	B	1230	CLA	9	0
20	A	1133	CLA	12	0
23	A	5002	LHG	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
26	B	5004	LMG	2	0
30	3	611	CHL	5	0
20	B	1216	CLA	6	0
25	C	3003	SF4	2	0
30	4	611	CHL	1	0
22	A	4008	BCR	4	0
20	B	1208	CLA	2	0
20	F	1302	CLA	8	0
22	L	4019	BCR	28	0
22	I	4020	BCR	10	0
20	2	604	CLA	26	0
28	F	5005	DGD	18	0
22	A	4011	BCR	7	0
21	B	2002	PQN	3	0
20	4	612	CLA	16	0
20	A	1115	CLA	9	0
20	A	1110	CLA	8	0
22	1	504	BCR	4	0
22	B	4004	BCR	4	0
22	K	4001	BCR	7	0
20	A	1012	CLA	10	0
20	B	1225	CLA	8	0
20	1	613	CLA	8	0
20	2	606	CLA	17	0
20	A	1105	CLA	10	0
20	A	1121	CLA	1	0
30	2	611	CHL	5	0
29	2	501	LUT	12	0
20	3	617	CLA	13	0
20	A	1125	CLA	12	0
20	L	1502	CLA	19	0
22	B	4006	BCR	5	0
23	2	801	LHG	7	0
20	A	1108	CLA	3	0
20	B	1238	CLA	9	0
32	2	807	3PH	5	0
20	2	603	CLA	8	0
22	J	4012	BCR	5	0
22	K	4002	BCR	8	0
20	A	1127	CLA	11	0
25	C	3002	SF4	1	0
22	H	4021	BCR	18	0

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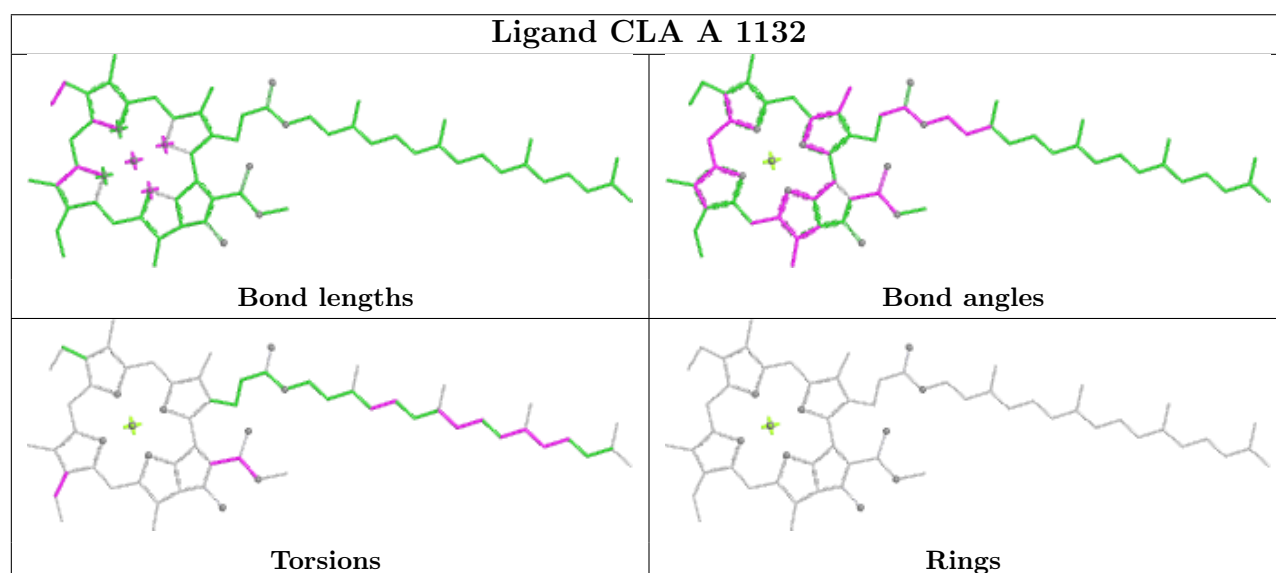
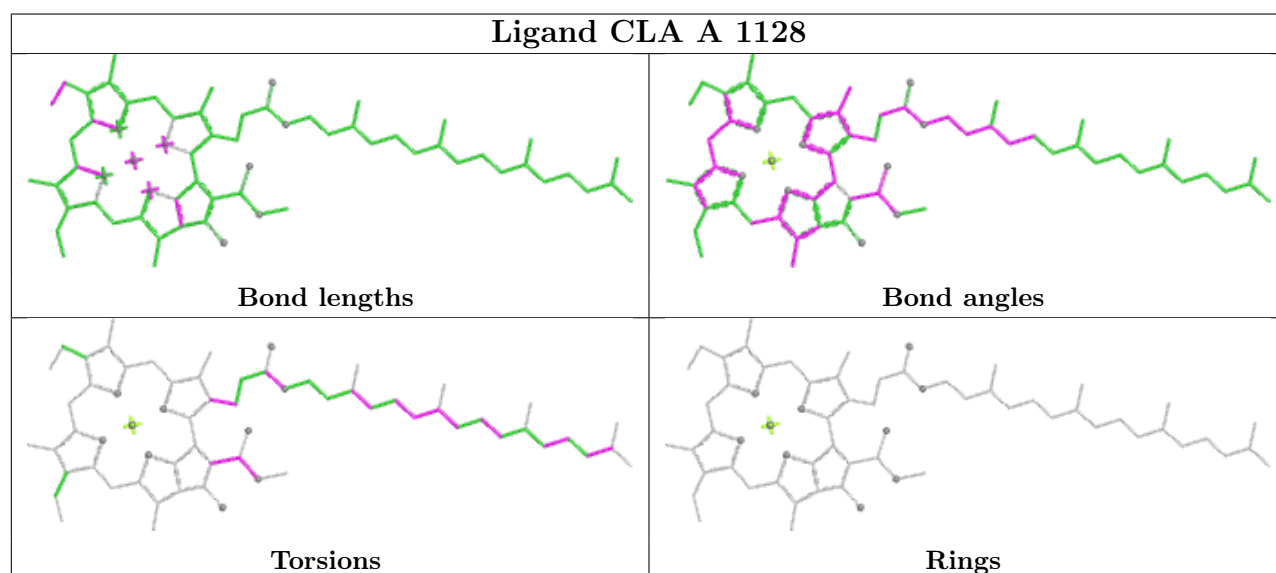
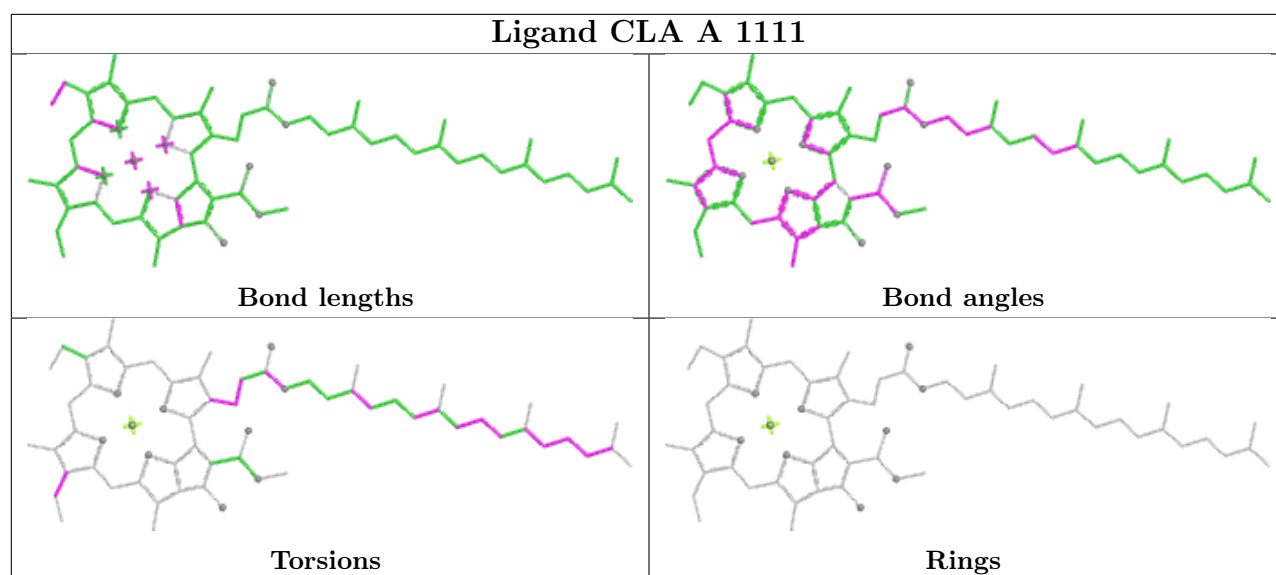
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20	A	1141	CLA	3	0
20	B	1240	CLA	7	0
20	B	1021	CLA	9	0
20	F	1301	CLA	13	0
20	4	609	CLA	17	0
20	K	1402	CLA	1	0
20	1	611	CLA	9	0
20	A	1122	CLA	7	0
20	B	1234	CLA	9	0
30	2	610	CHL	17	0
20	B	1211	CLA	10	0
20	2	601	CLA	9	0
20	A	1109	CLA	12	0
23	B	5001	LHG	2	0
20	B	1223	CLA	7	0
20	A	1124	CLA	4	0
28	4	802	DGD	4	0
24	2	808	LMT	7	0
20	A	1130	CLA	3	0
20	3	608	CLA	5	0
20	A	1119	CLA	12	0
30	1	609	CHL	13	0
30	2	613	CHL	5	0
20	1	605	CLA	8	0
30	2	615	CHL	4	0
20	A	1118	CLA	2	0
20	A	1137	CLA	4	0
20	B	1212	CLA	4	0
22	2	503	BCR	23	0
20	A	1126	CLA	17	0
30	3	607	CHL	1	0
23	A	5001	LHG	1	0
20	B	1206	CLA	10	0
20	A	1140	CLA	7	0
20	4	603	CLA	18	0
29	3	502	LUT	4	0
22	A	4017	BCR	7	0
20	B	1201	CLA	4	0
20	A	1101	CLA	16	0
20	A	1120	CLA	3	0

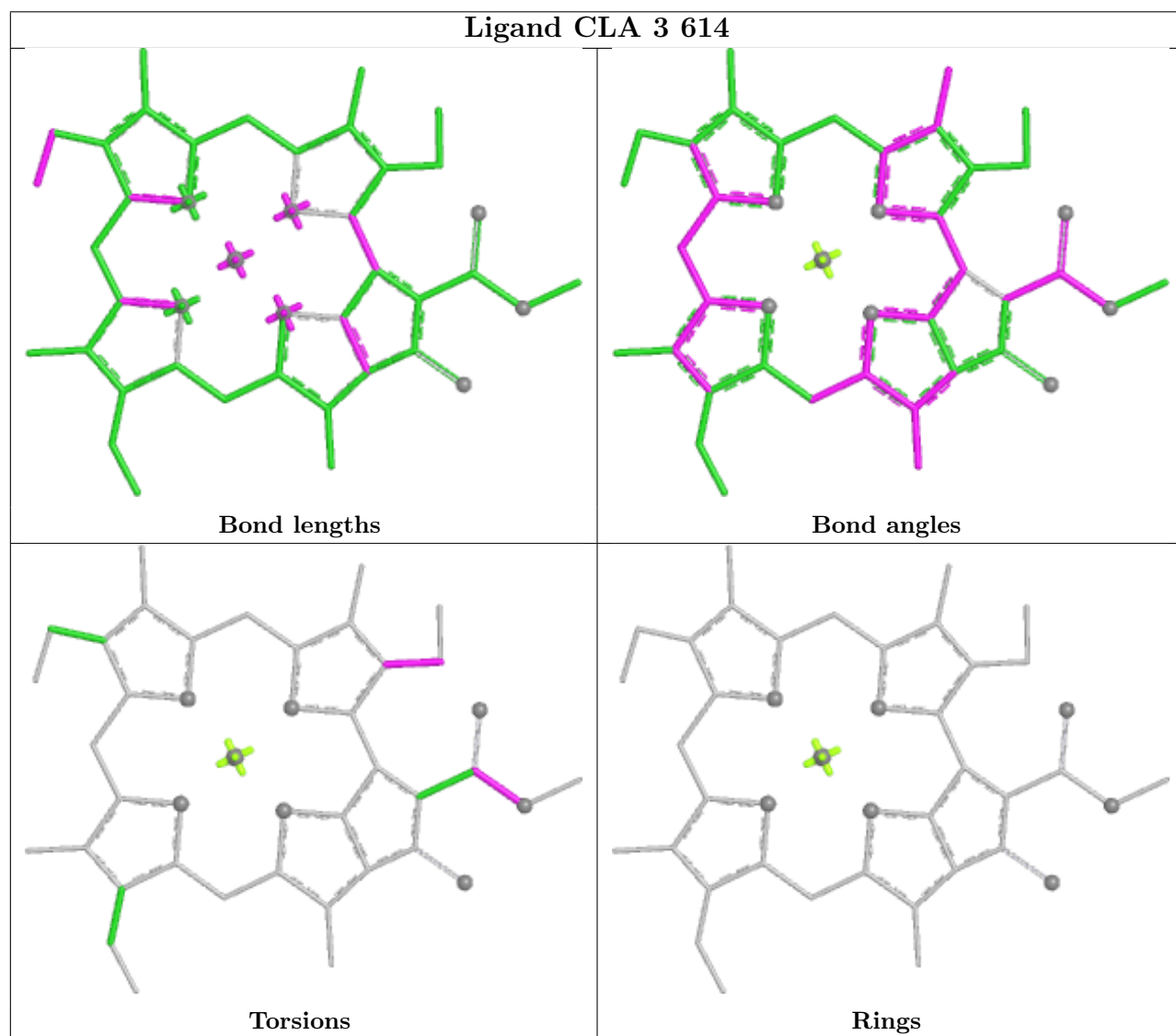
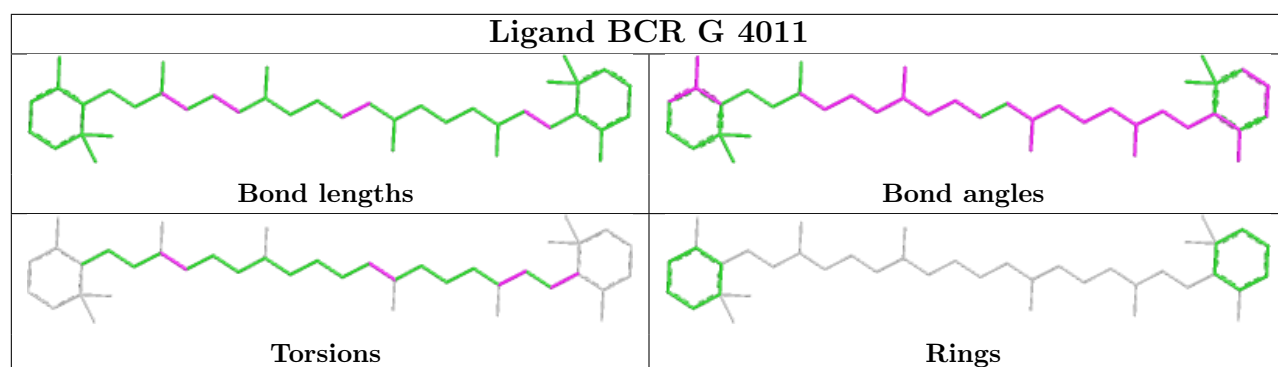
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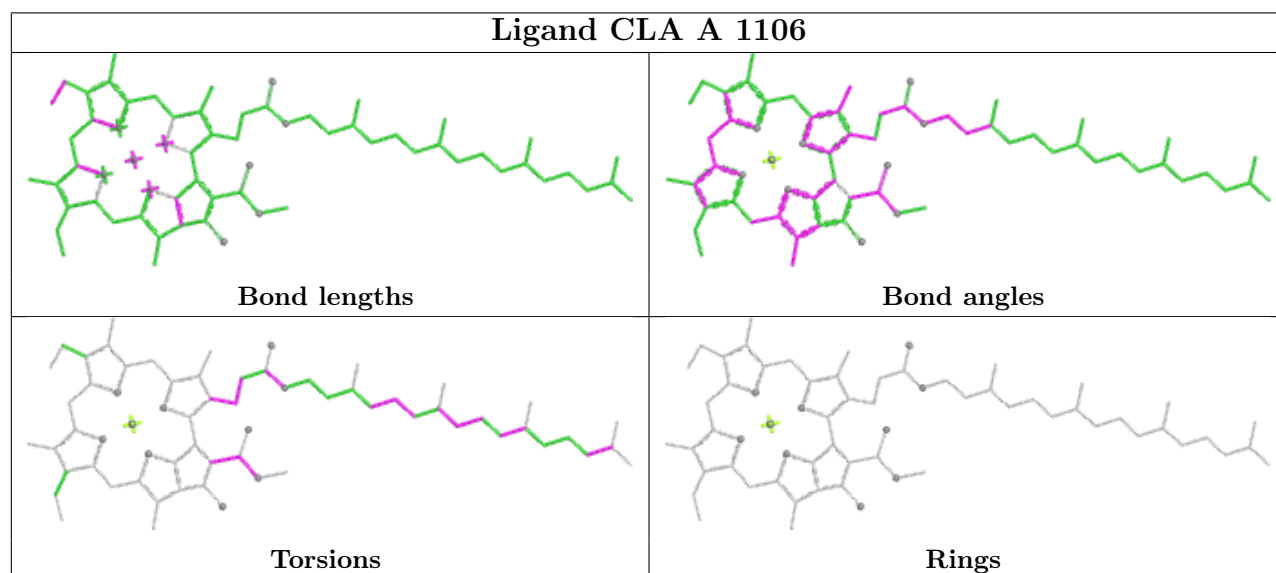
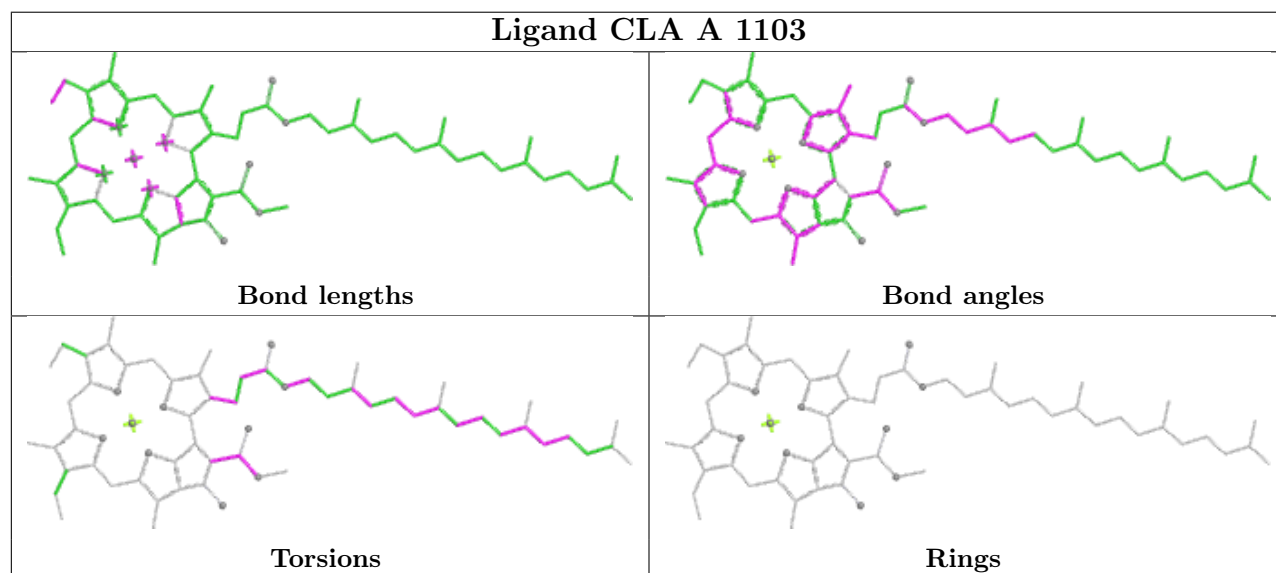
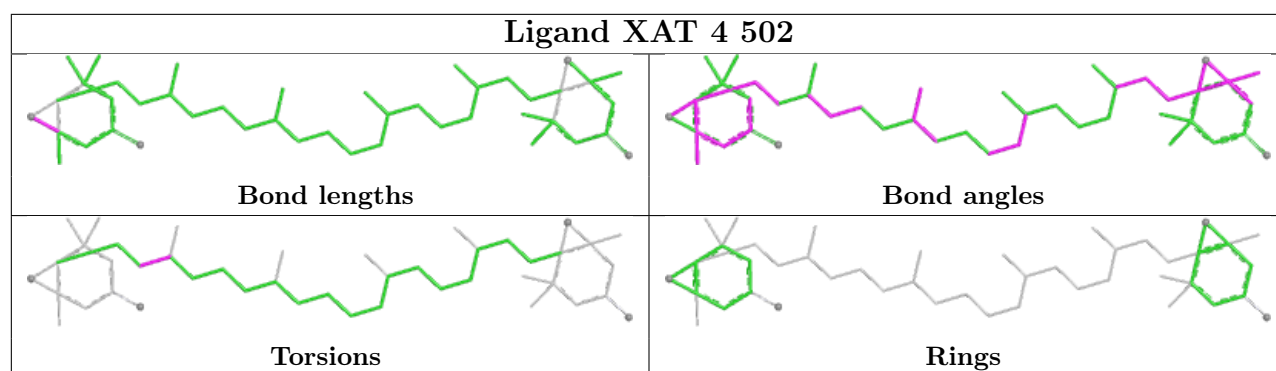
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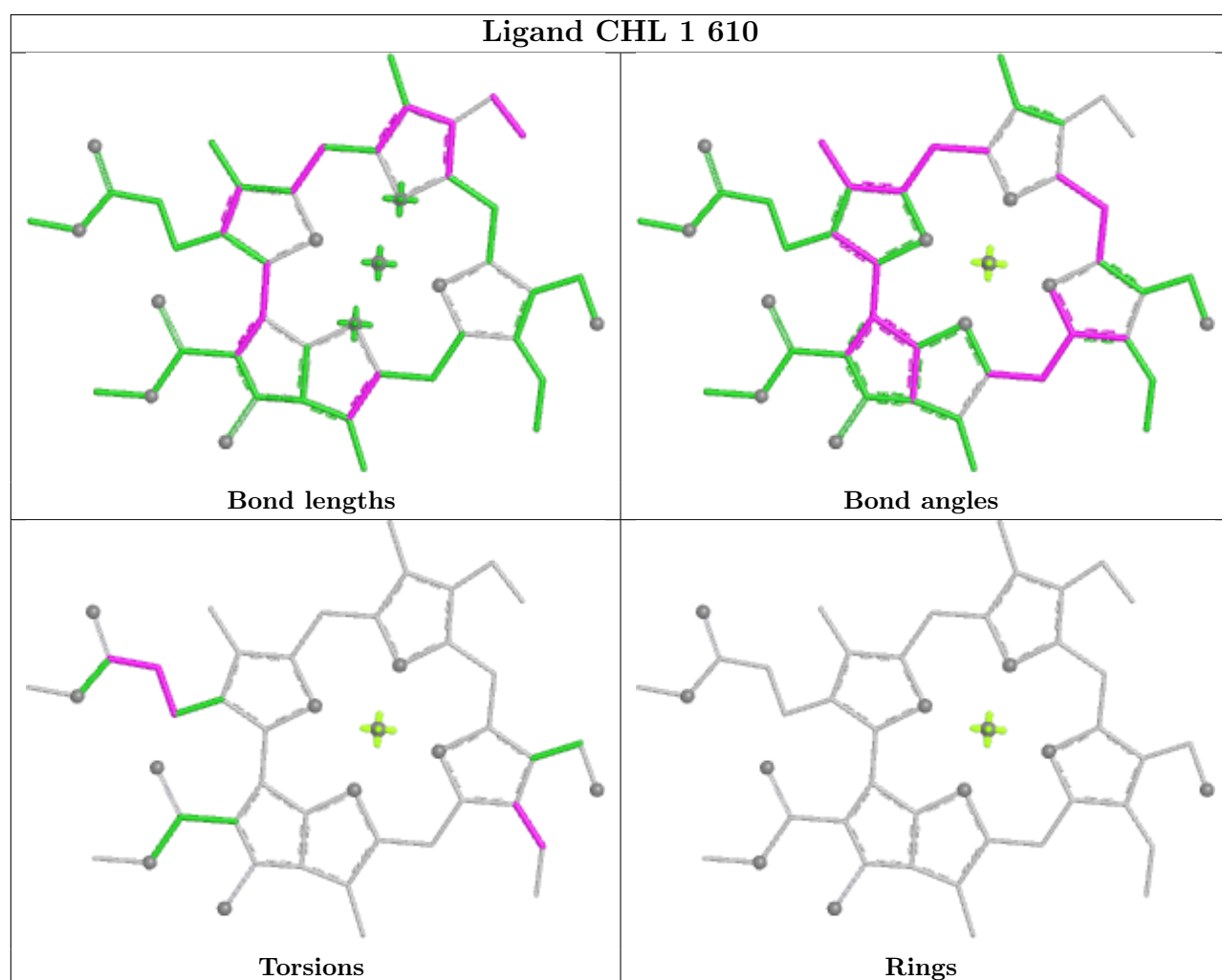
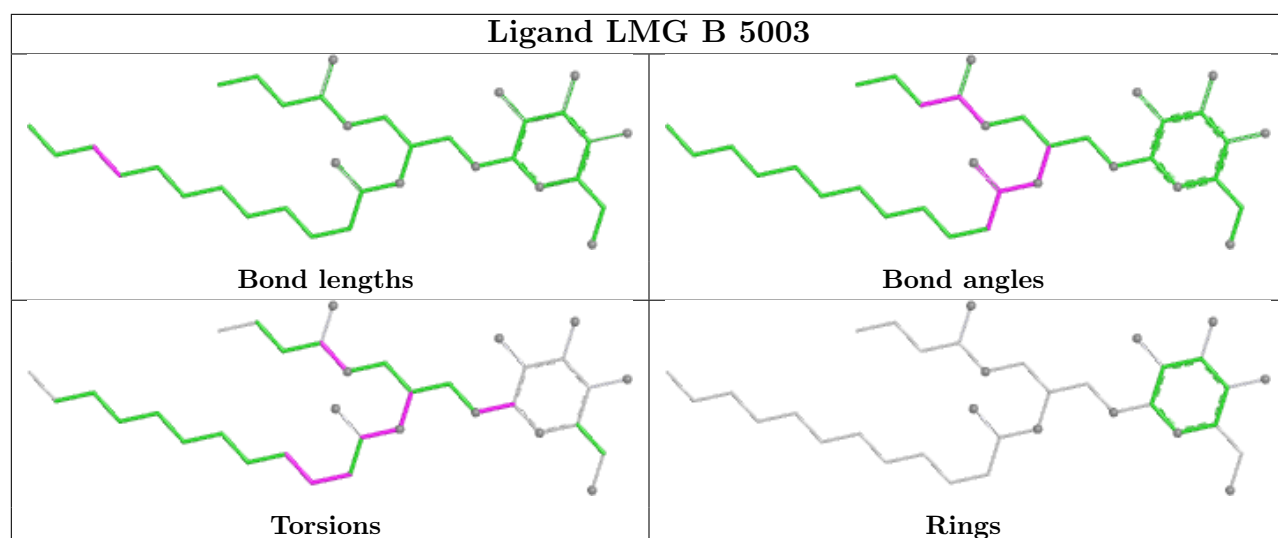
Mol	Chain	Res	Type	Clashes	Symm-Clashes
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29	J	4013	LUT	9	0
20	A	1102	CLA	5	0
20	A	1134	CLA	11	0
20	A	1135	CLA	4	0
20	B	1204	CLA	8	0
20	B	1219	CLA	2	0
20	H	1701	CLA	18	0
20	1	607	CLA	1	0
20	2	608	CLA	9	0
26	G	5002	LMG	8	0
20	2	605	CLA	9	0
26	G	5006	LMG	2	0
20	A	1138	CLA	4	0
20	3	603	CLA	9	0
20	B	1203	CLA	4	0
20	4	602	CLA	17	0
20	A	1114	CLA	2	0

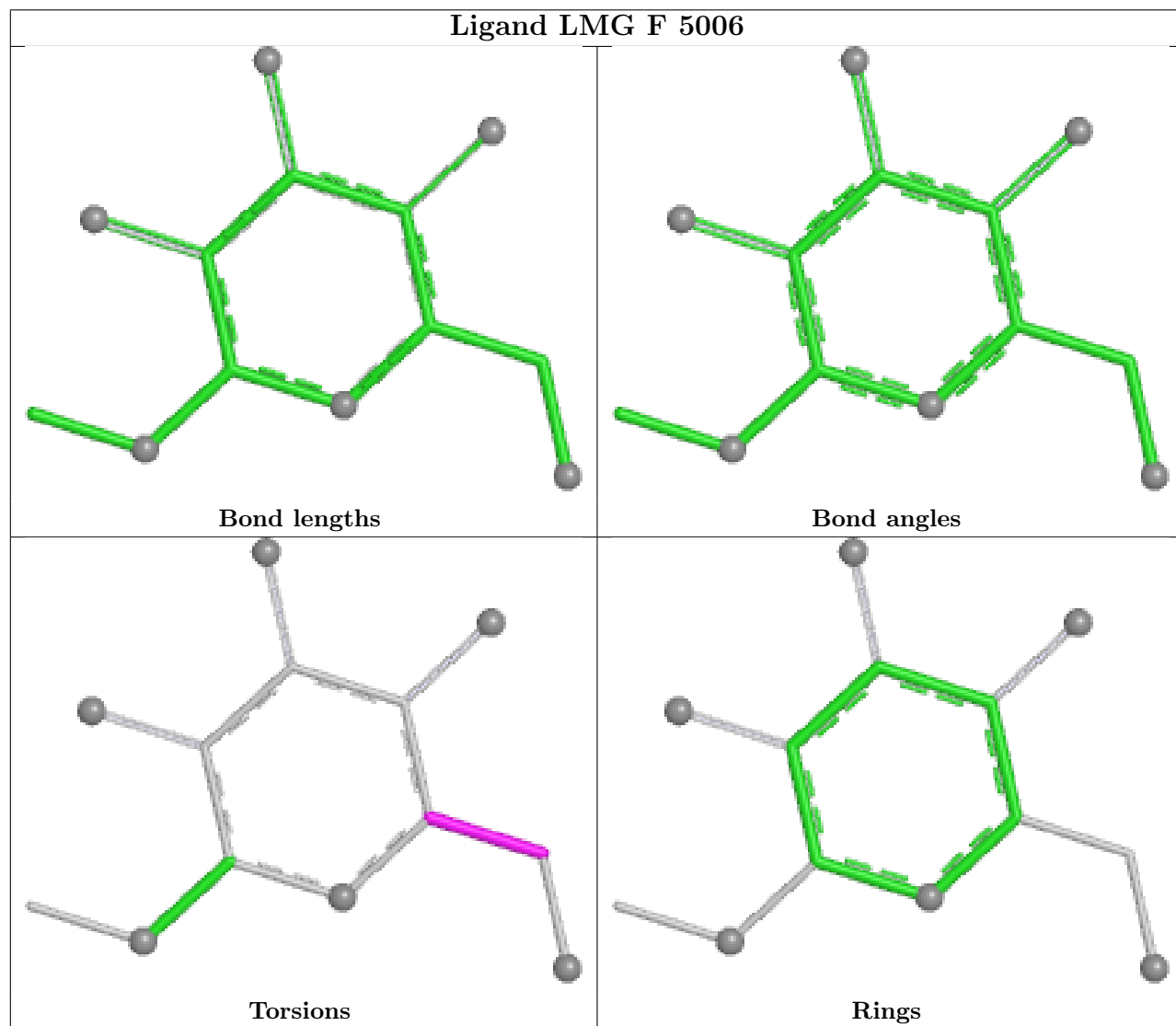
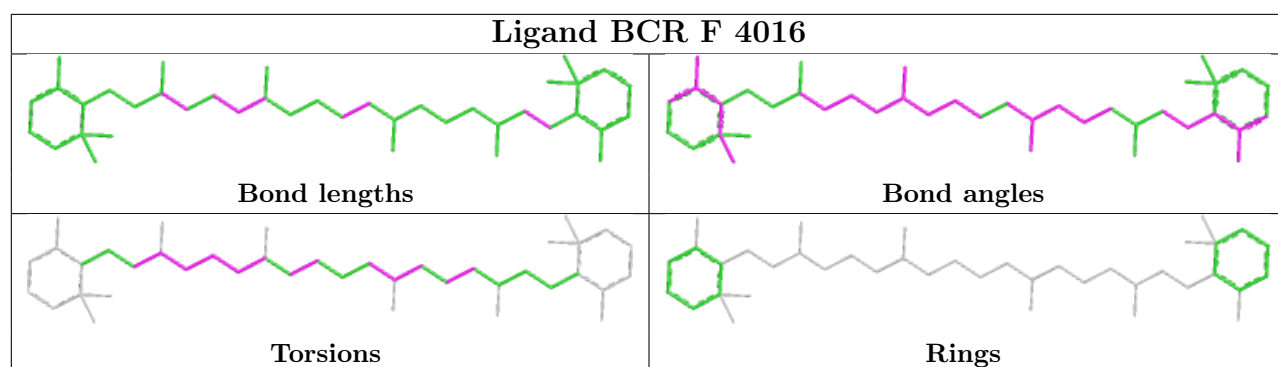
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

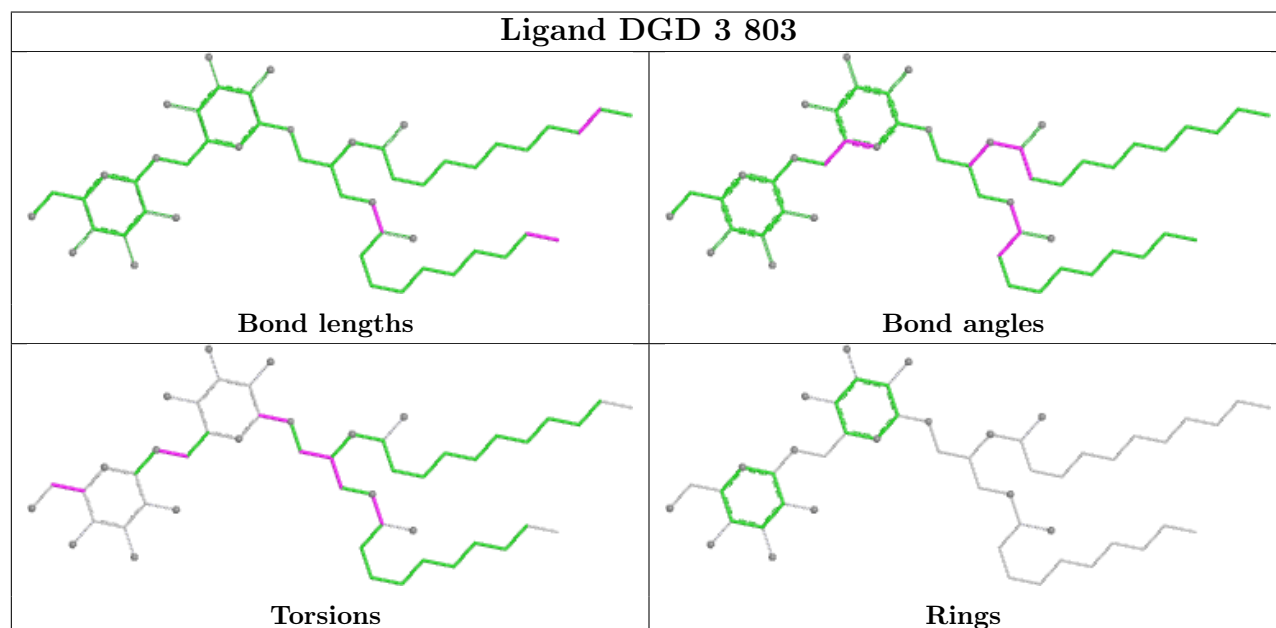
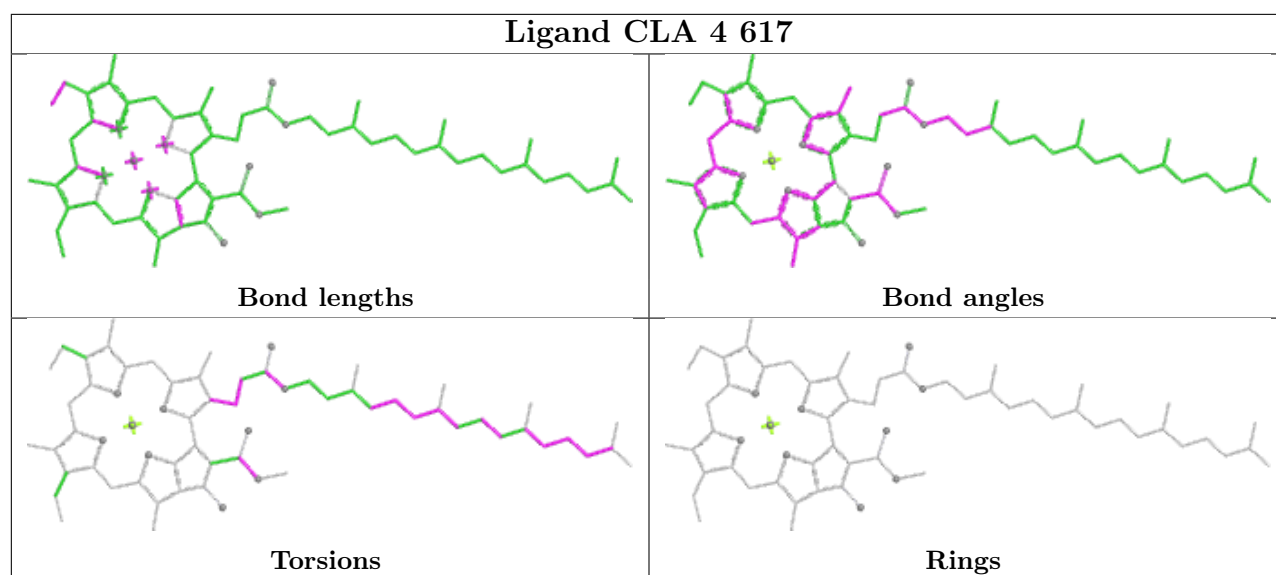


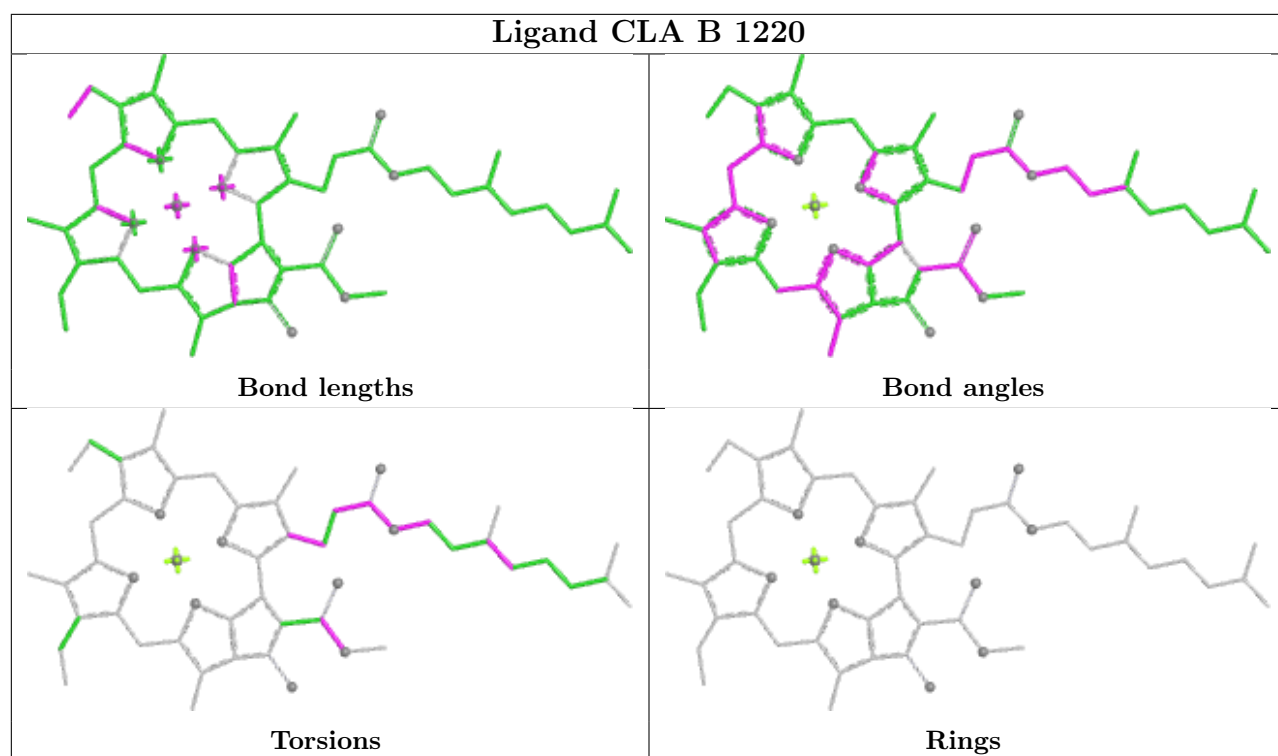




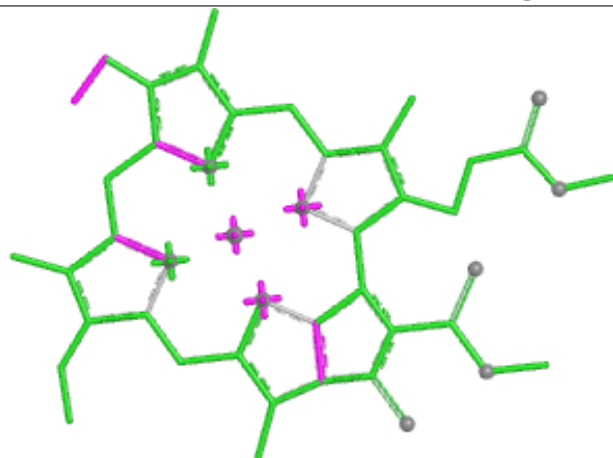




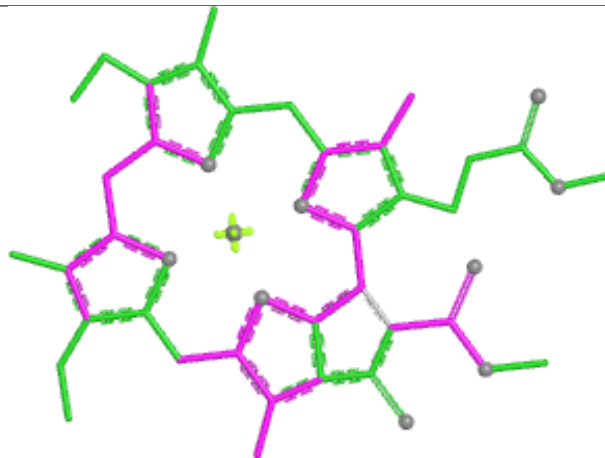




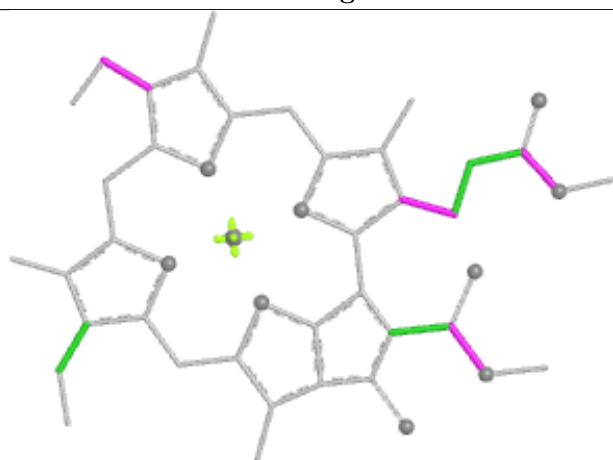
Ligand CLA 1 602



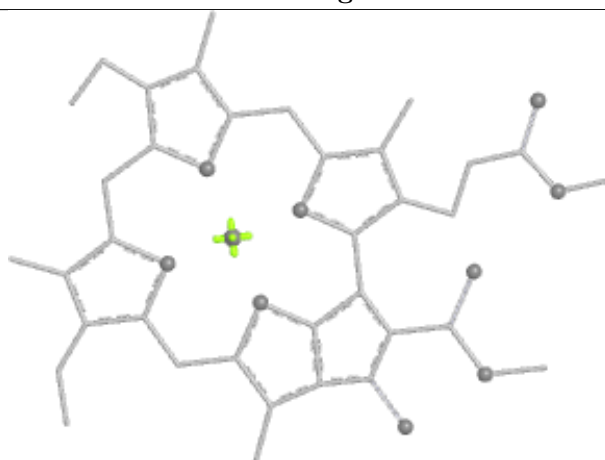
Bond lengths



Bond angles

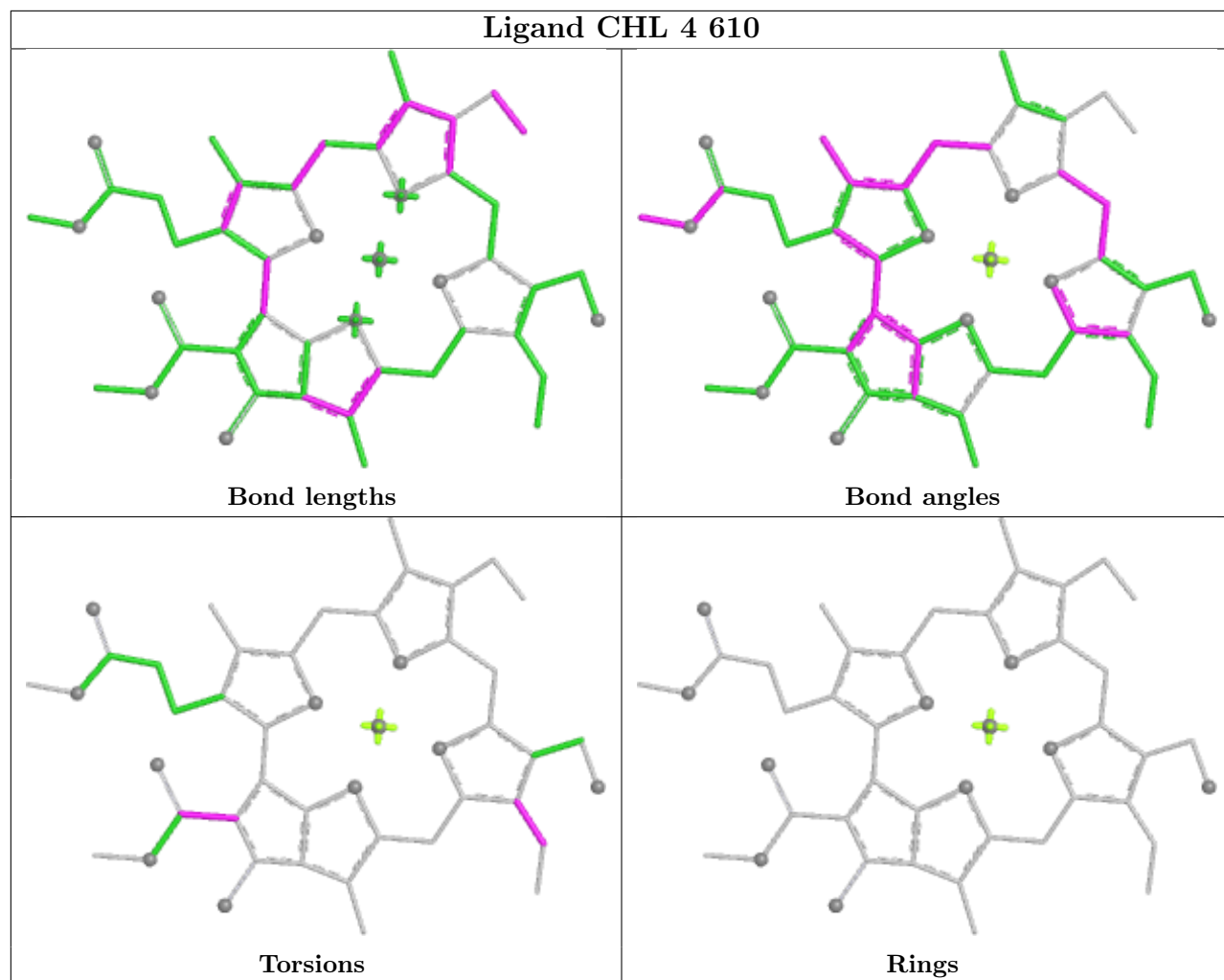


Torsions

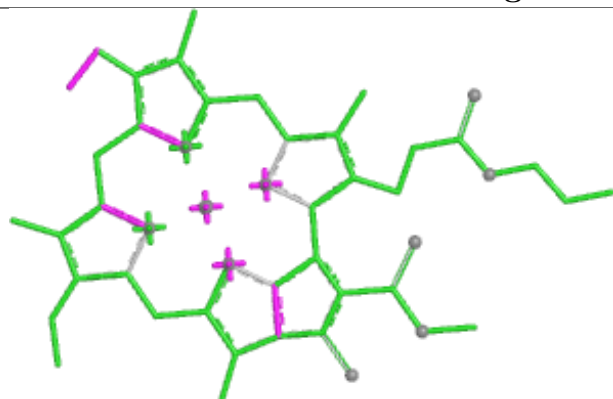


Rings

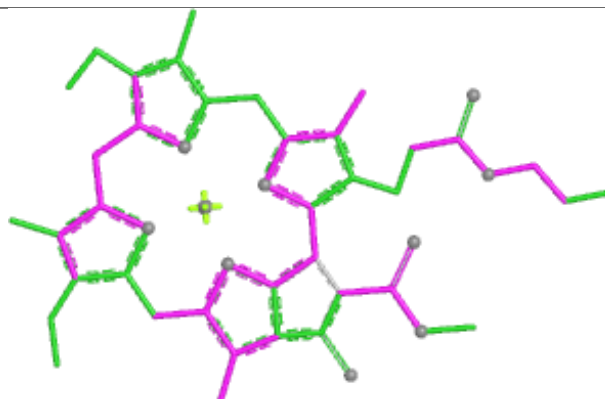
Ligand CHL 4 610



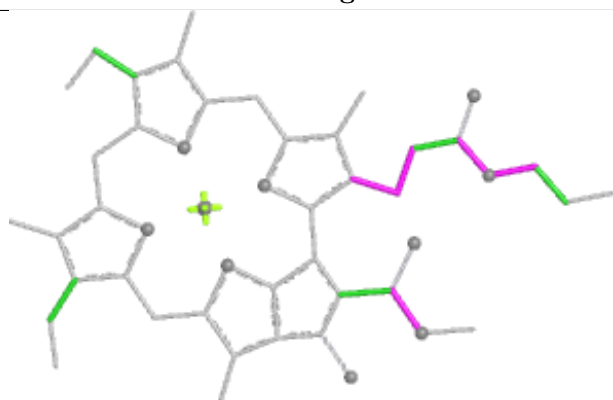
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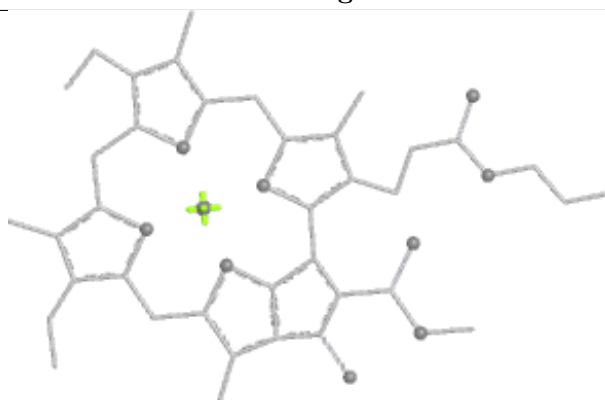
Bond lengths



Bond angles

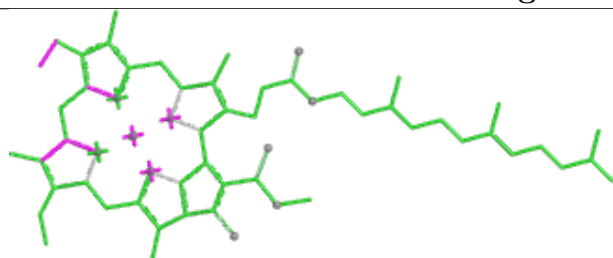


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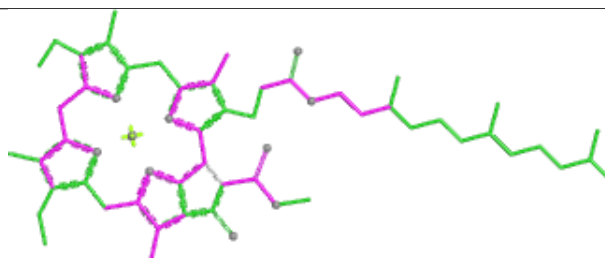


Rings

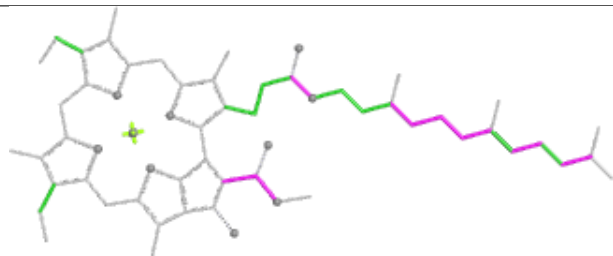
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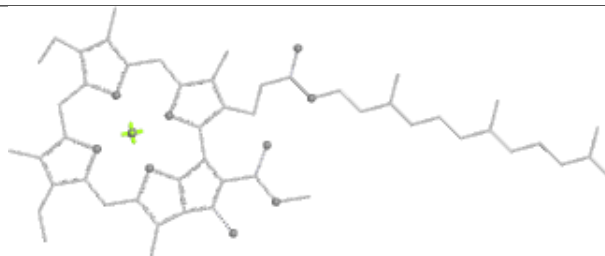
Bond lengths



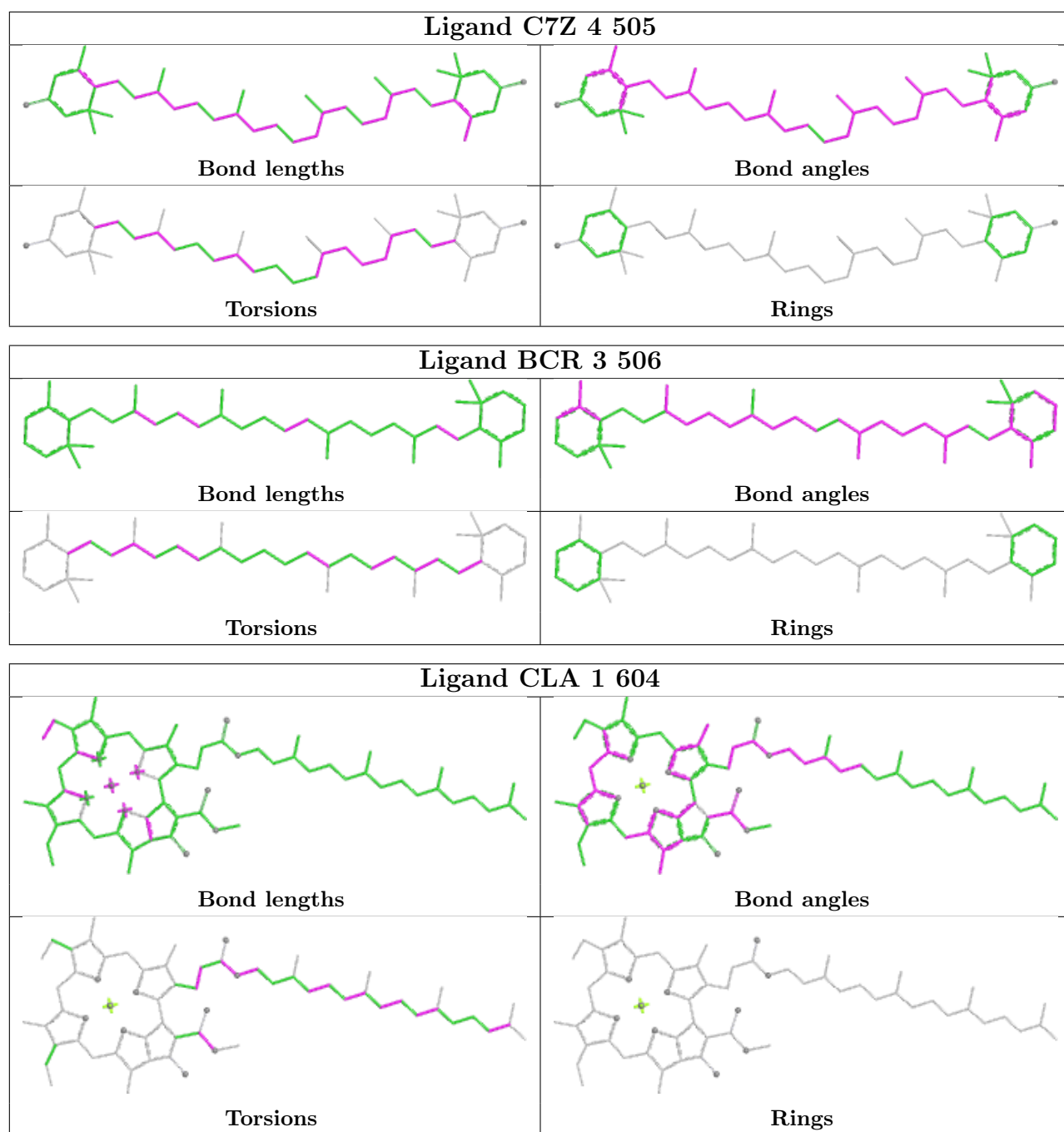
Bond angles

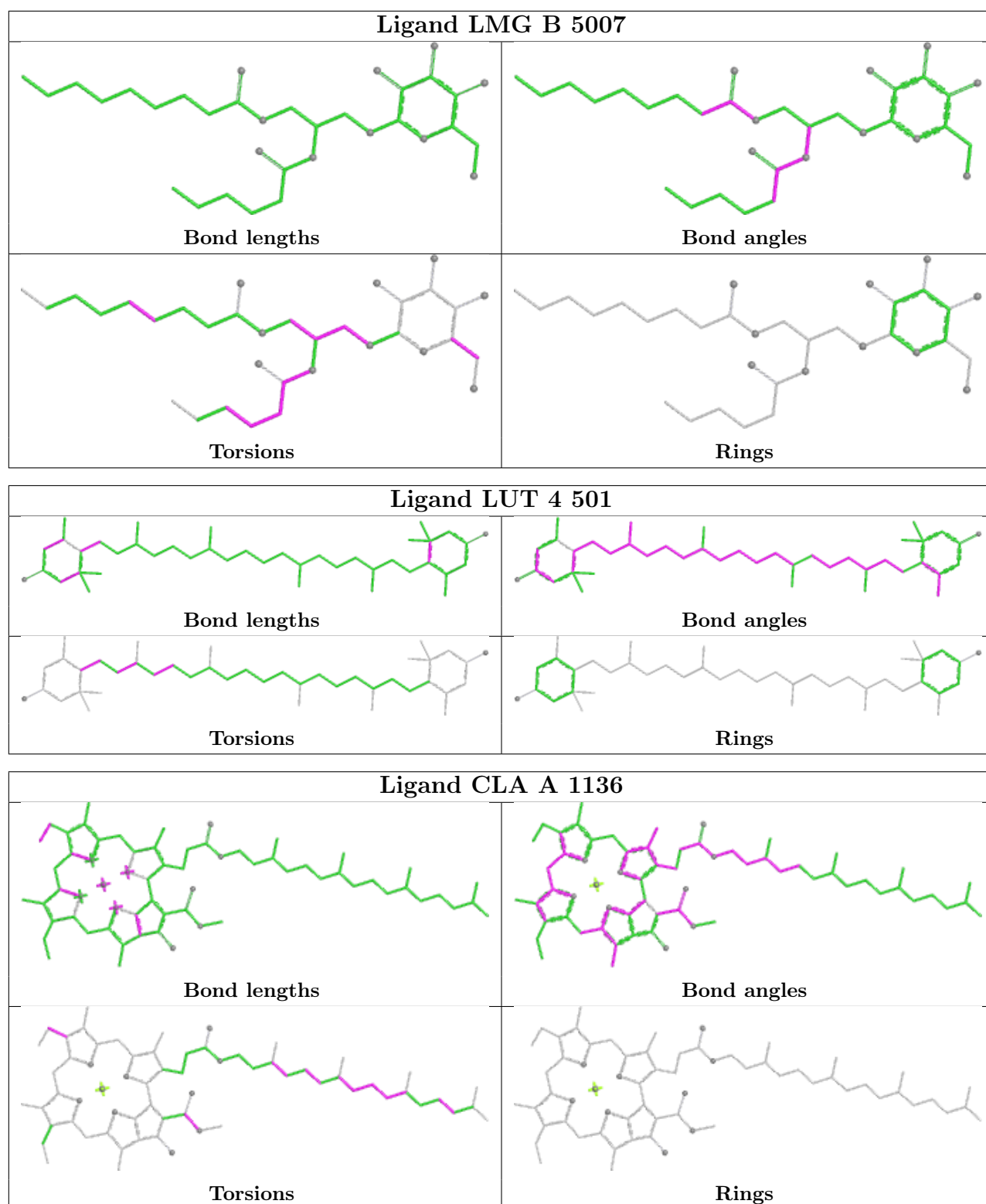


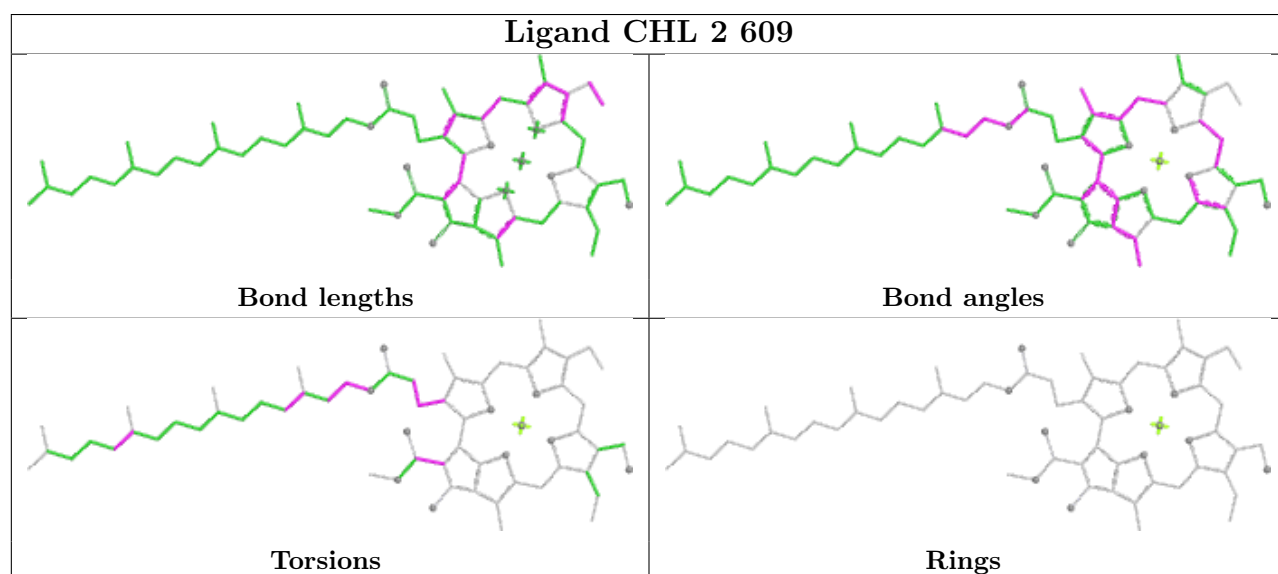
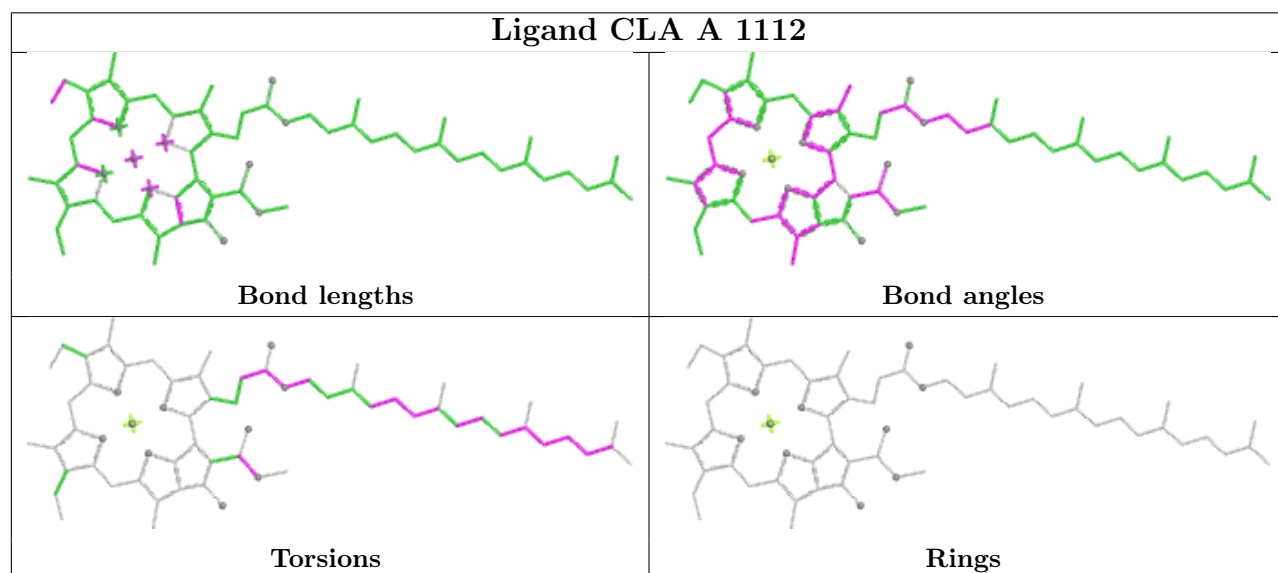
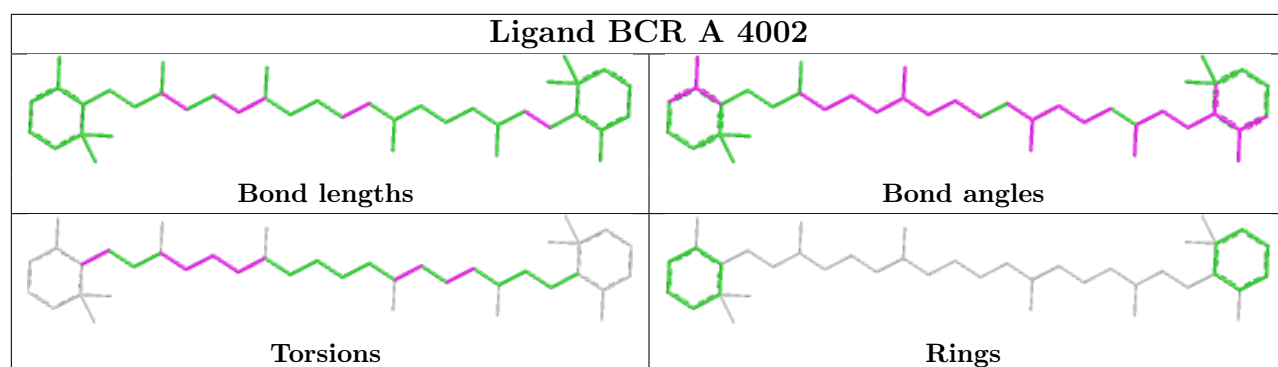
Torsions

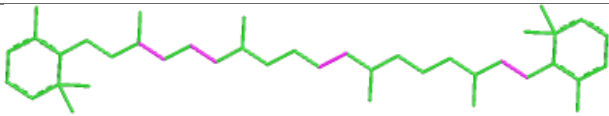
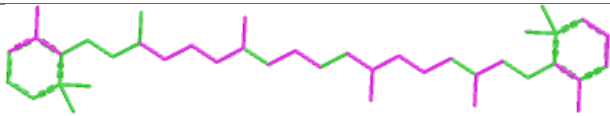
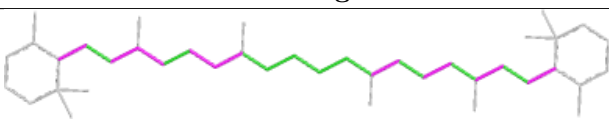
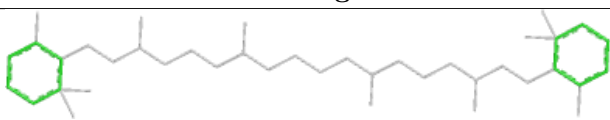


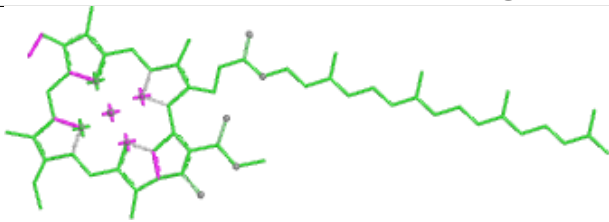
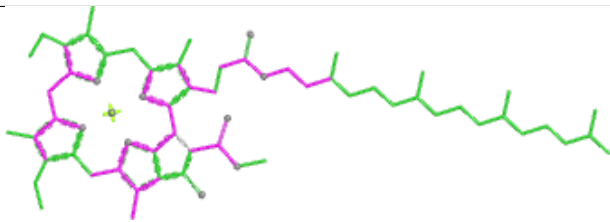
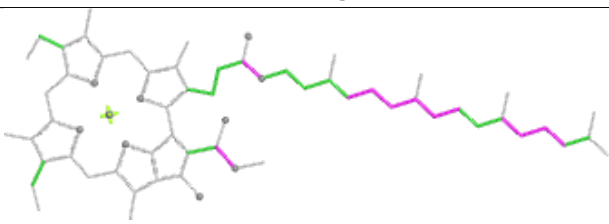
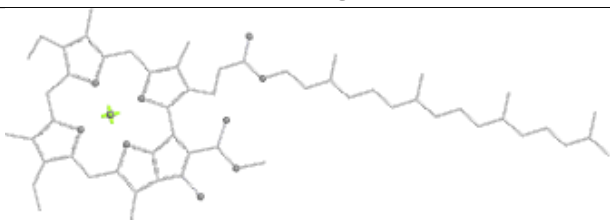
Rings

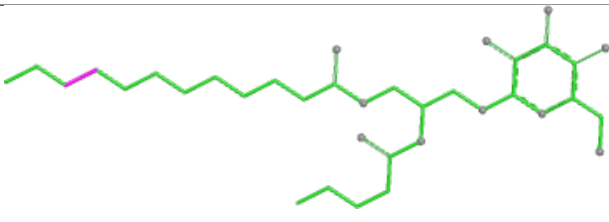
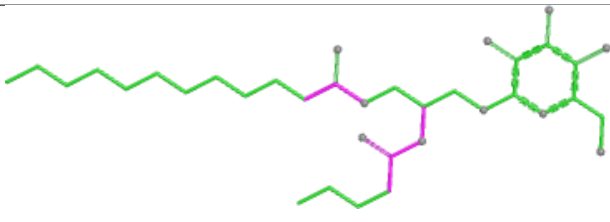
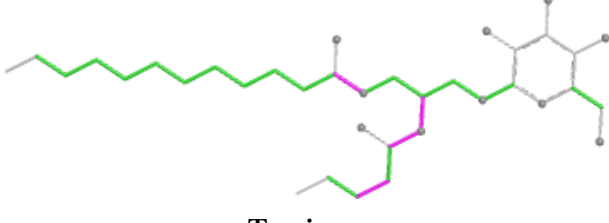
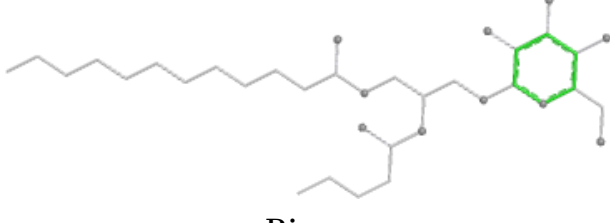


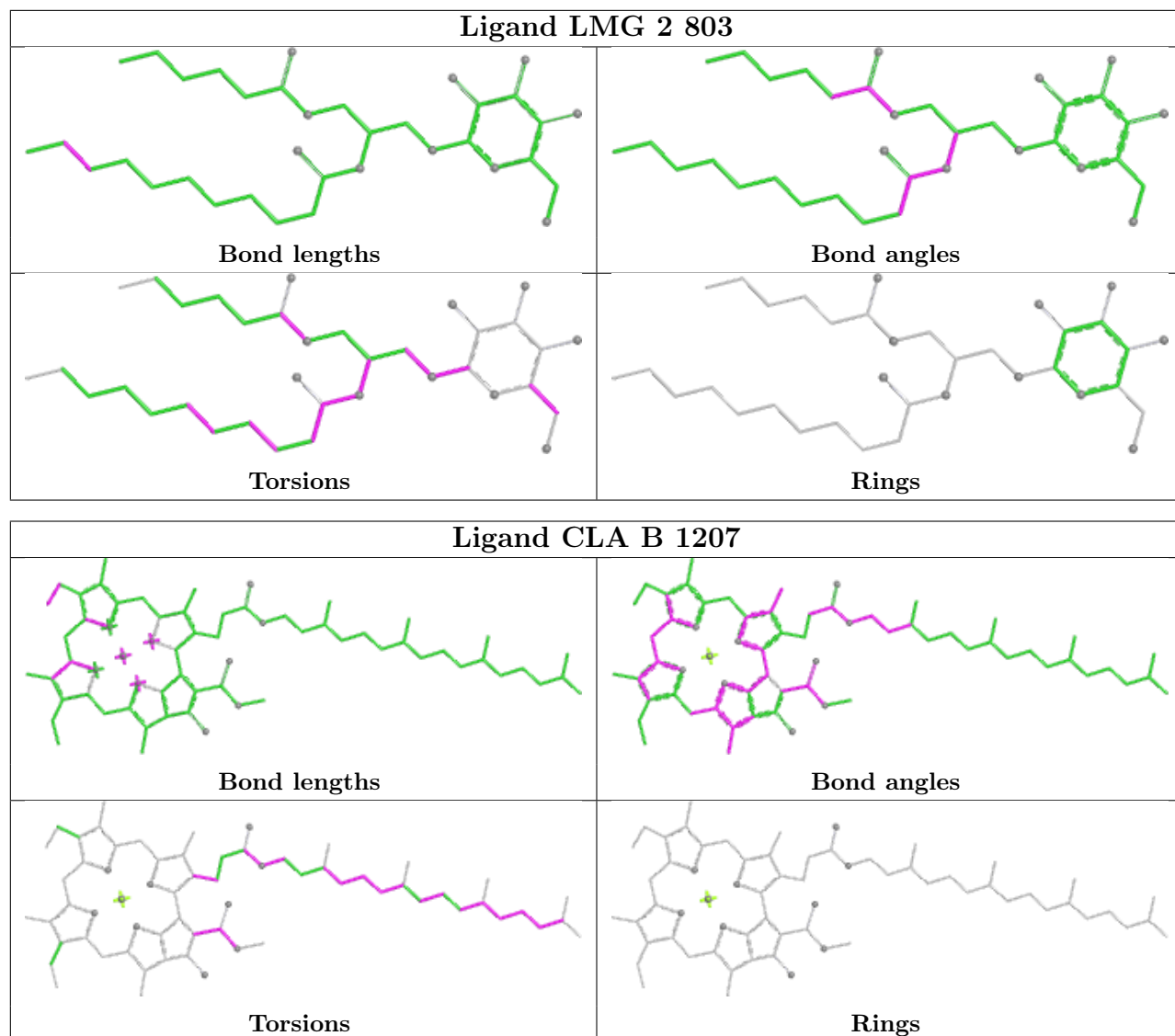




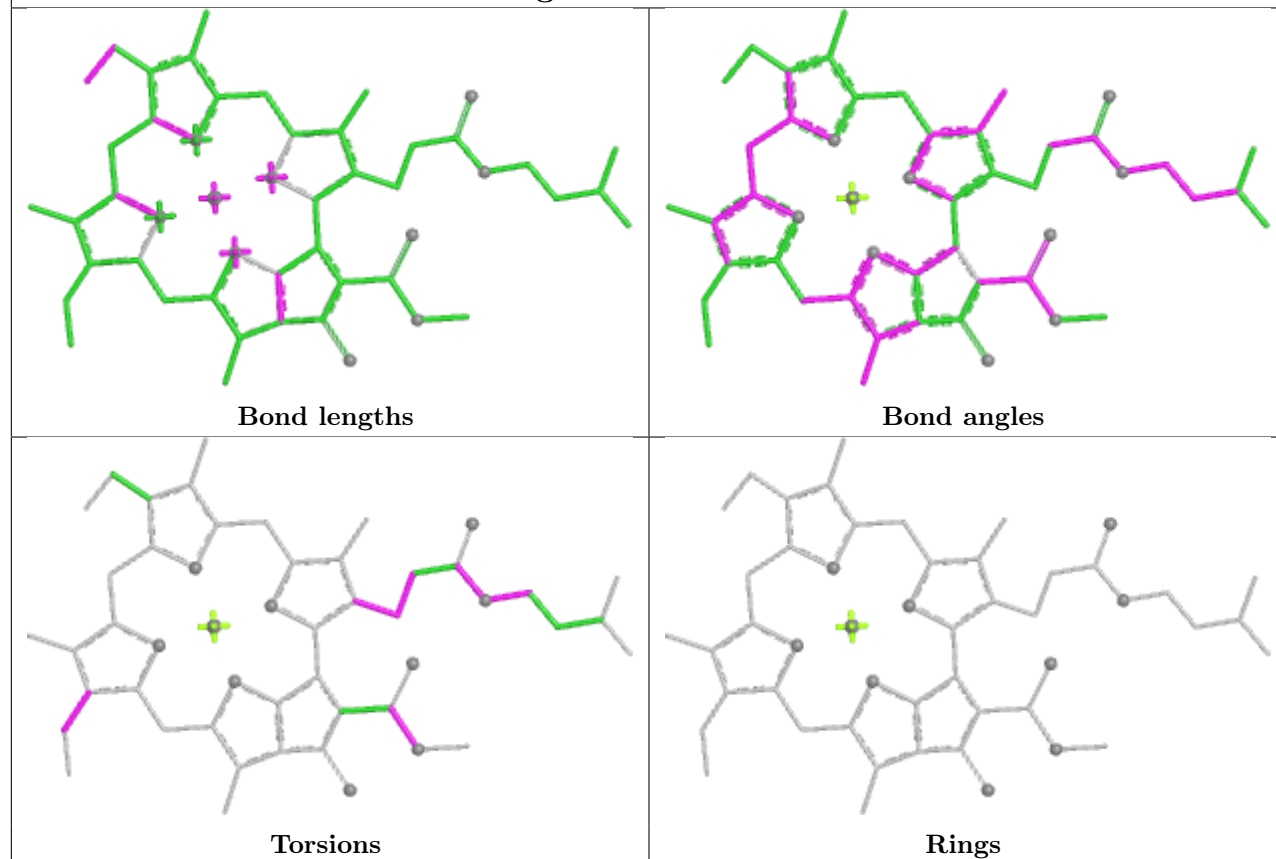
Ligand BCR 1 503	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA A 1129	
	
Bond lengths	Bond angles
	
Torsions	Rings

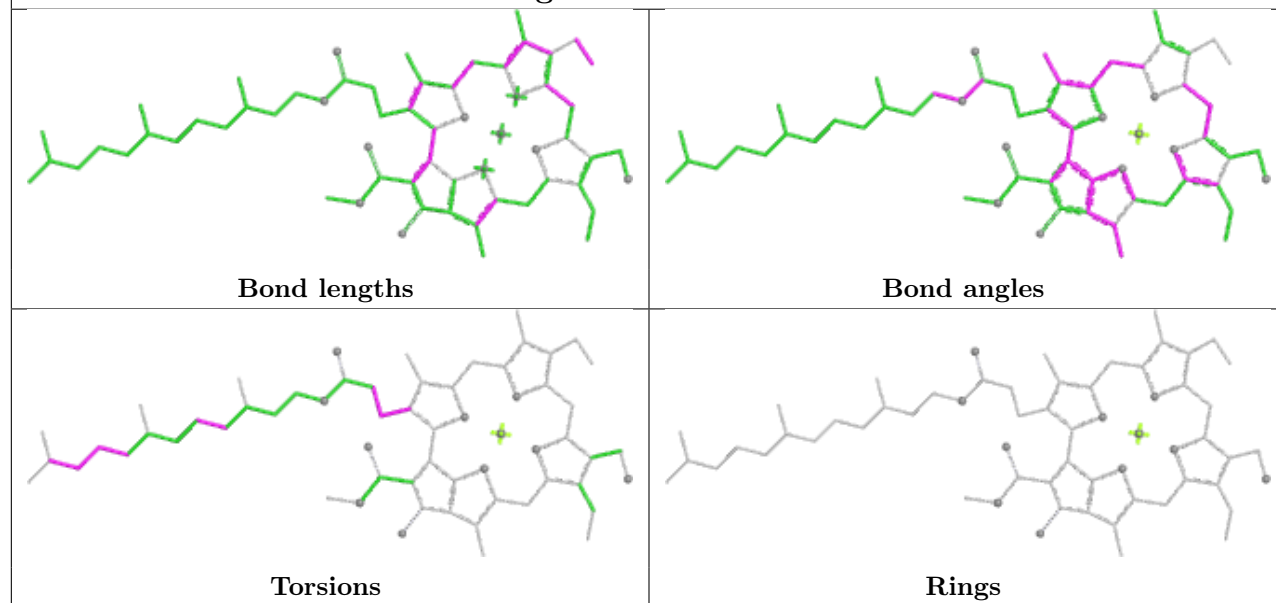
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Bond lengths	Bond angles
	
Torsions	Rings

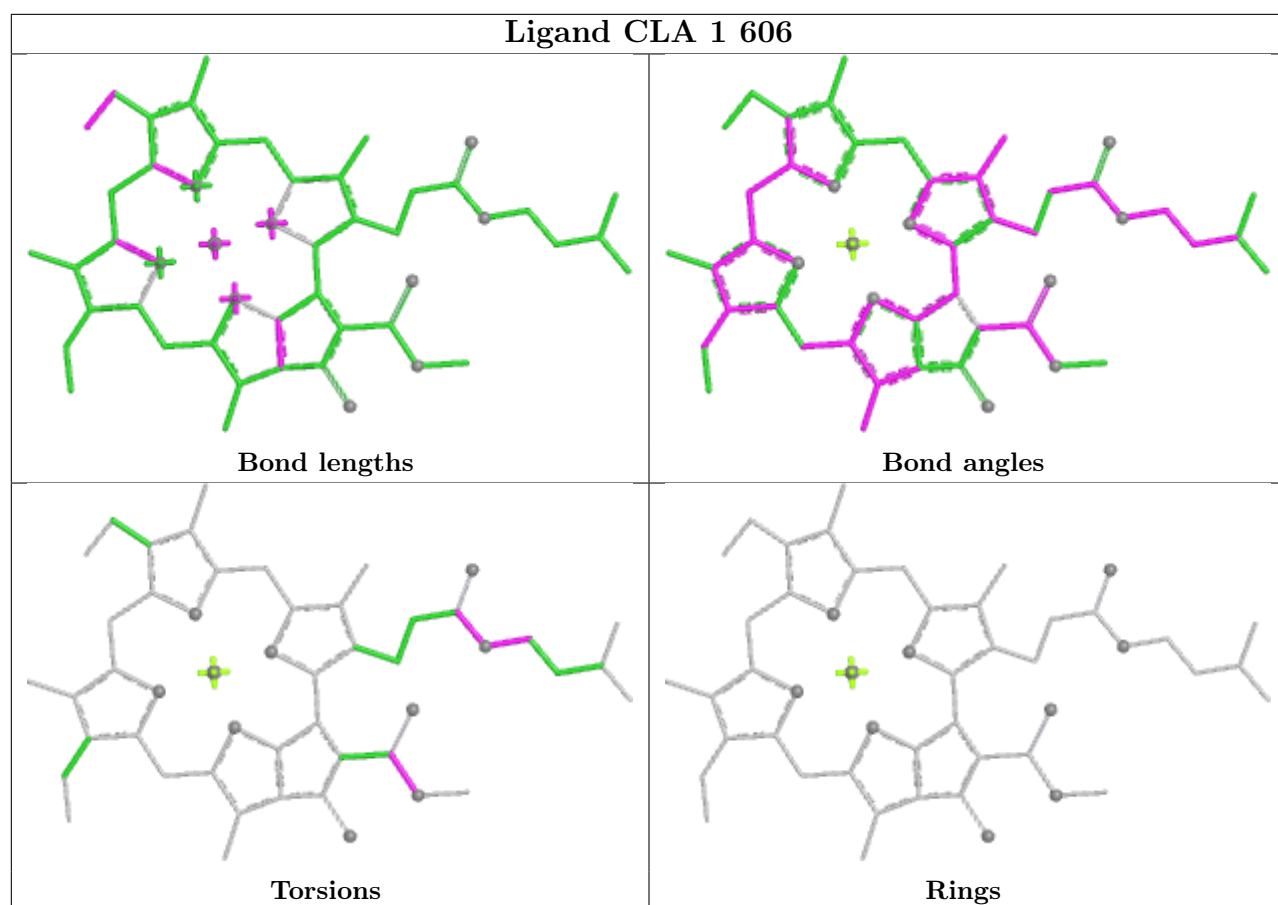


Ligand CLA 3 612

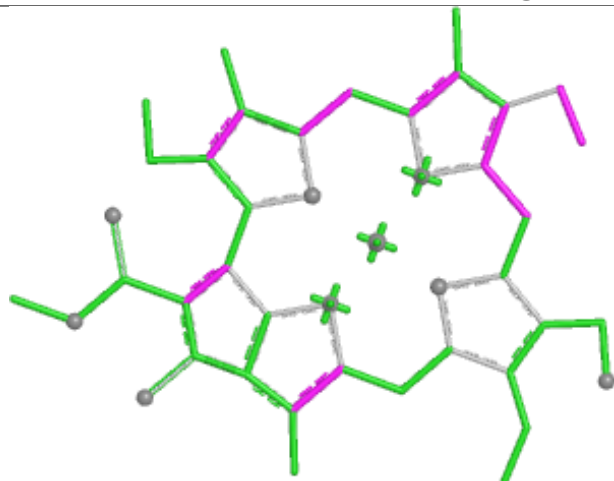


Ligand CHL 4 613

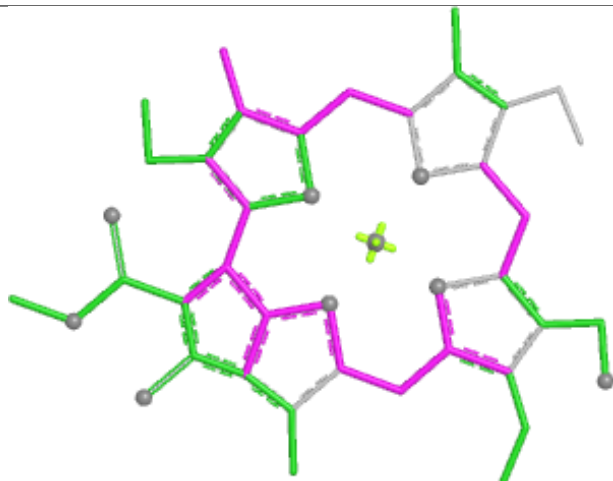




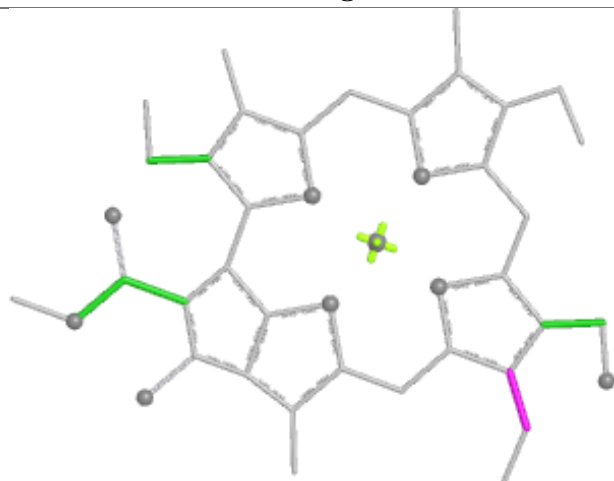
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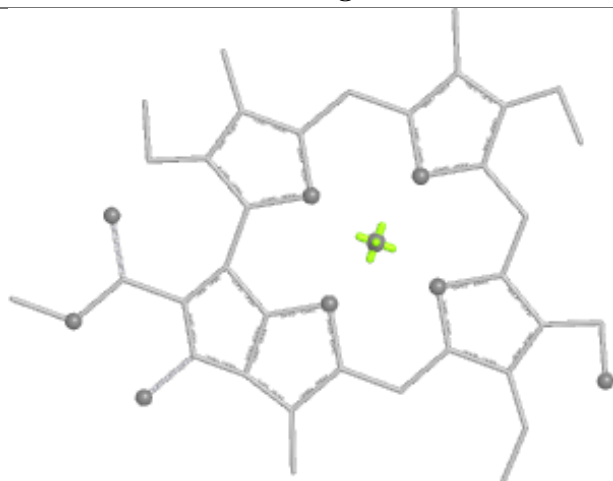
Bond lengths



Bond angles

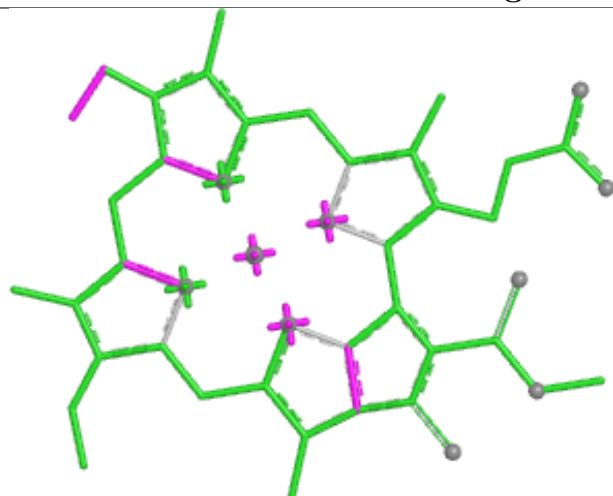


Torsions

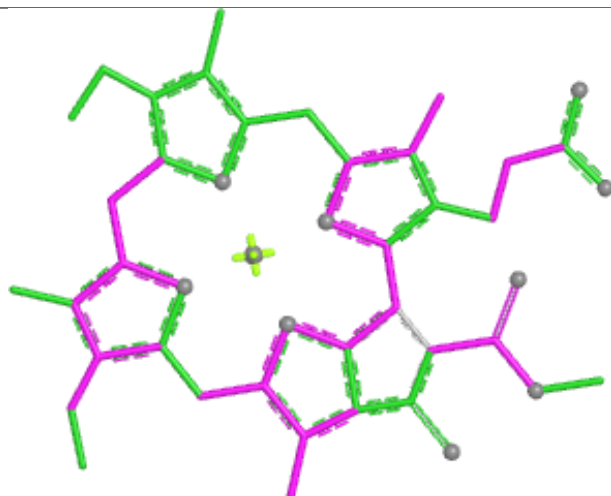


Rings

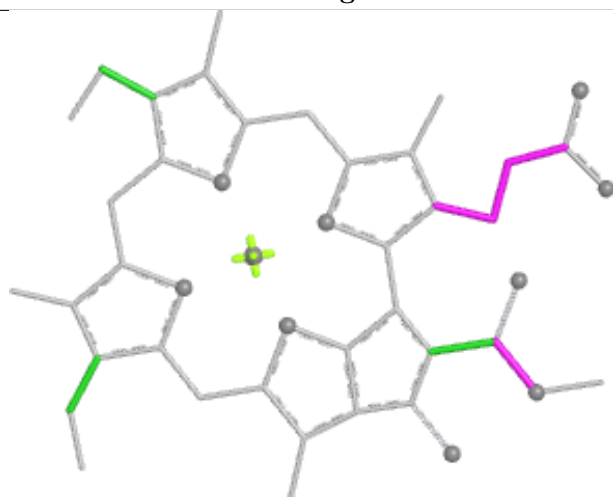
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Bond lengths



Bond angles

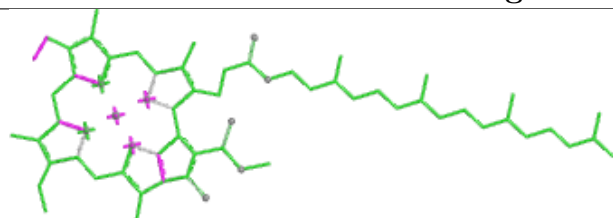


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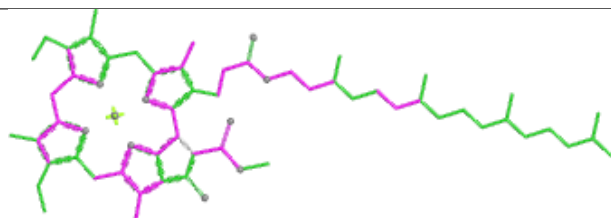


Rings

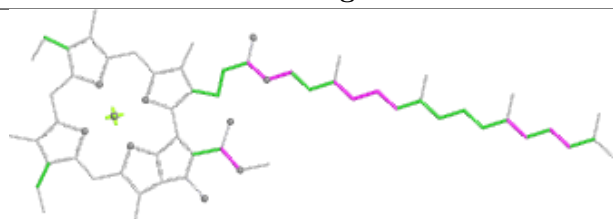
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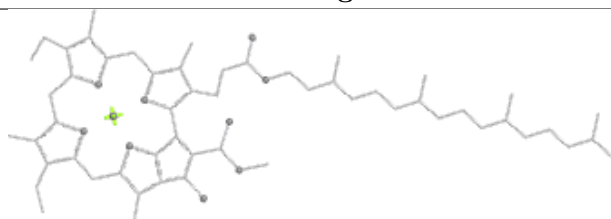
Bond lengths



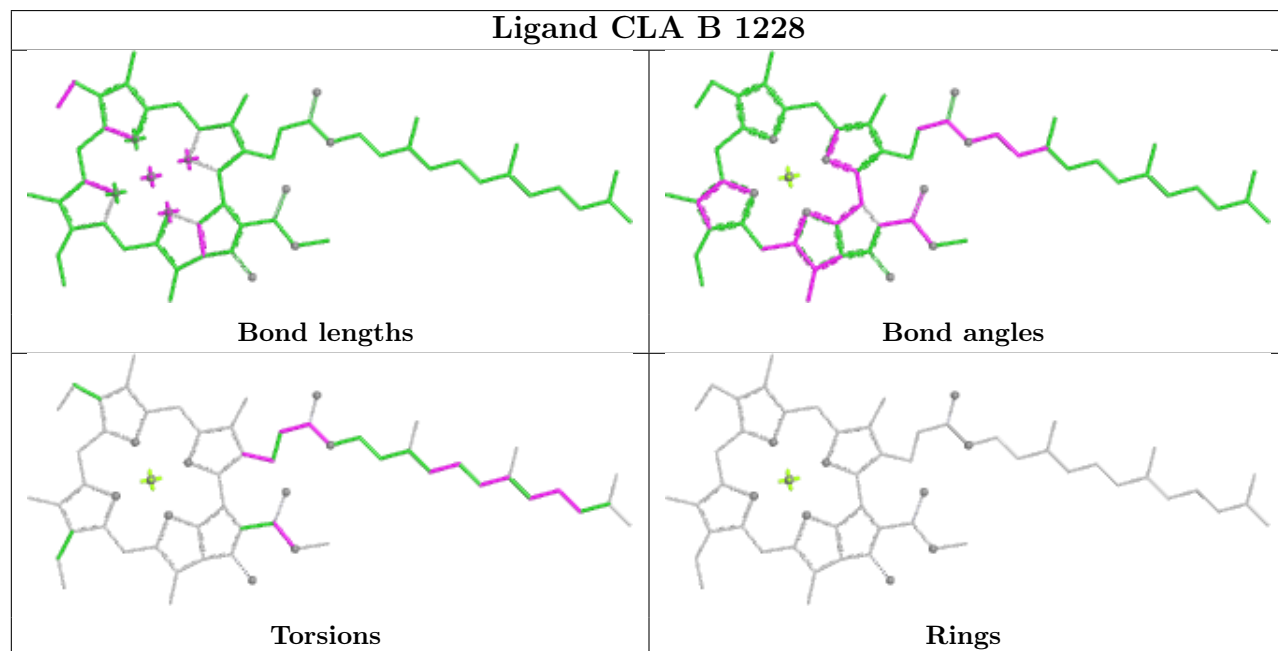
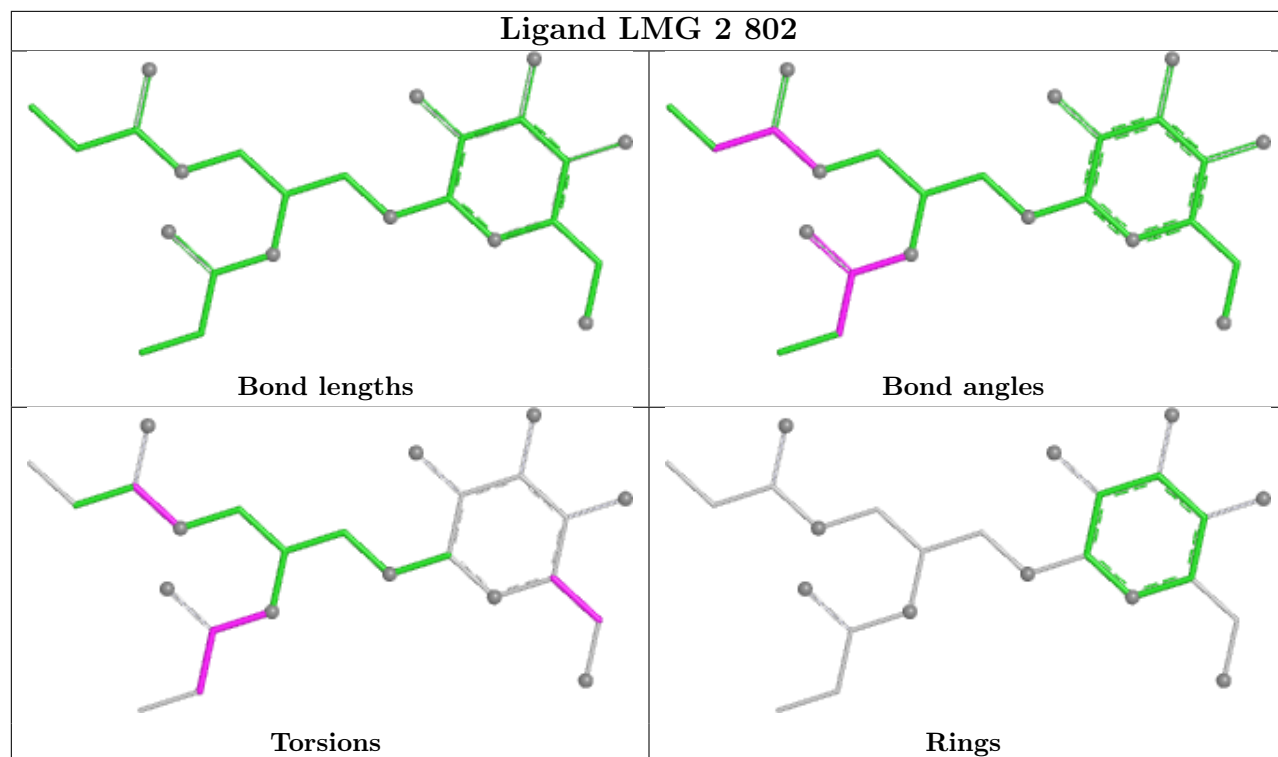
Bond angles

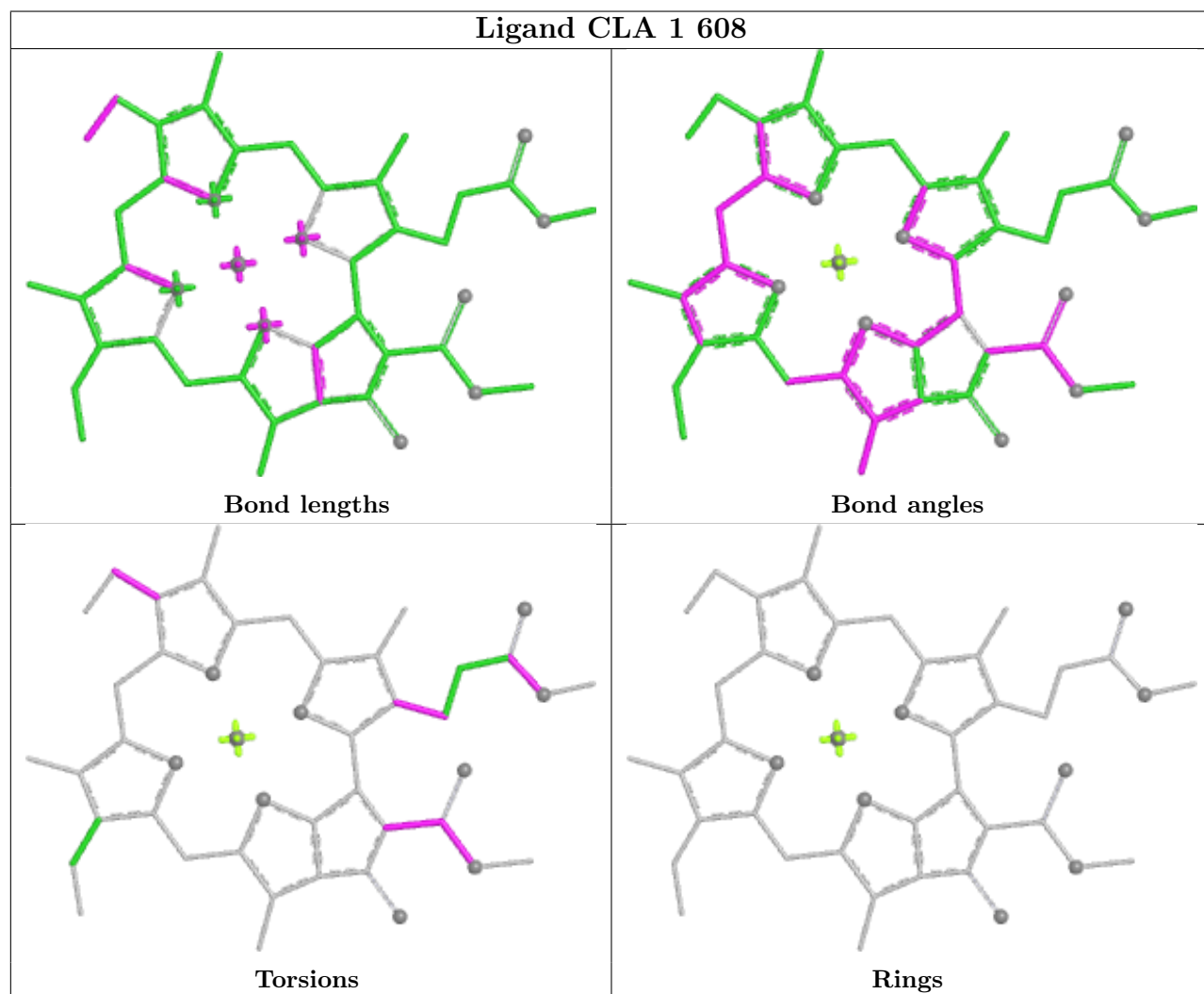
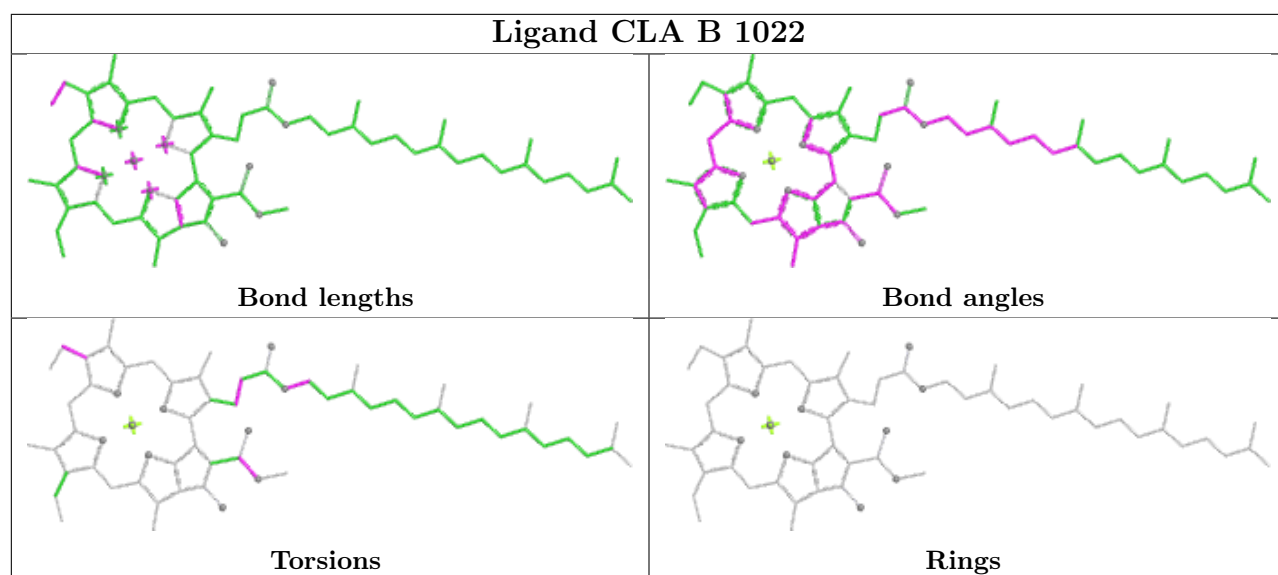


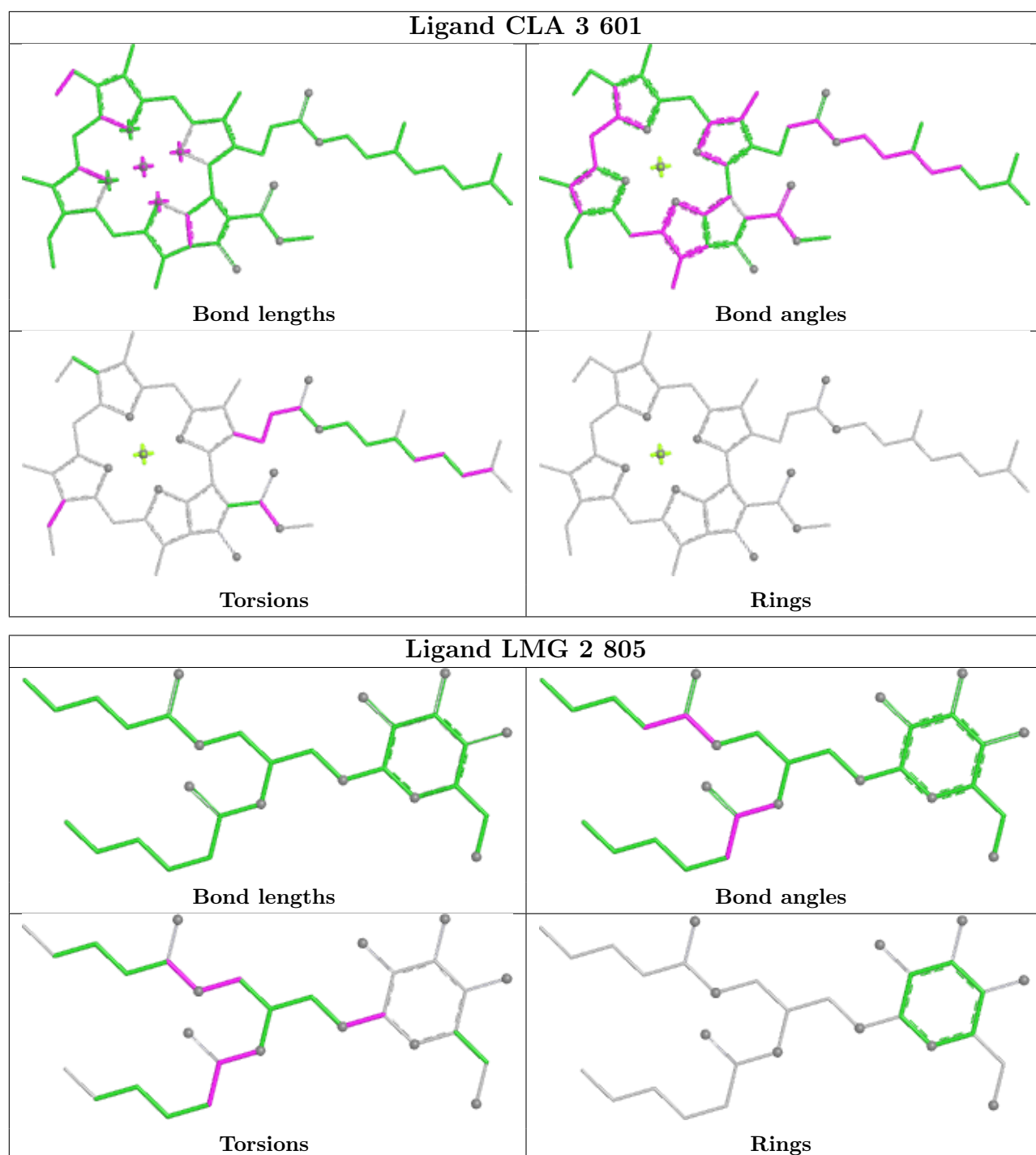
Torsions

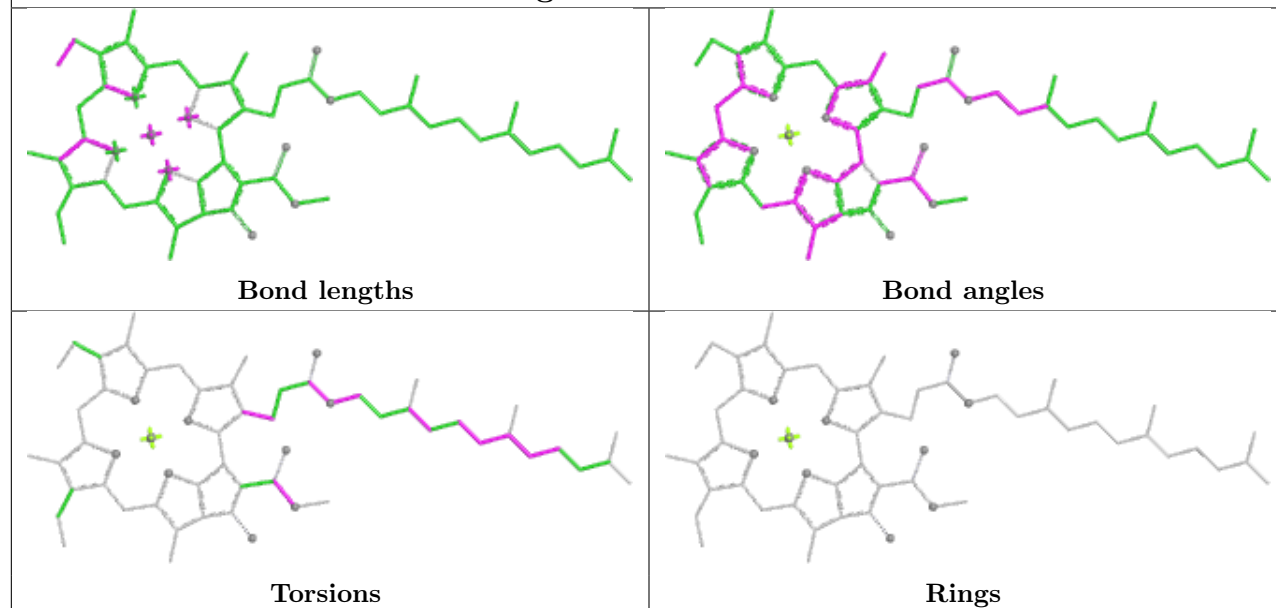
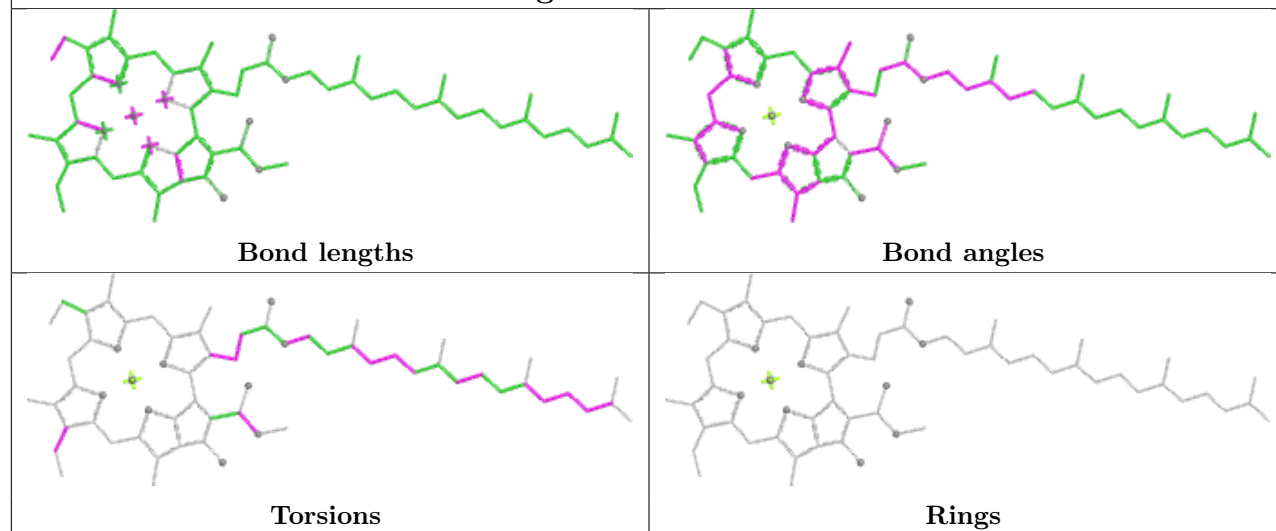


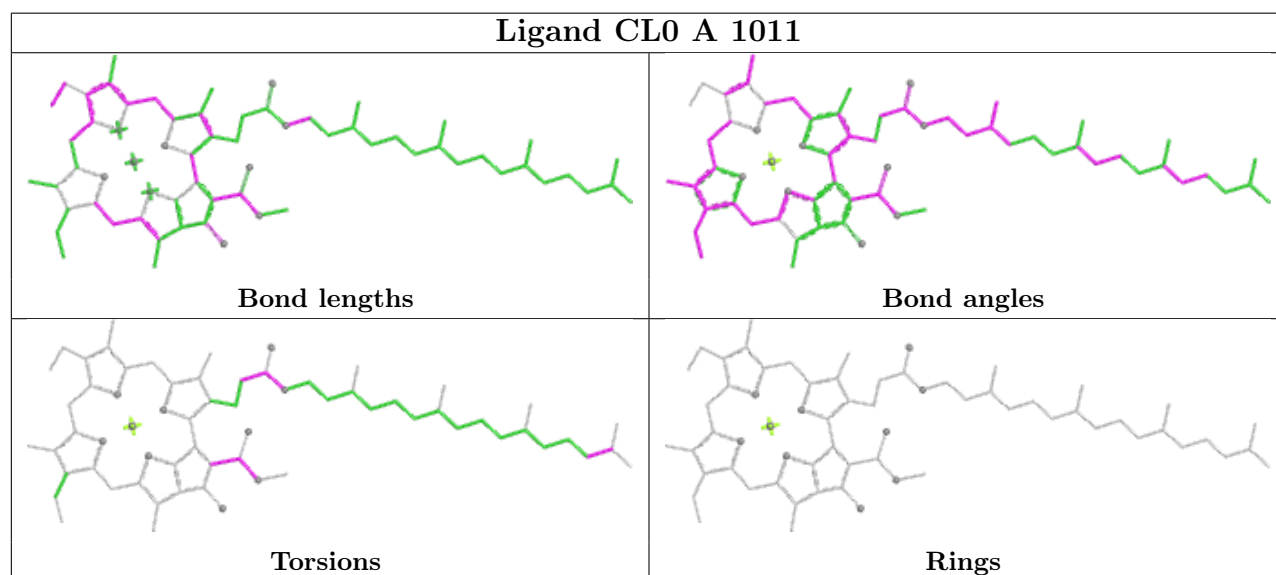
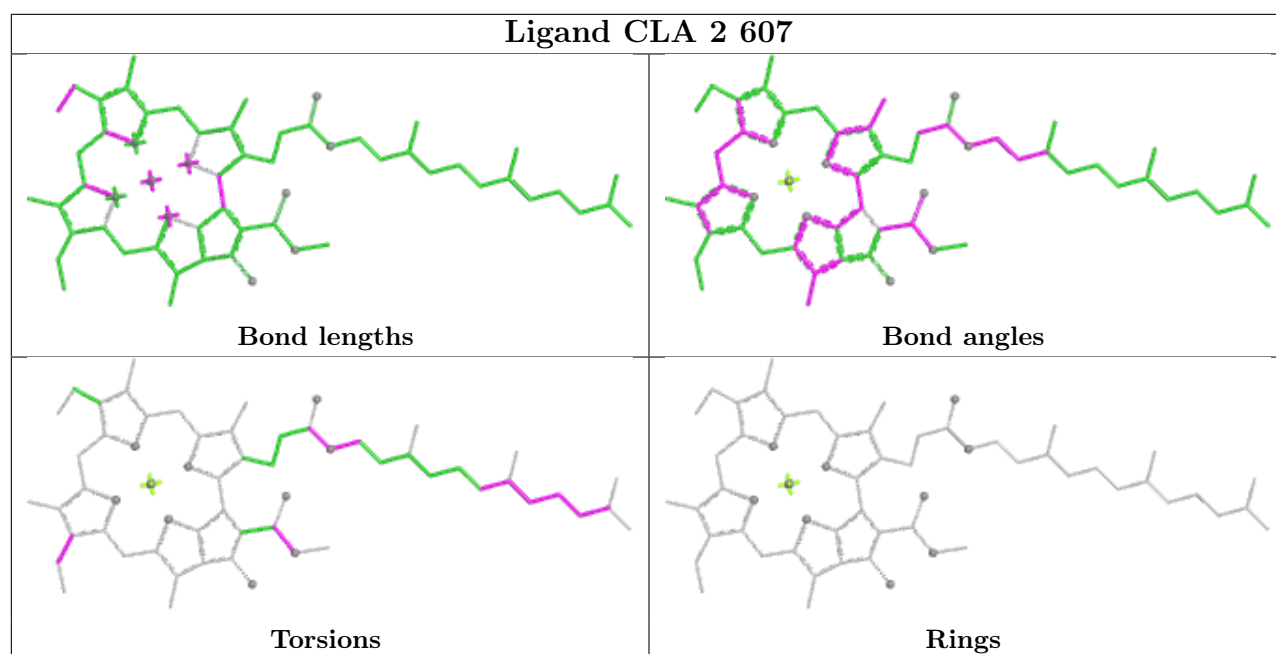
Rings

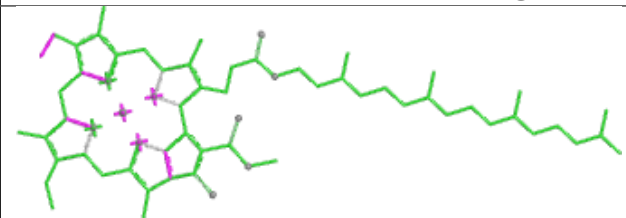
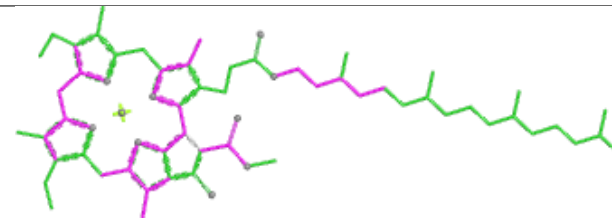
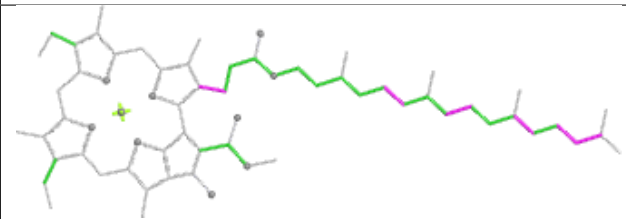
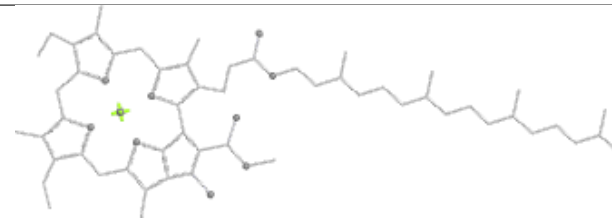


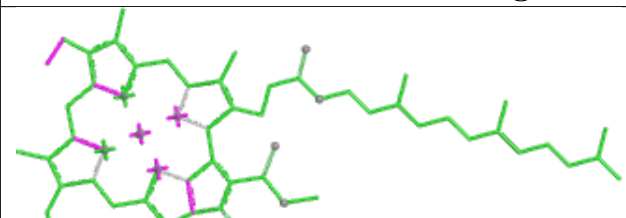
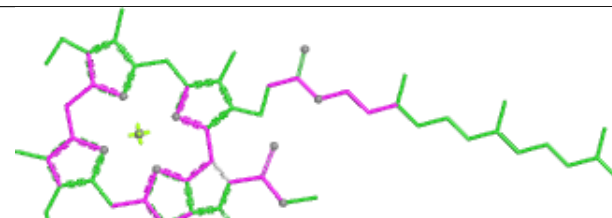
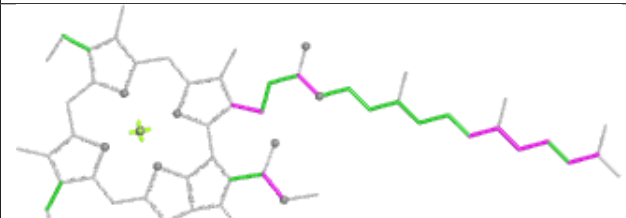
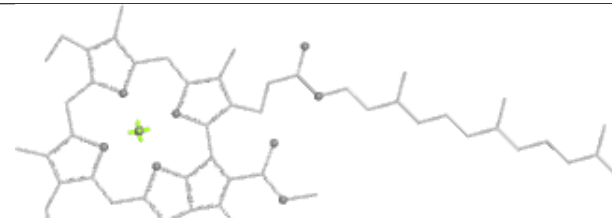


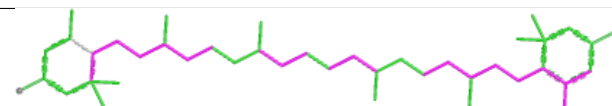
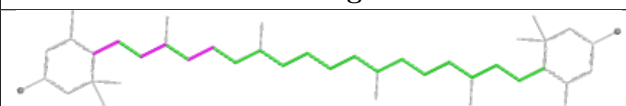
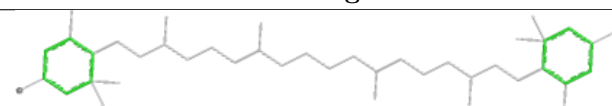


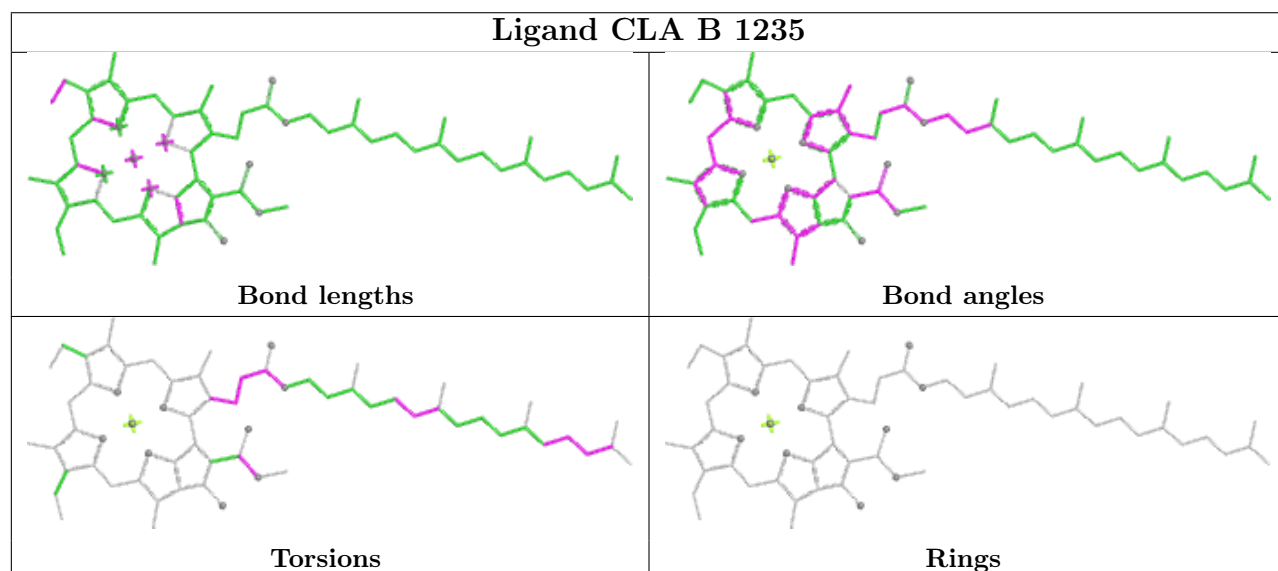
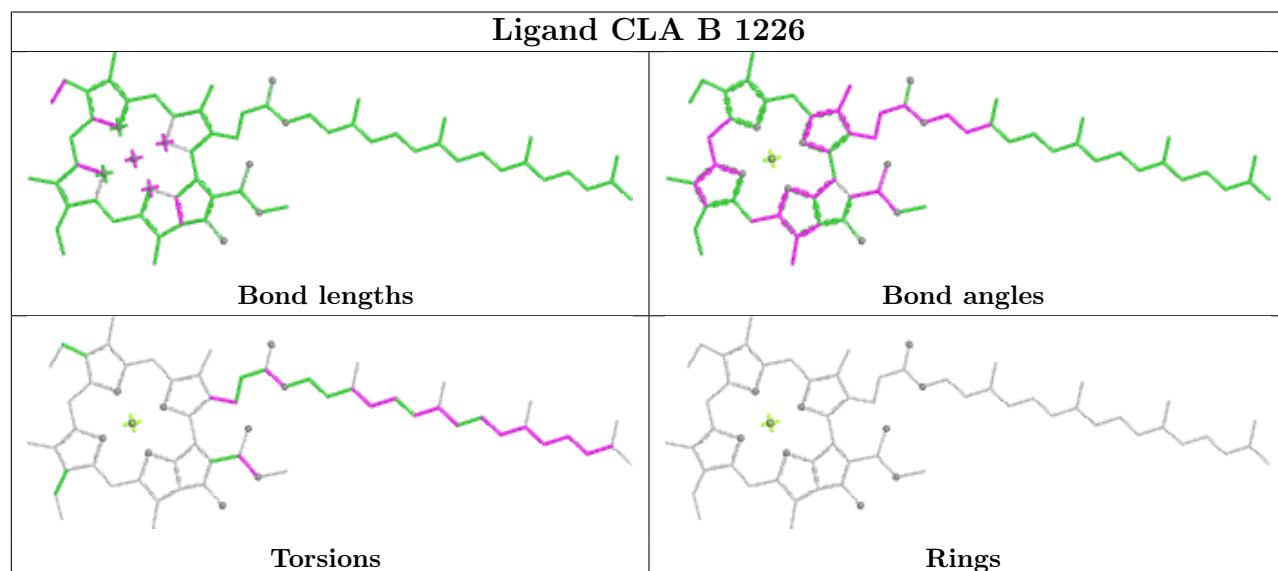
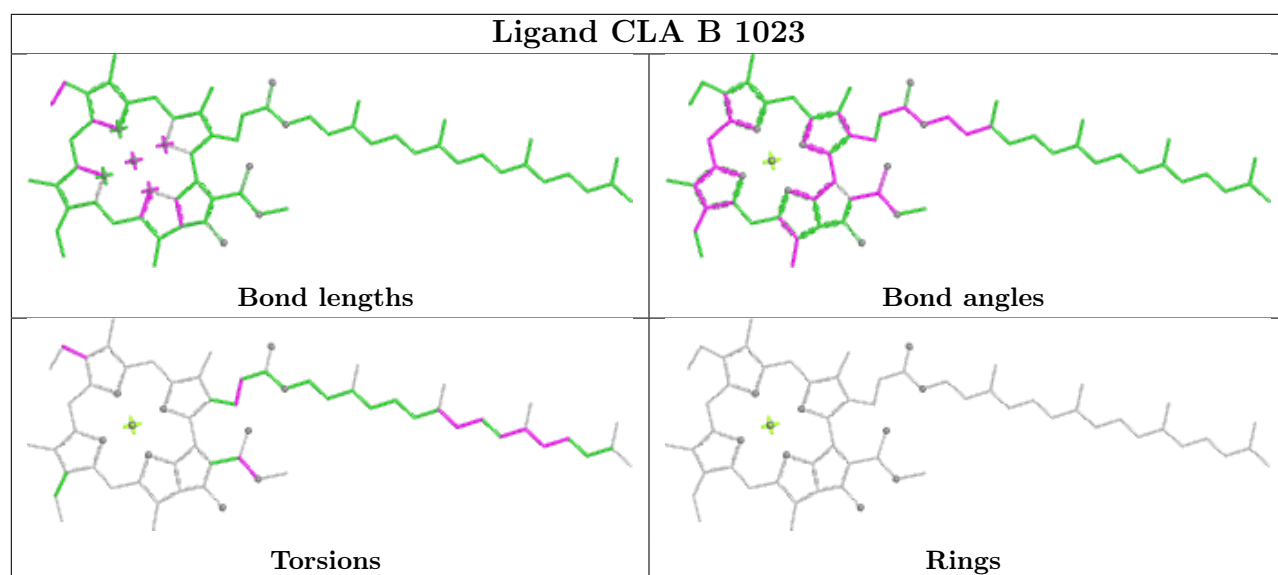
Ligand CLA 4 601**Ligand CLA 1 601**

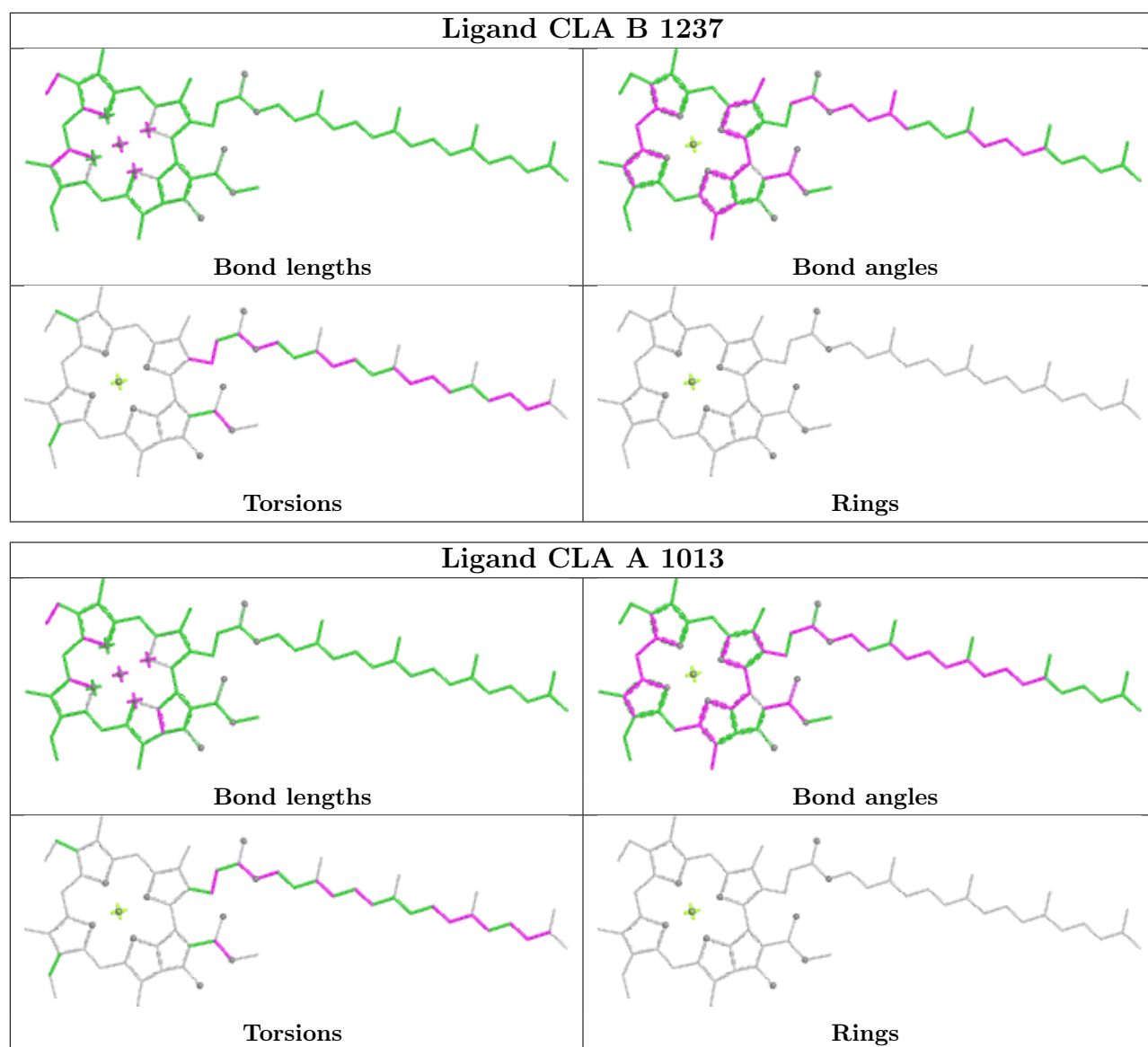


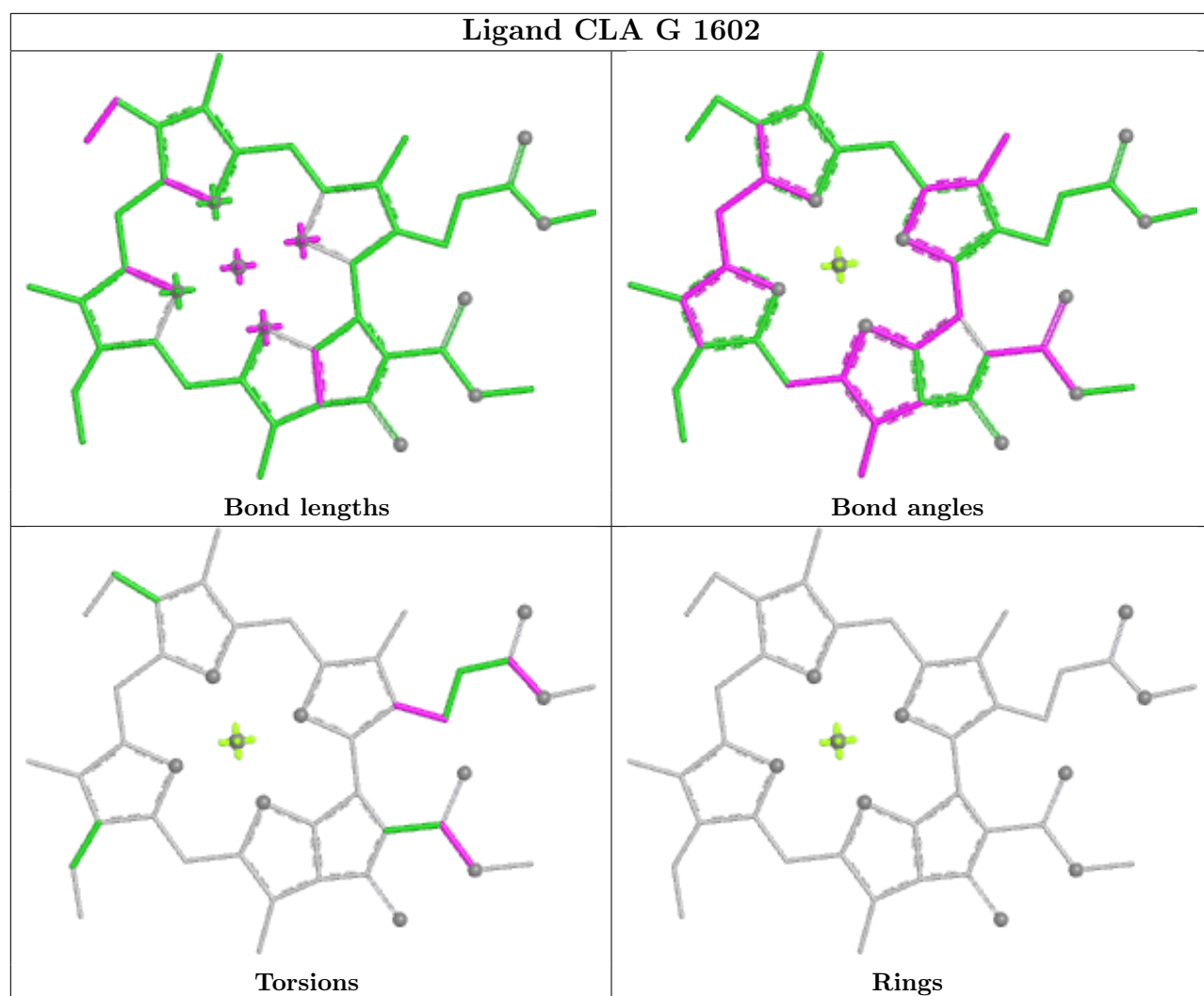
Ligand CLA B 1218	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA B 1231	
	
Bond lengths	Bond angles
	
Torsions	Rings

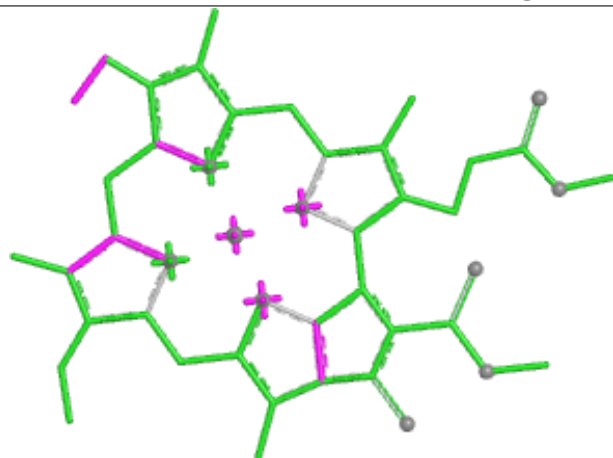
Ligand LUT 1 502	
	
Bond lengths	Bond angles
	
Torsions	Rings



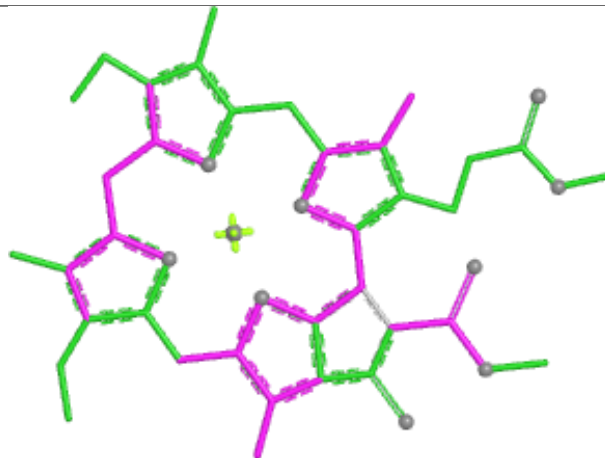




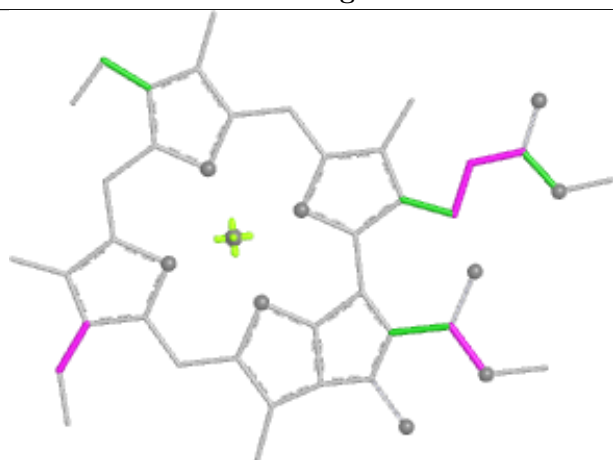
Ligand CLA 4 608



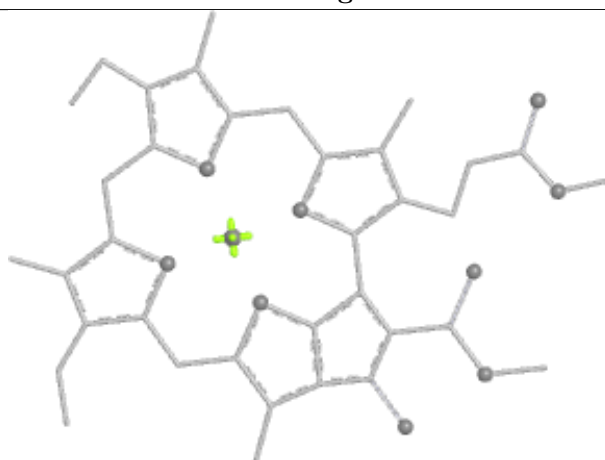
Bond lengths



Bond angles

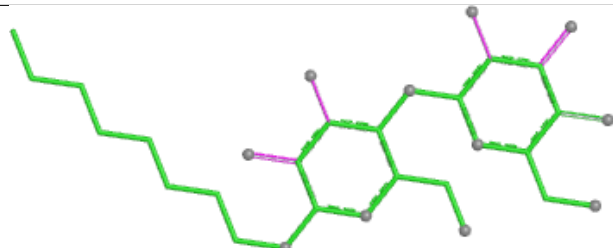


Torsions

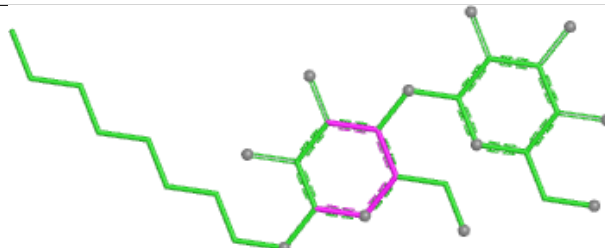


Rings

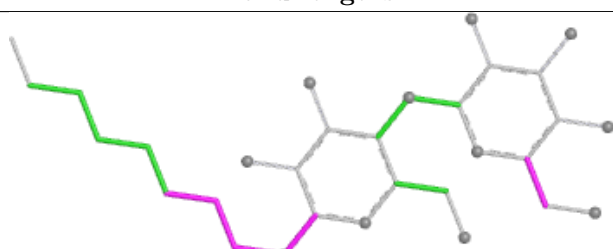
Ligand LMT G 5005



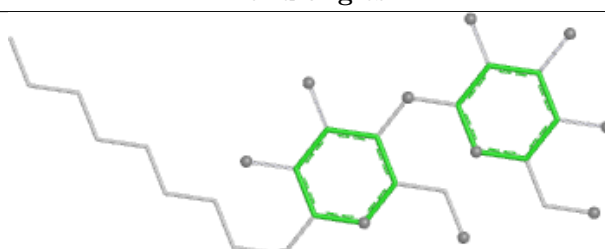
Bond lengths



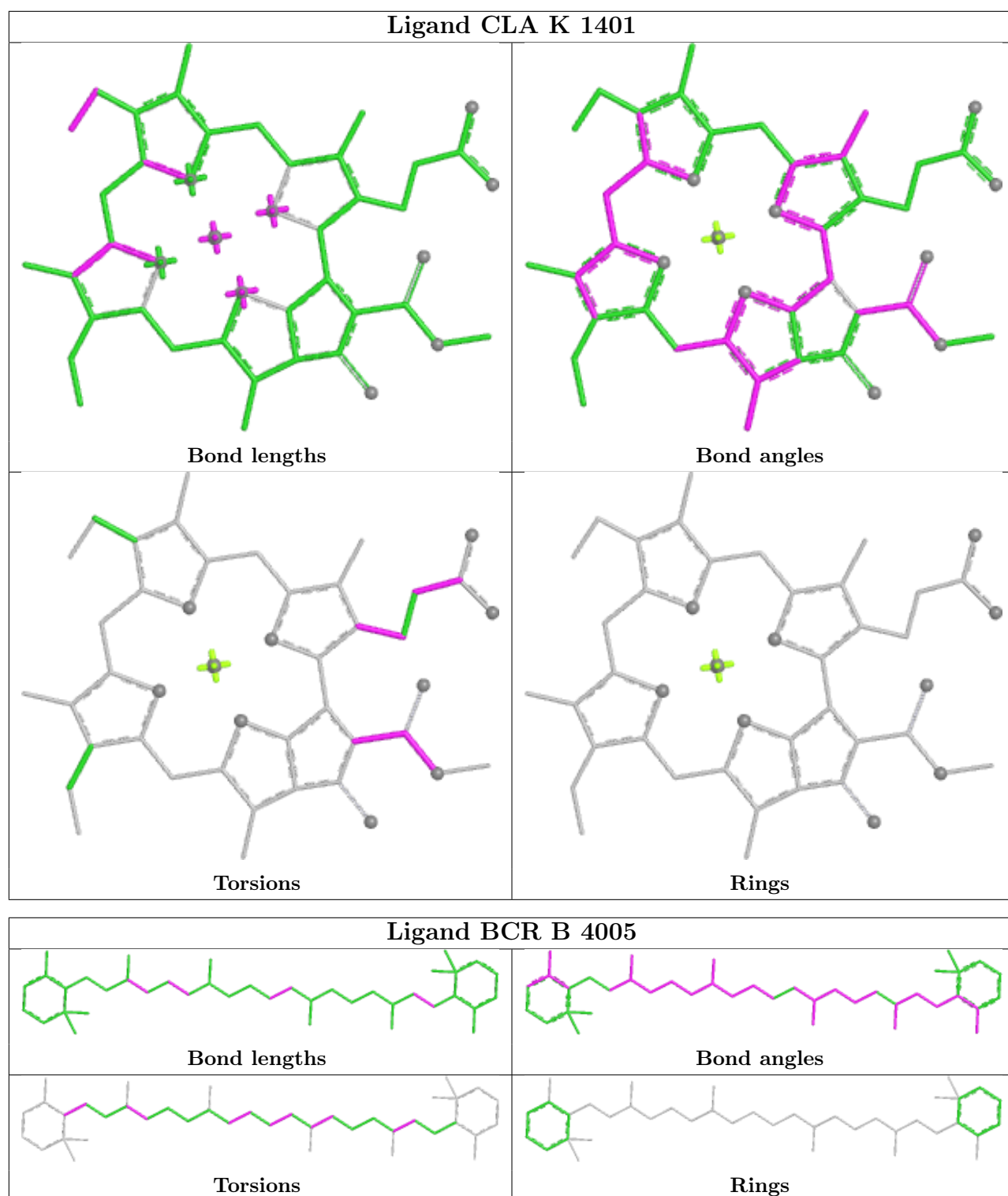
Bond angles

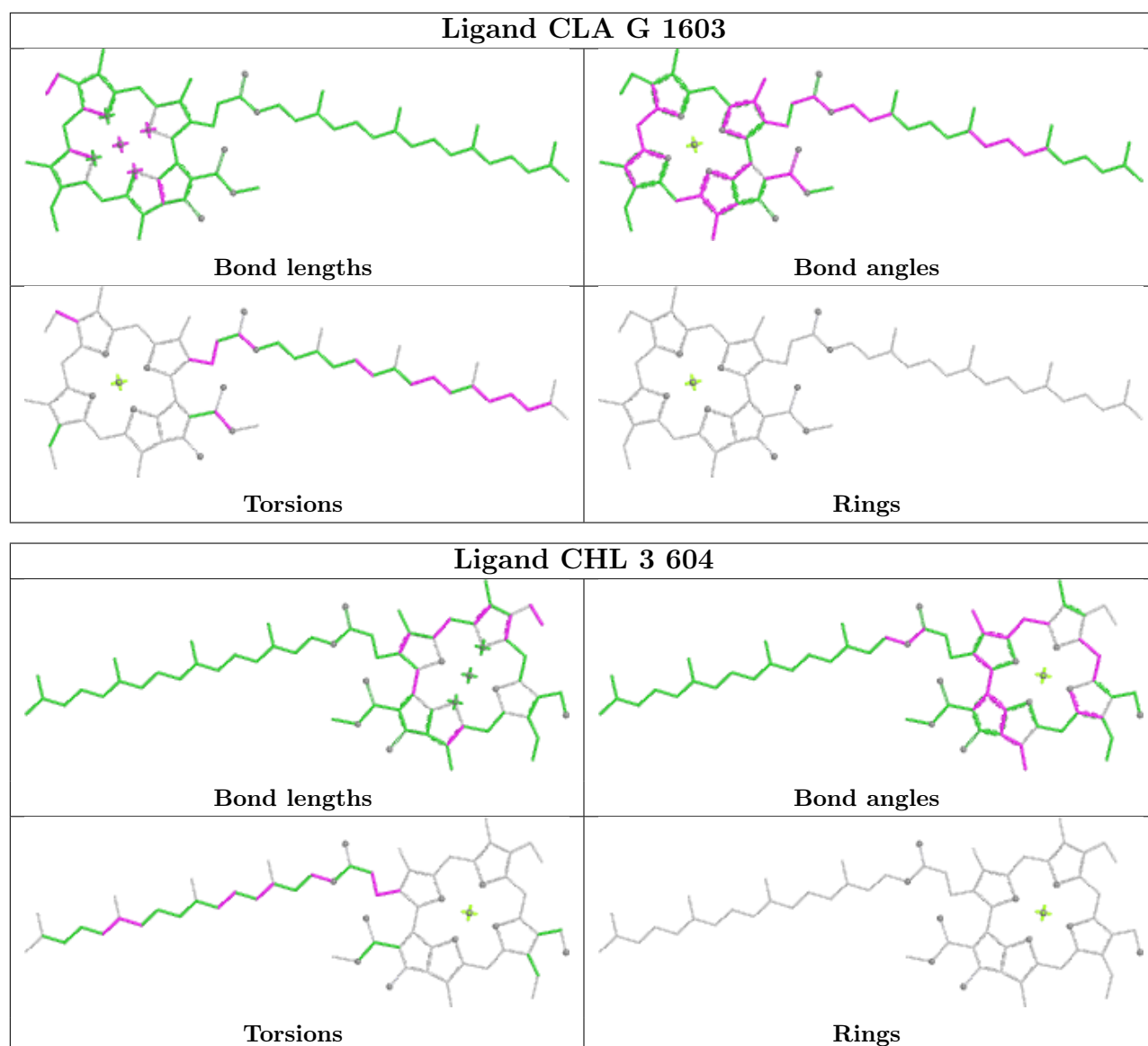


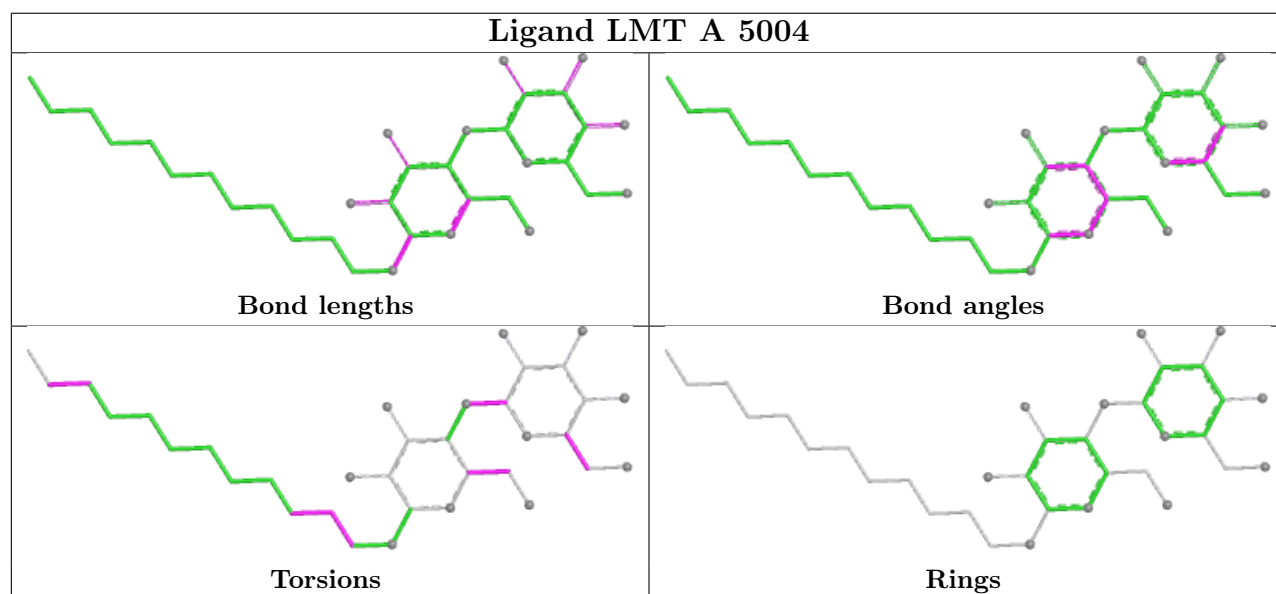
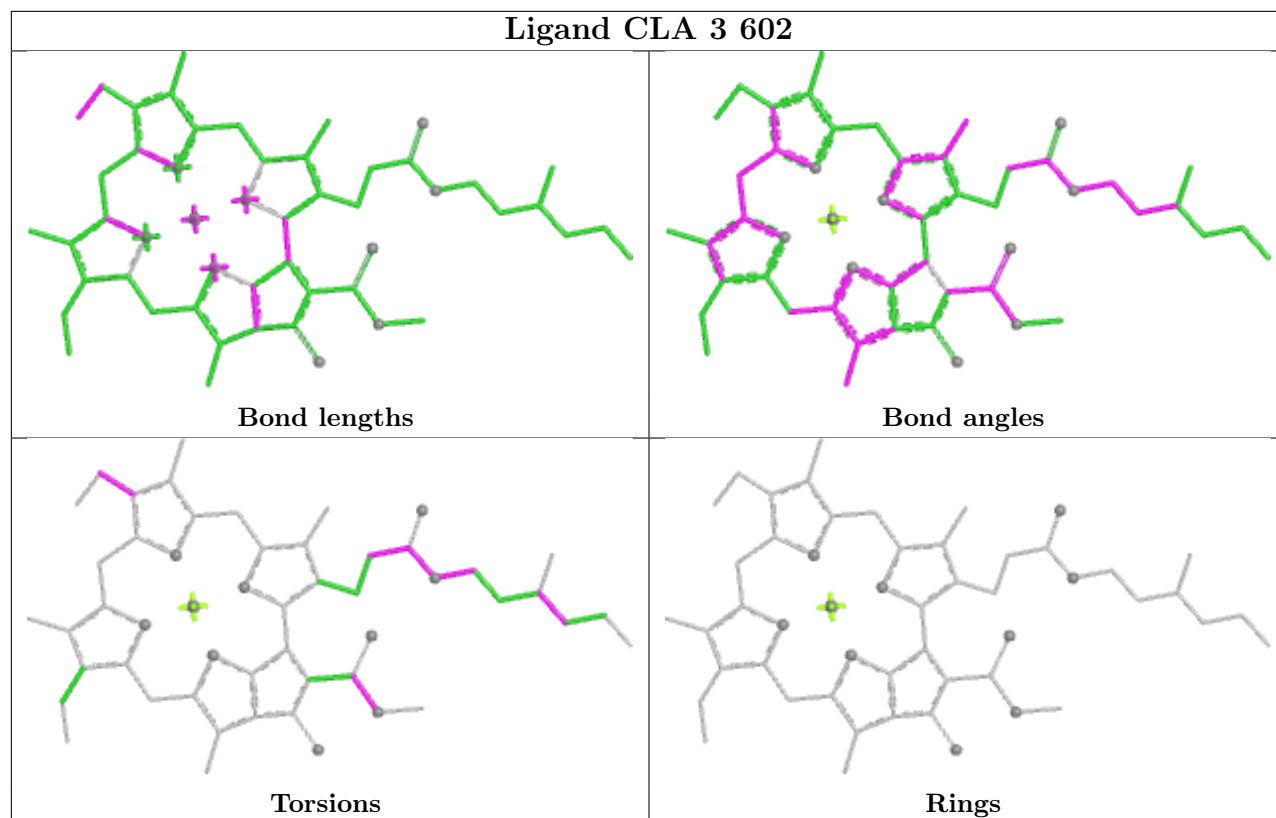
Torsions

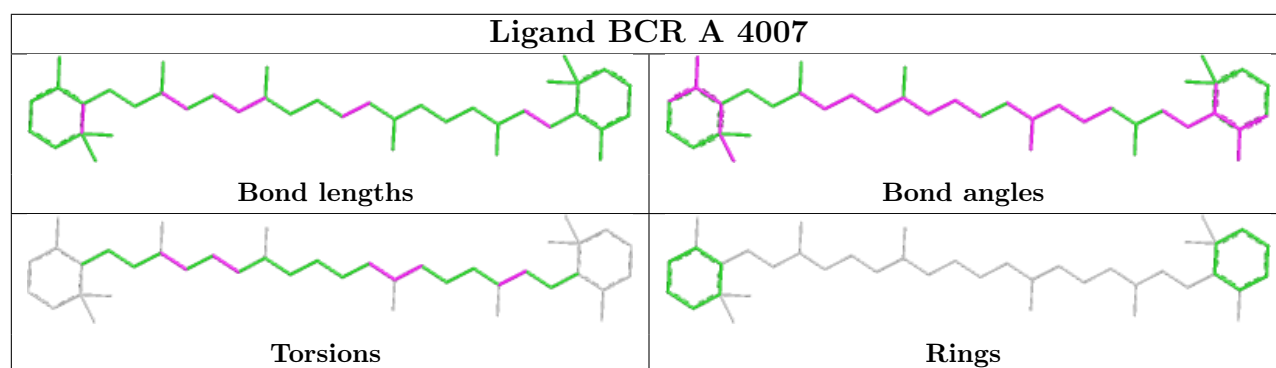
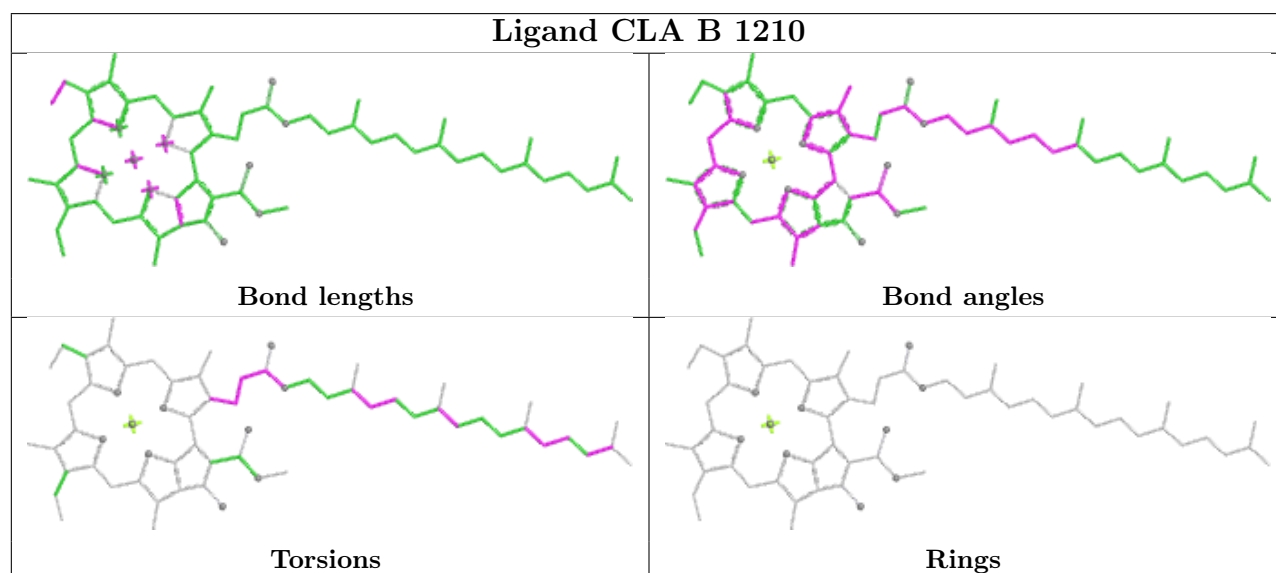
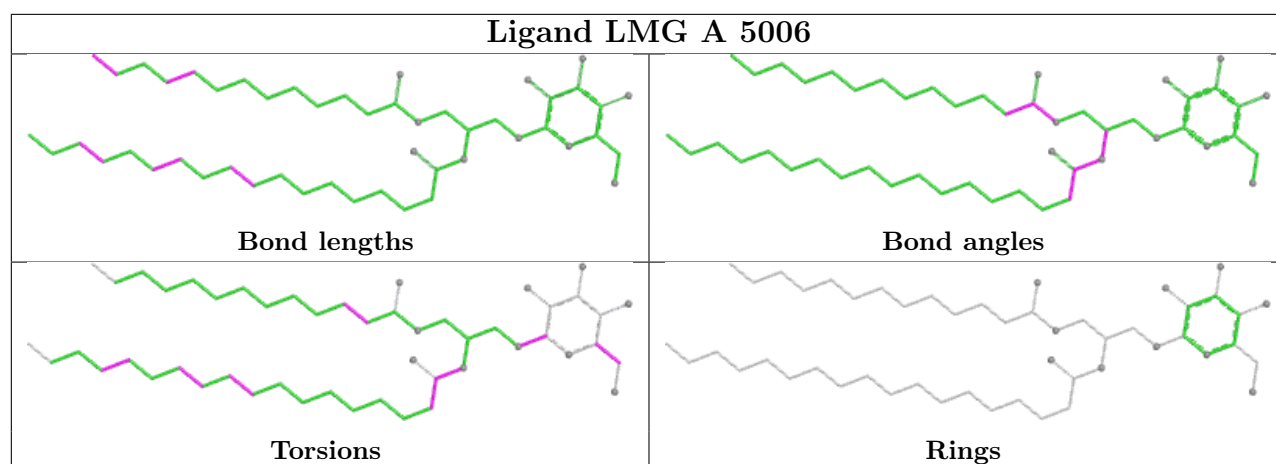


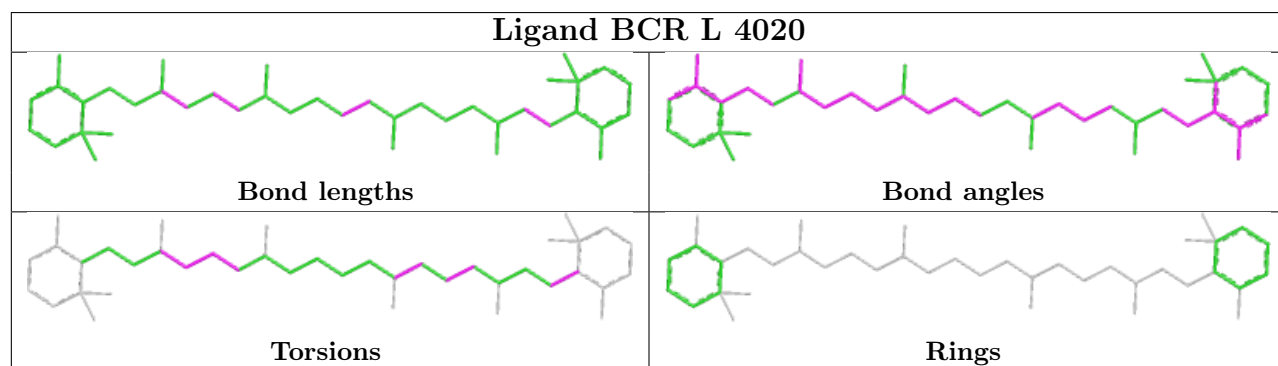
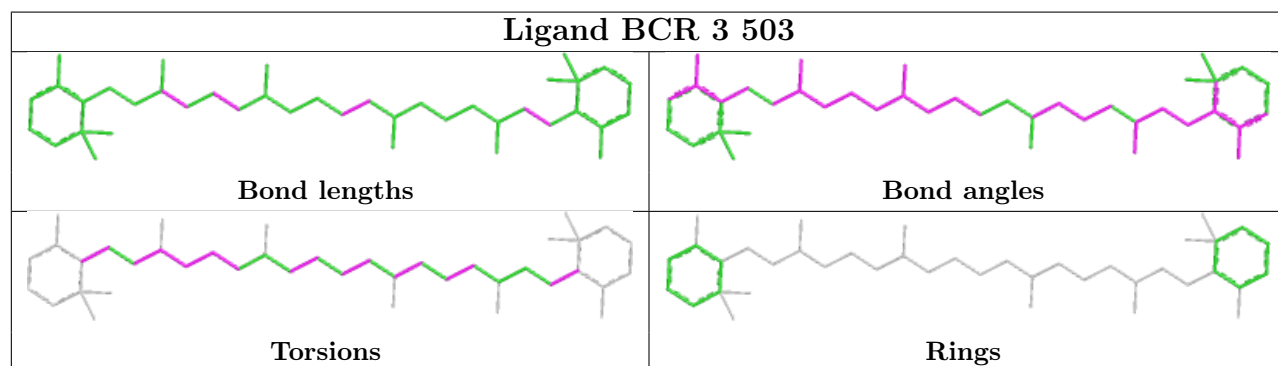
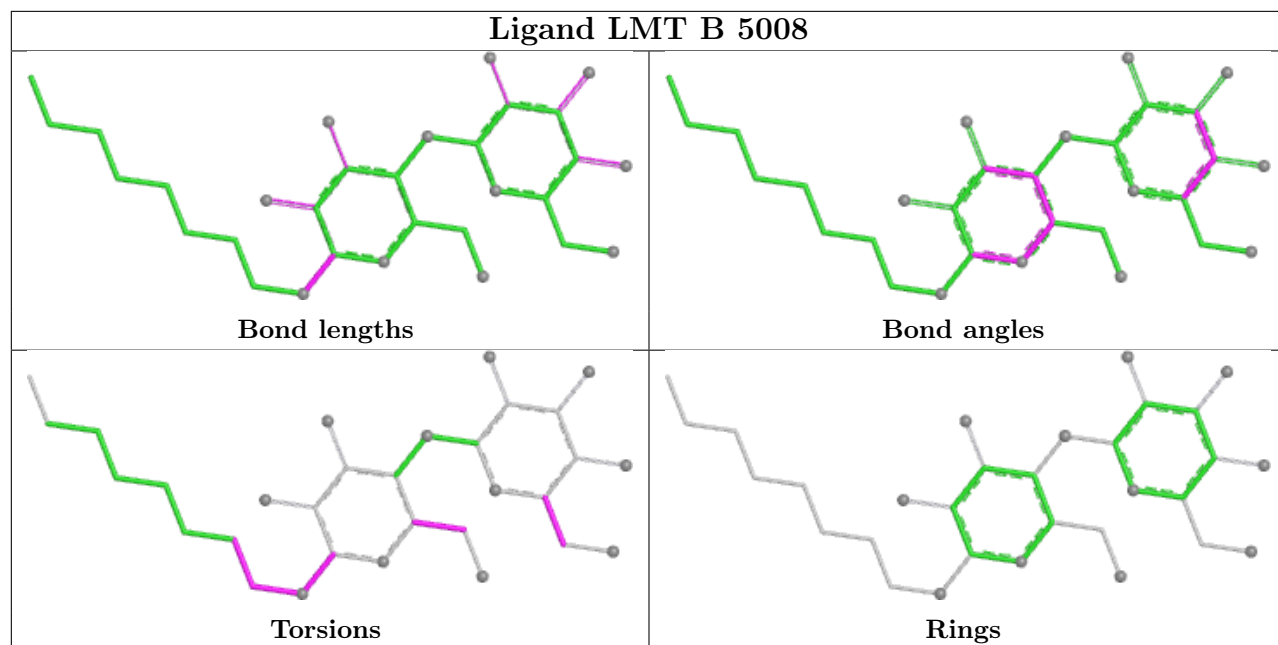
Rings

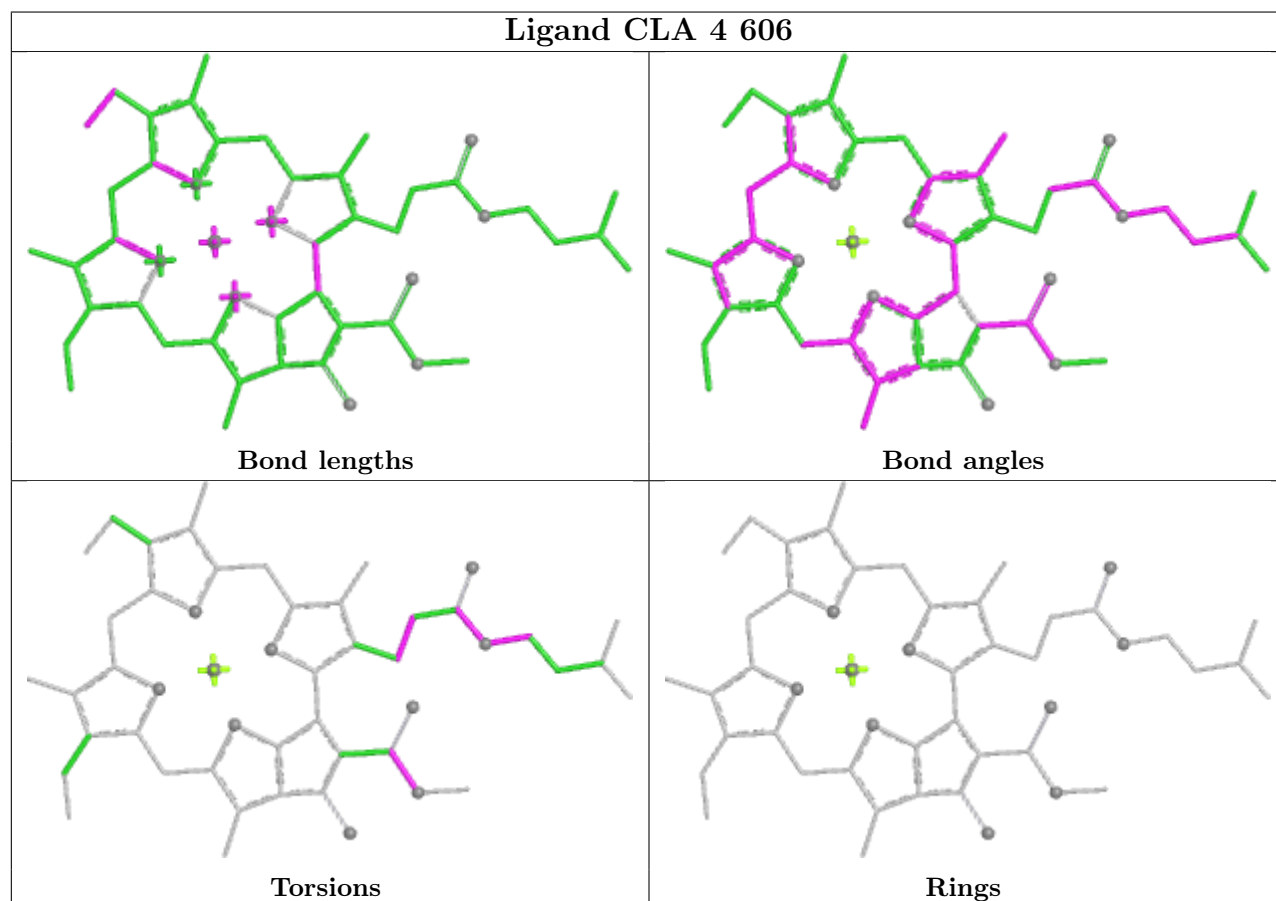
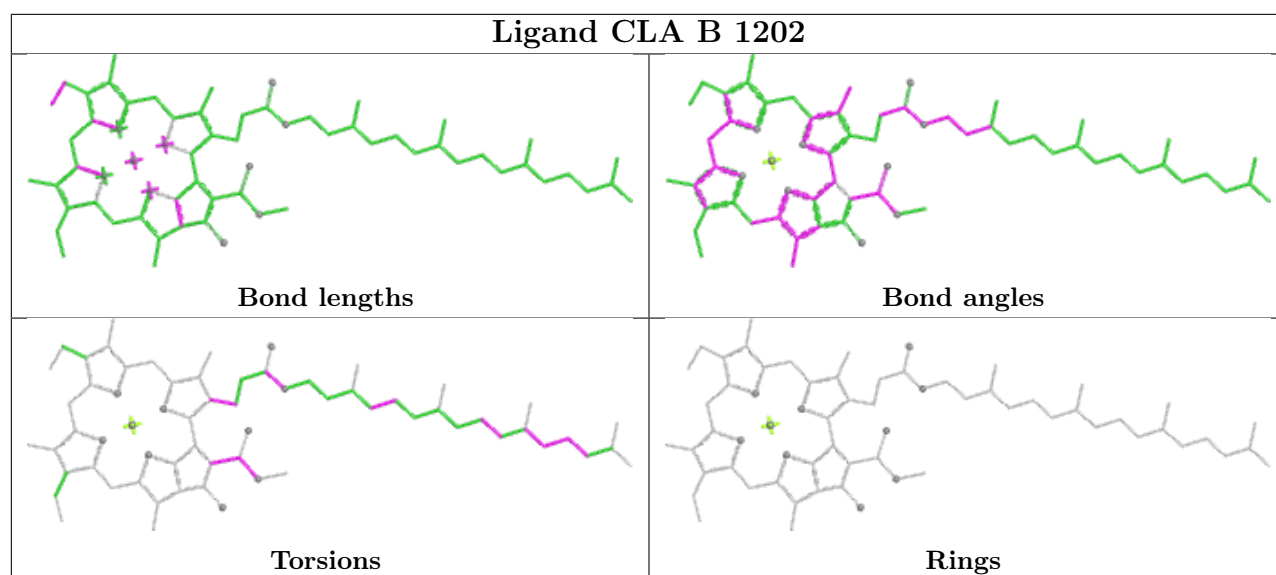


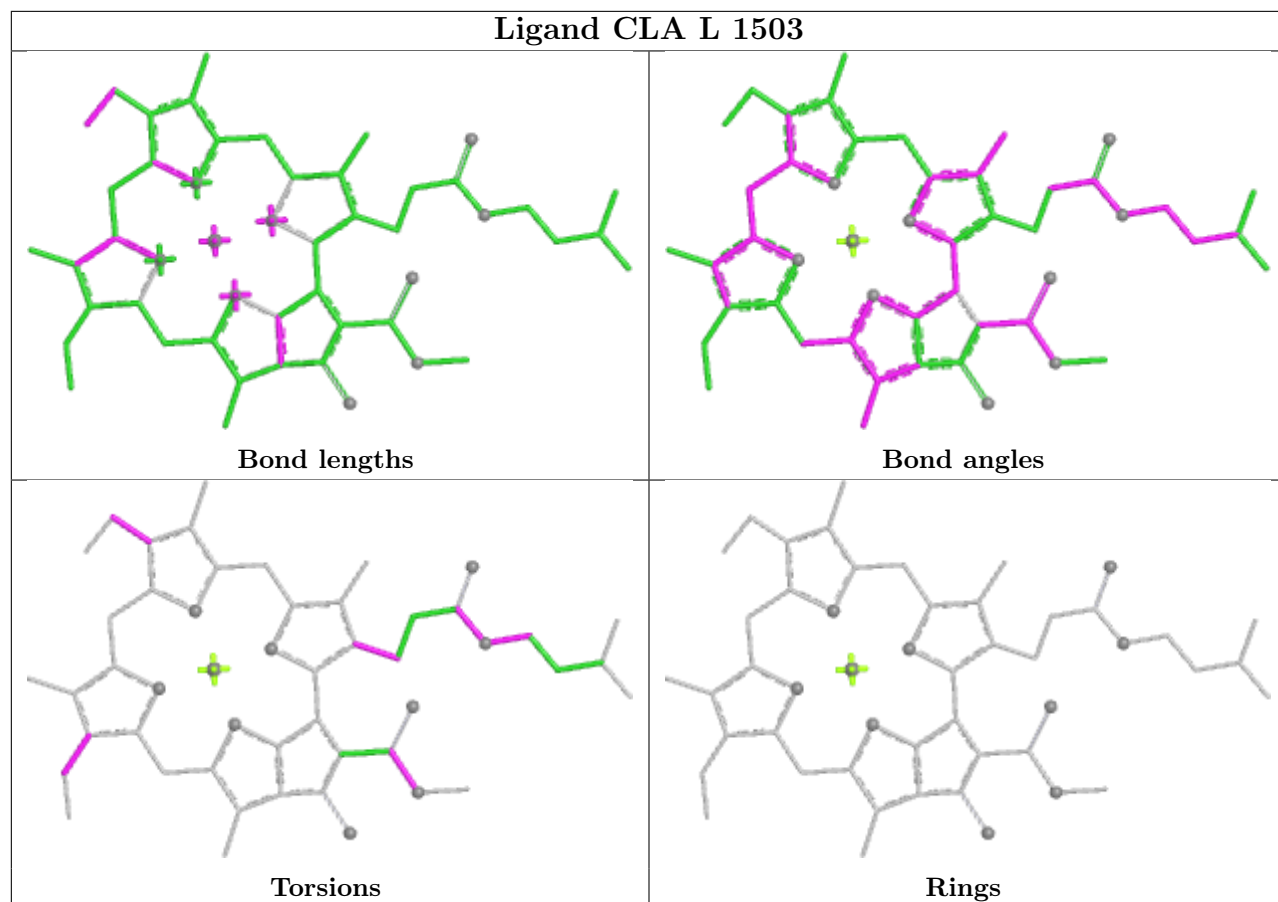


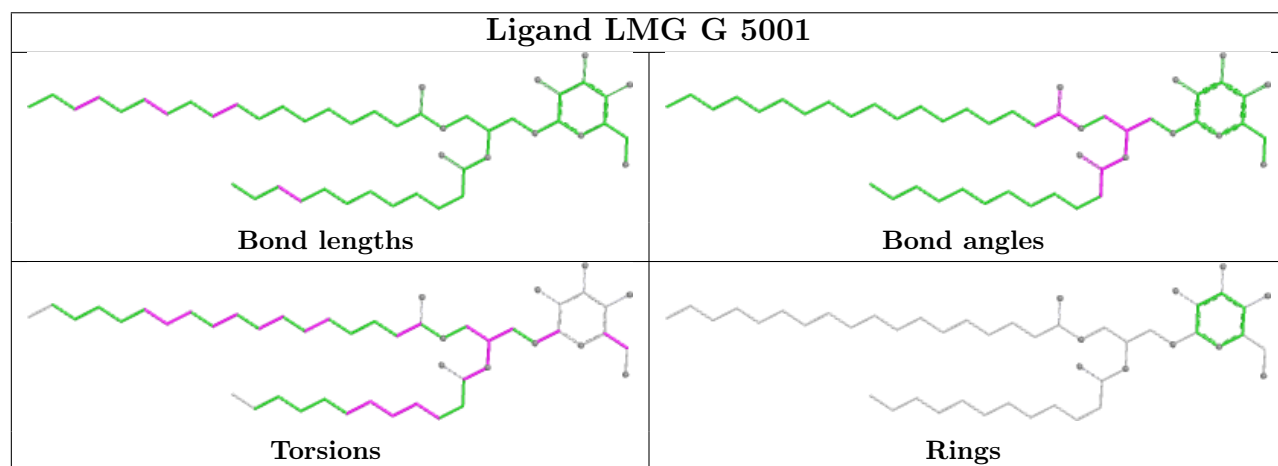
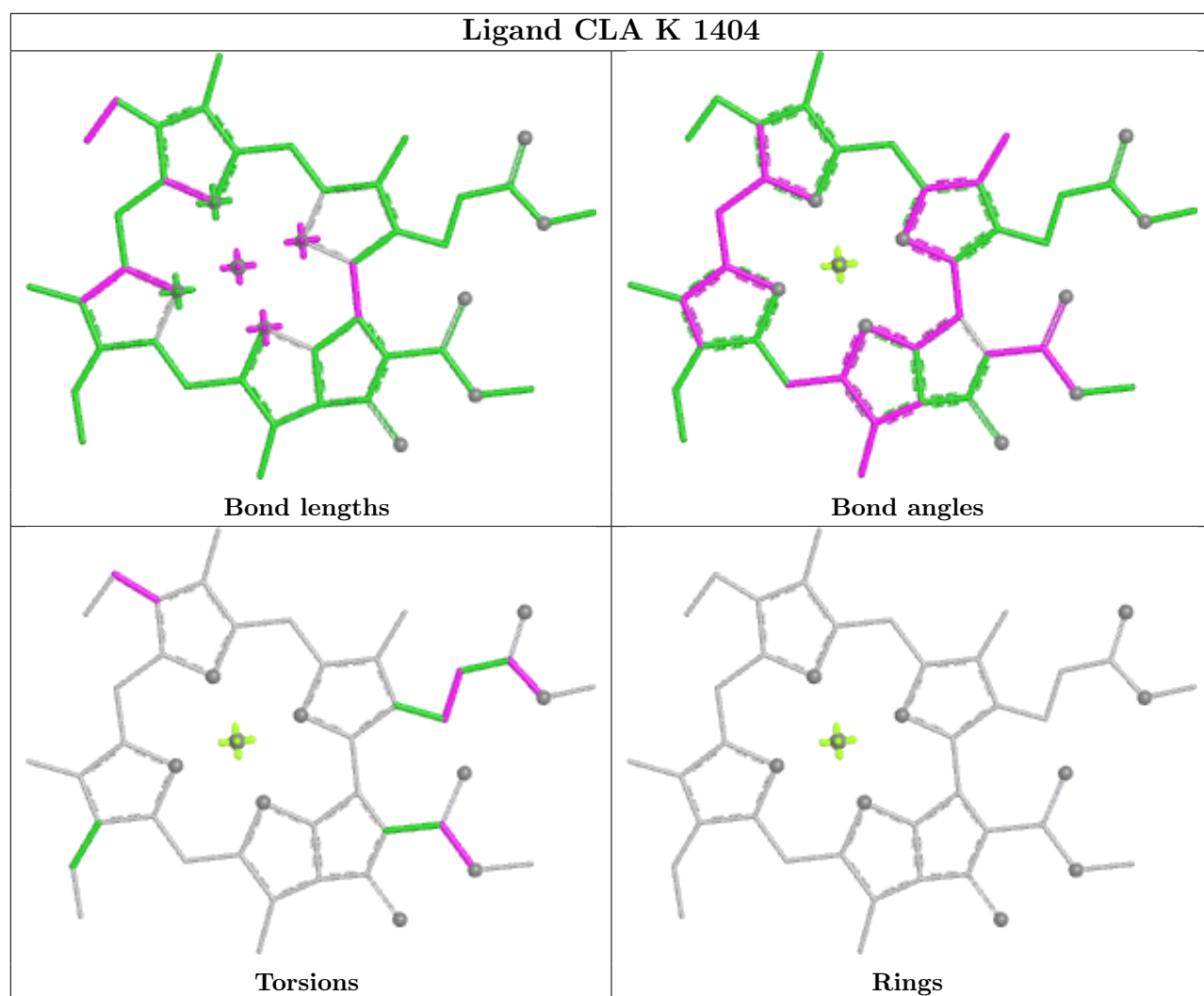


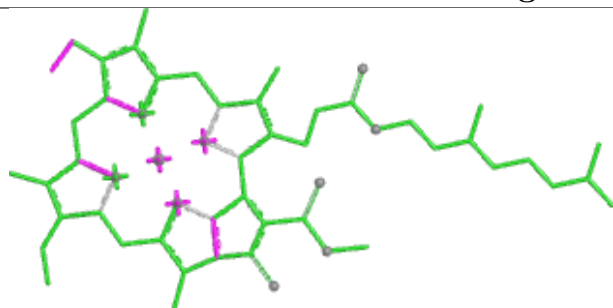




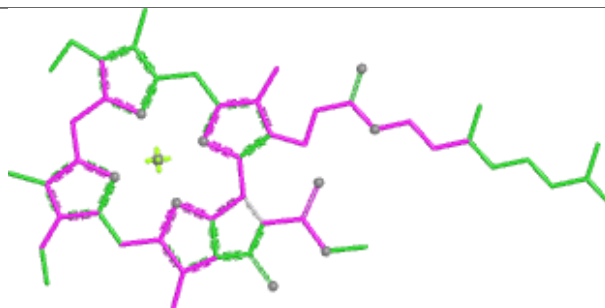




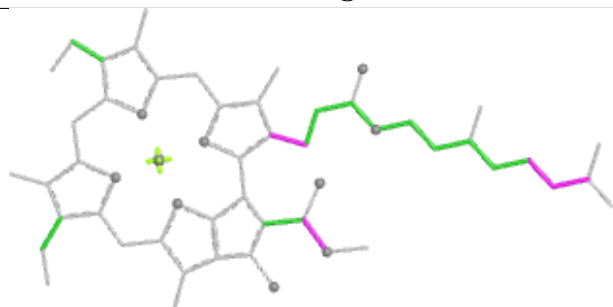


Ligand CLA 1 603

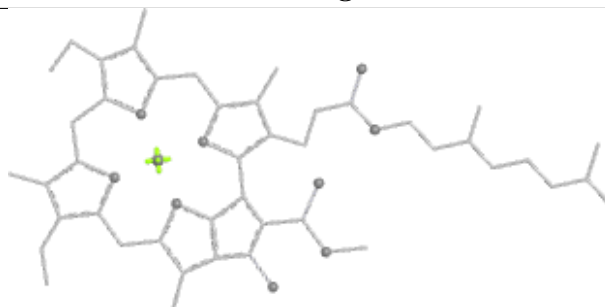
Bond lengths



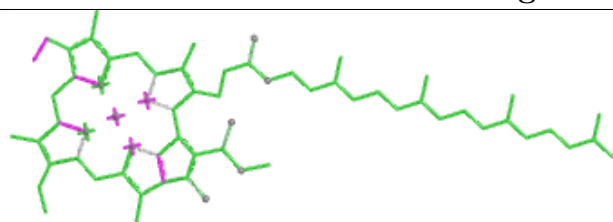
Bond angles



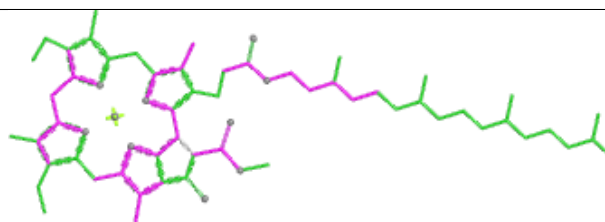
Torsions



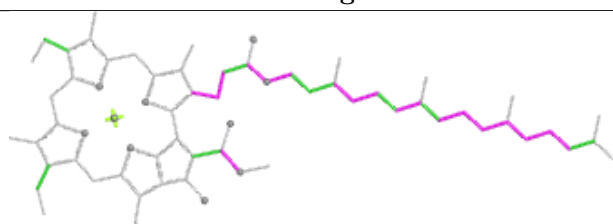
Rings

Ligand CLA 3 610

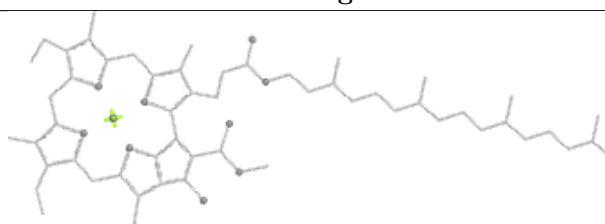
Bond lengths



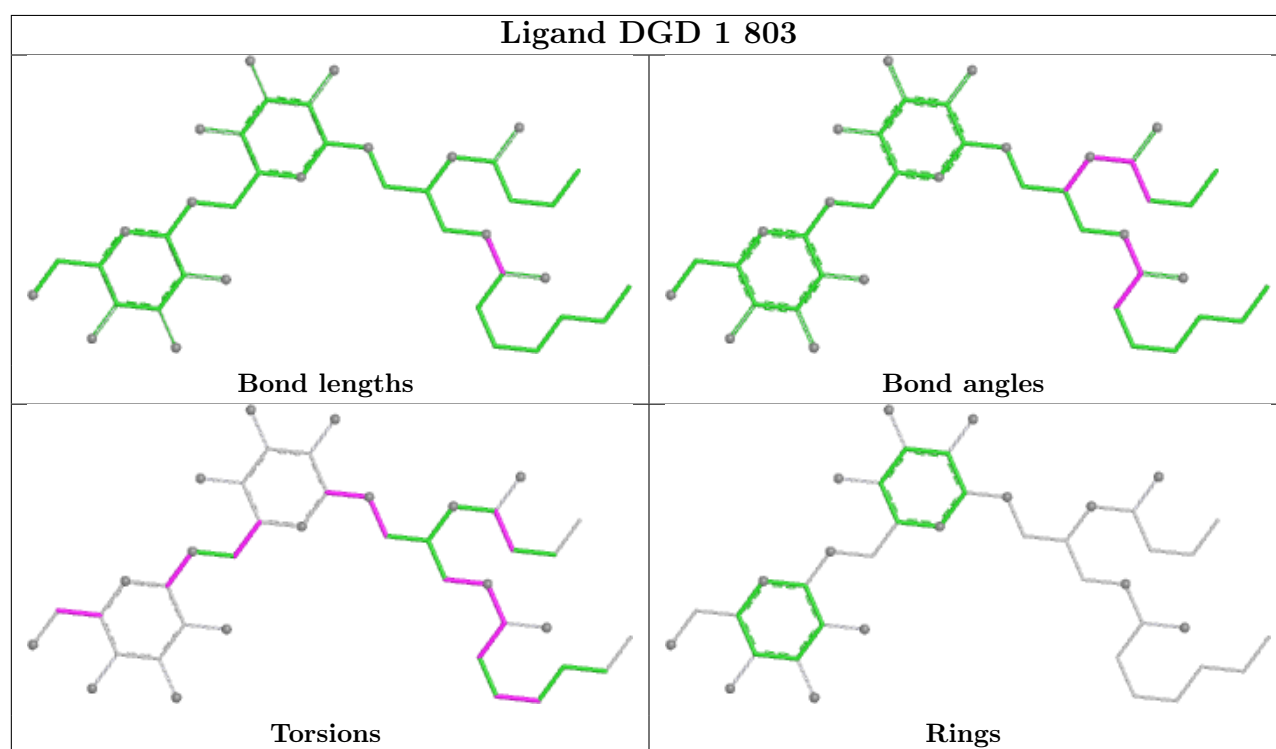
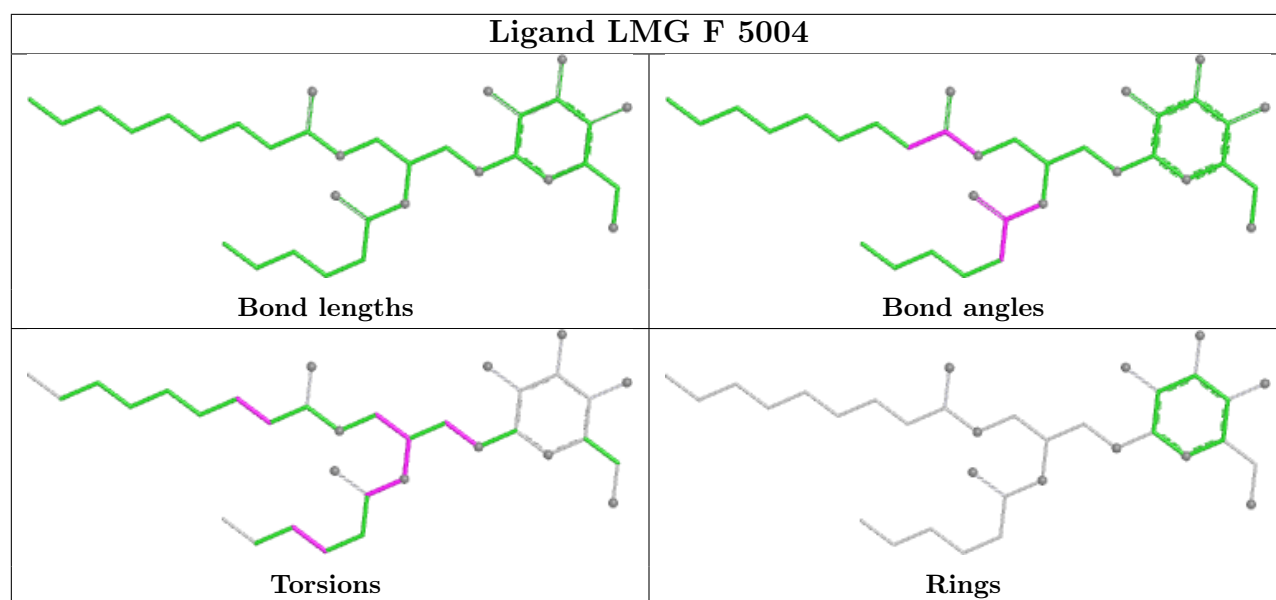
Bond angles

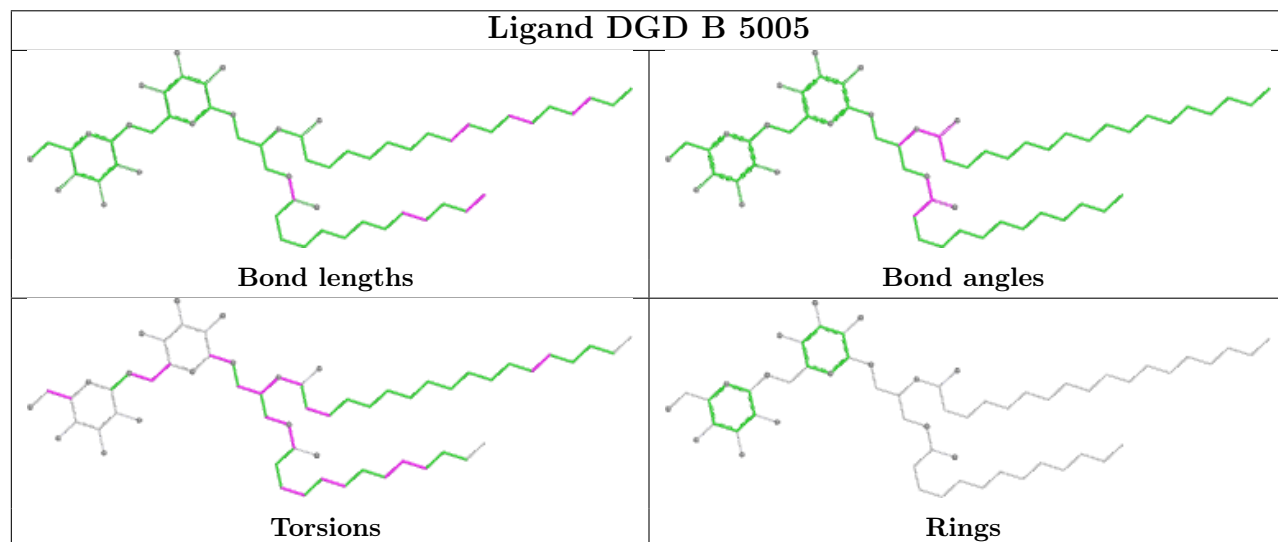
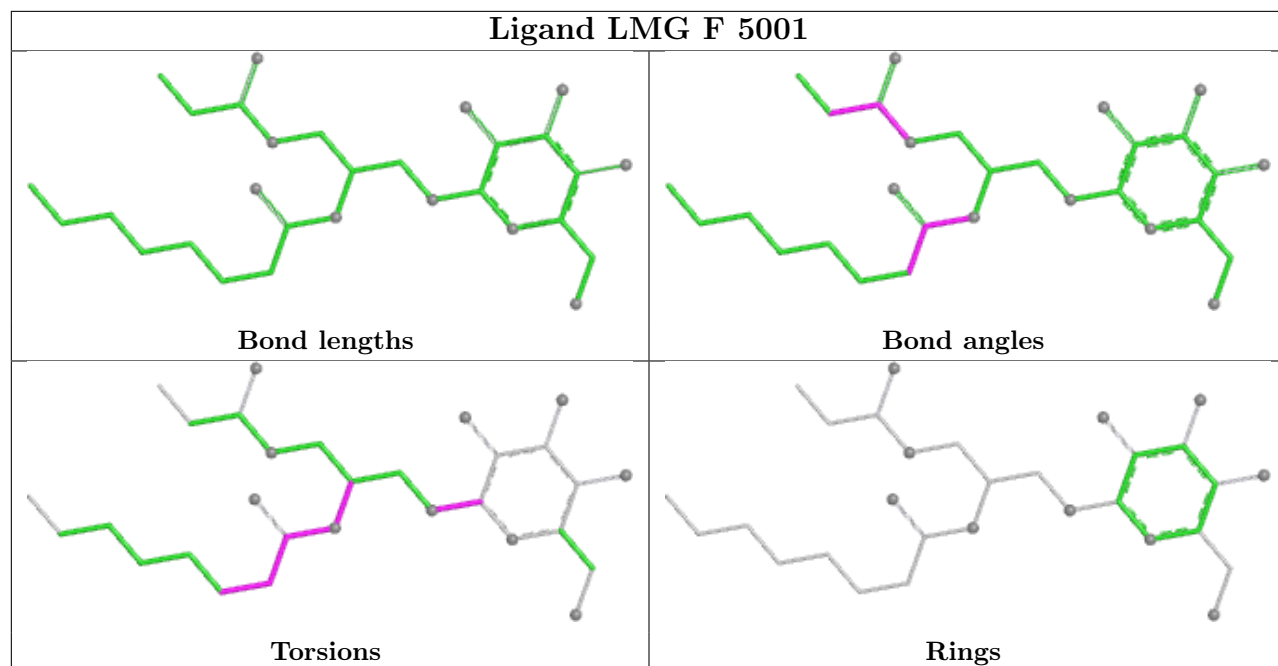


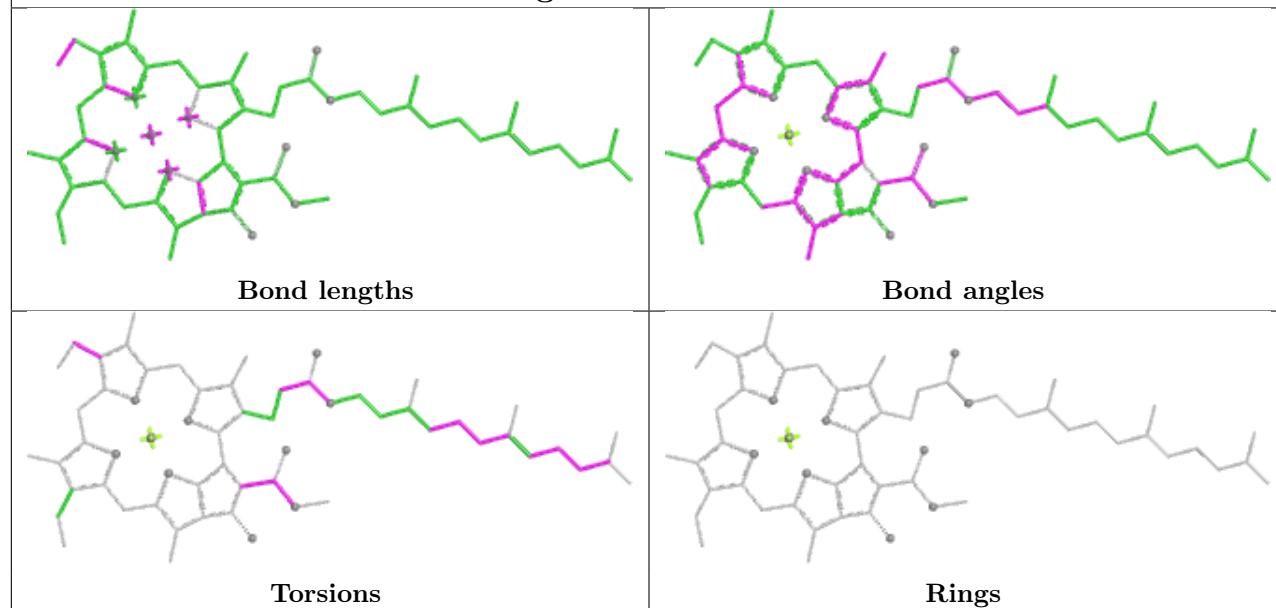
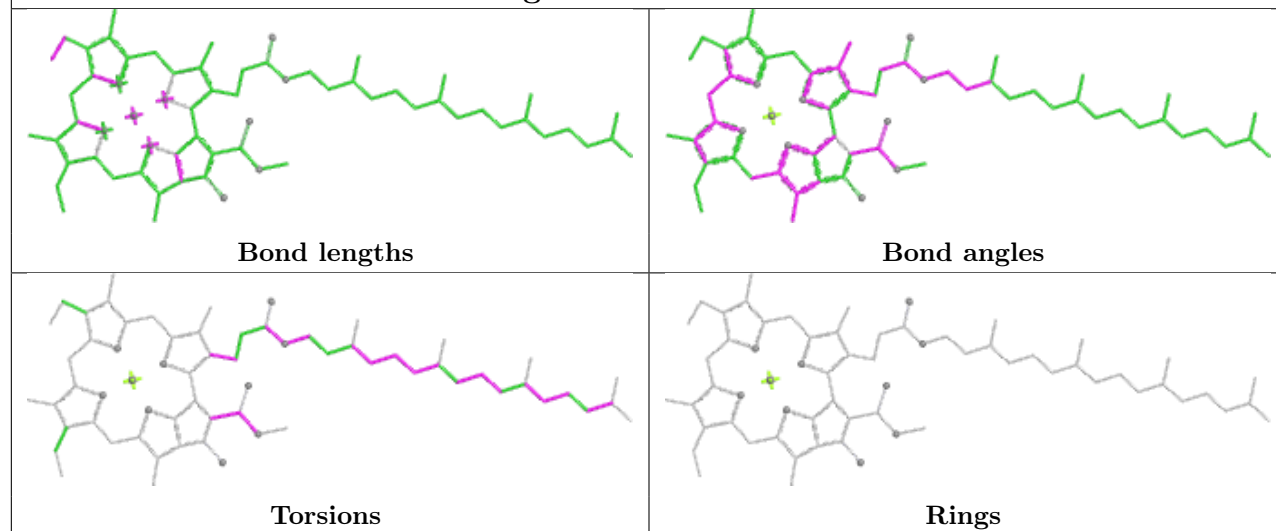
Torsions

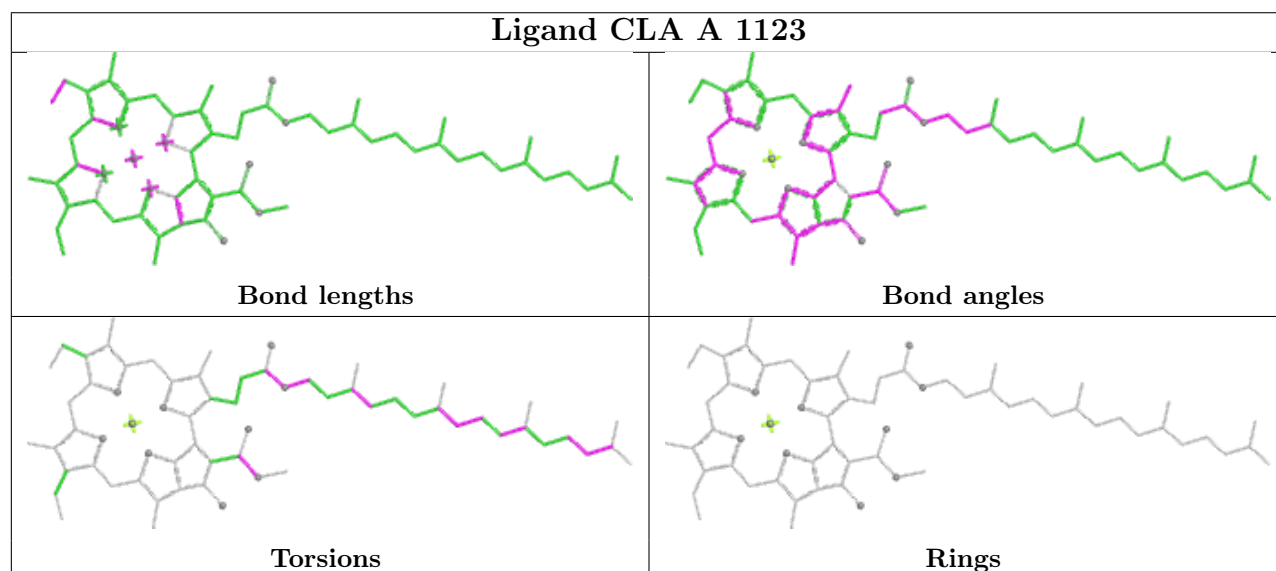
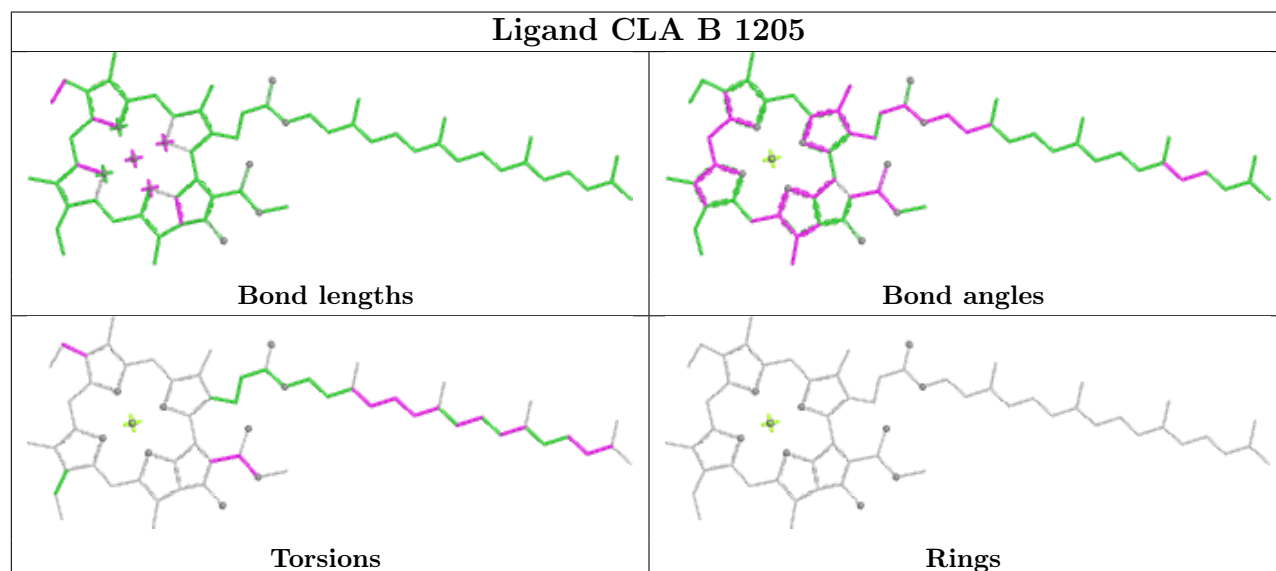
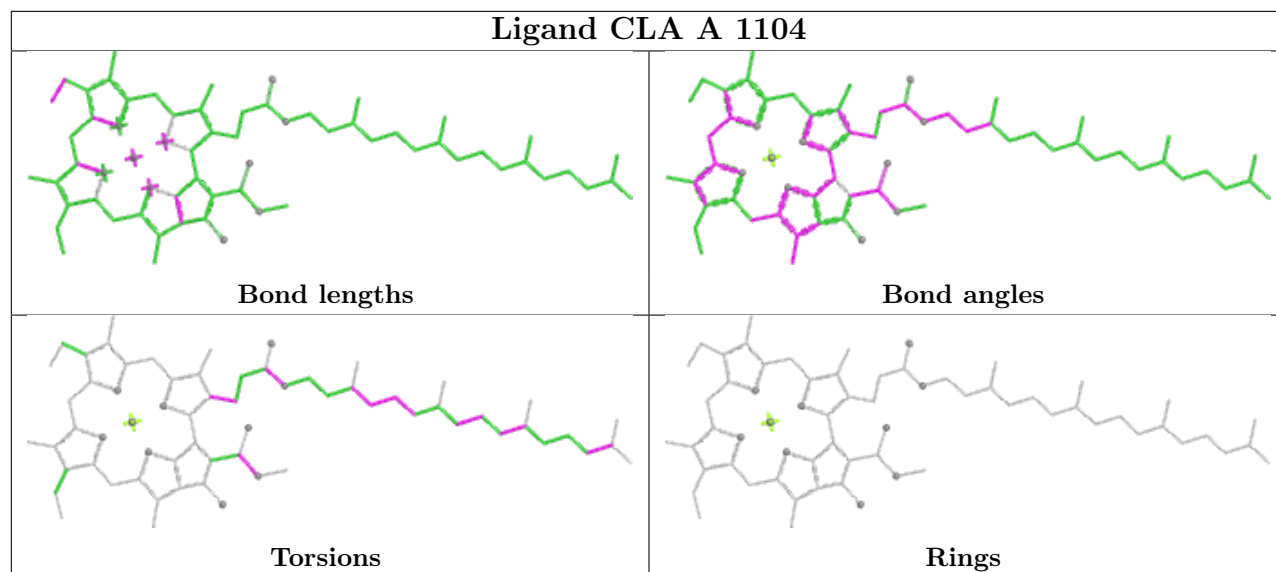


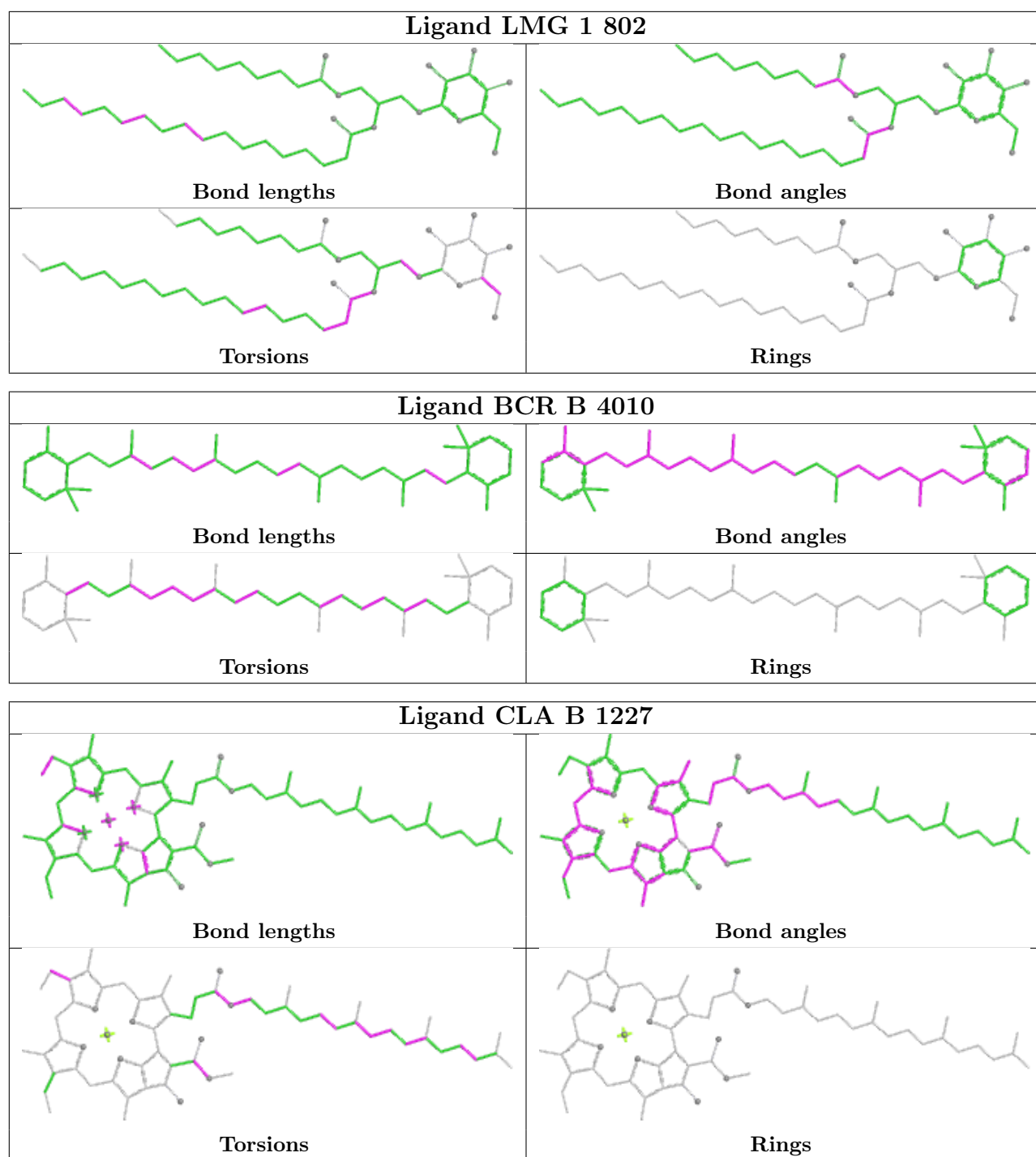
Rings

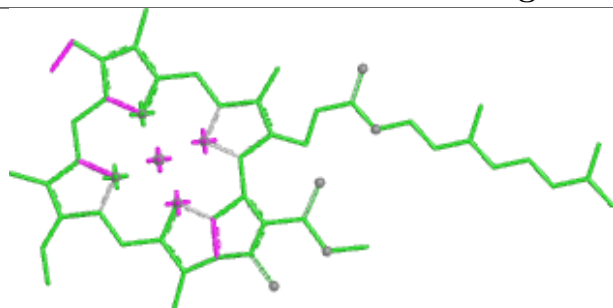




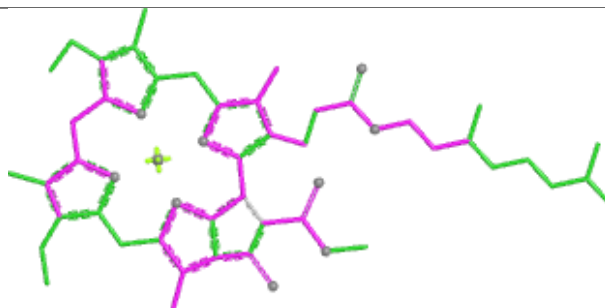
Ligand CLA 1 614**Ligand CLA B 1222**



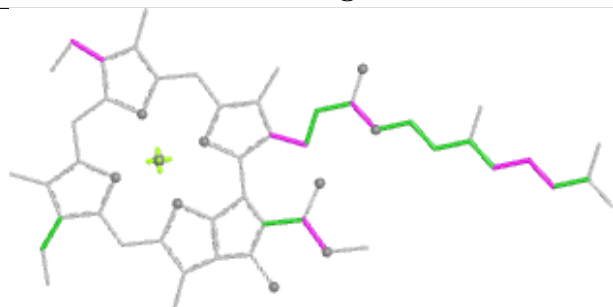


Ligand CLA 3 605

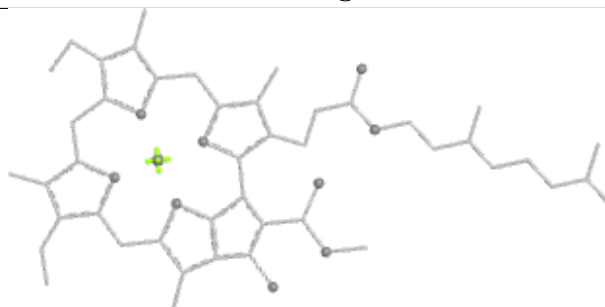
Bond lengths



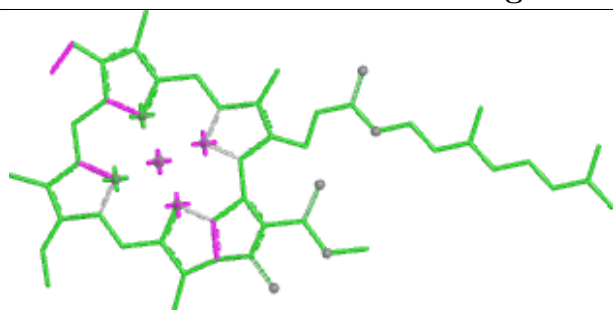
Bond angles



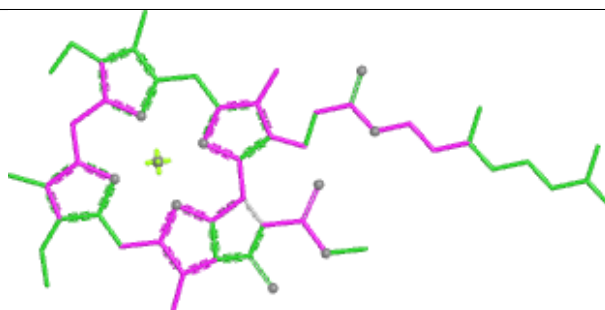
Torsions



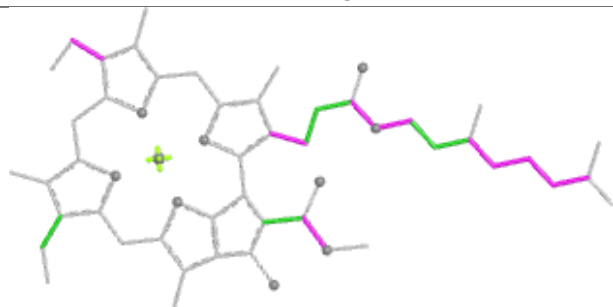
Rings

Ligand CLA B 1232

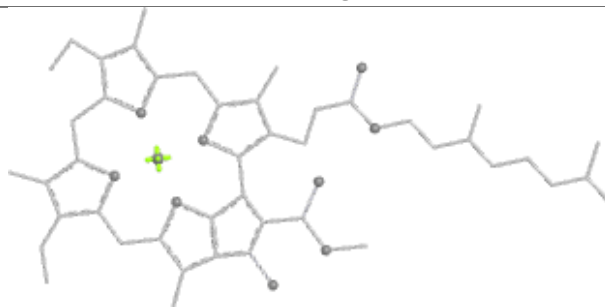
Bond lengths



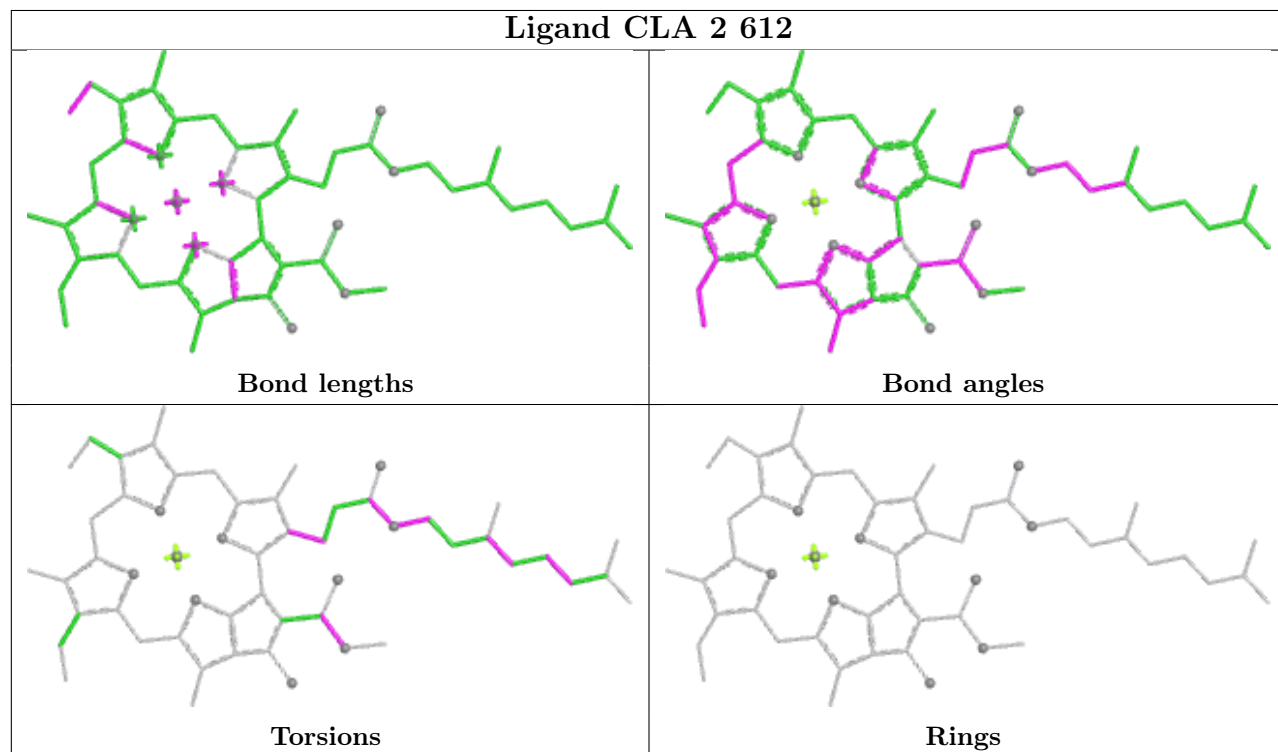
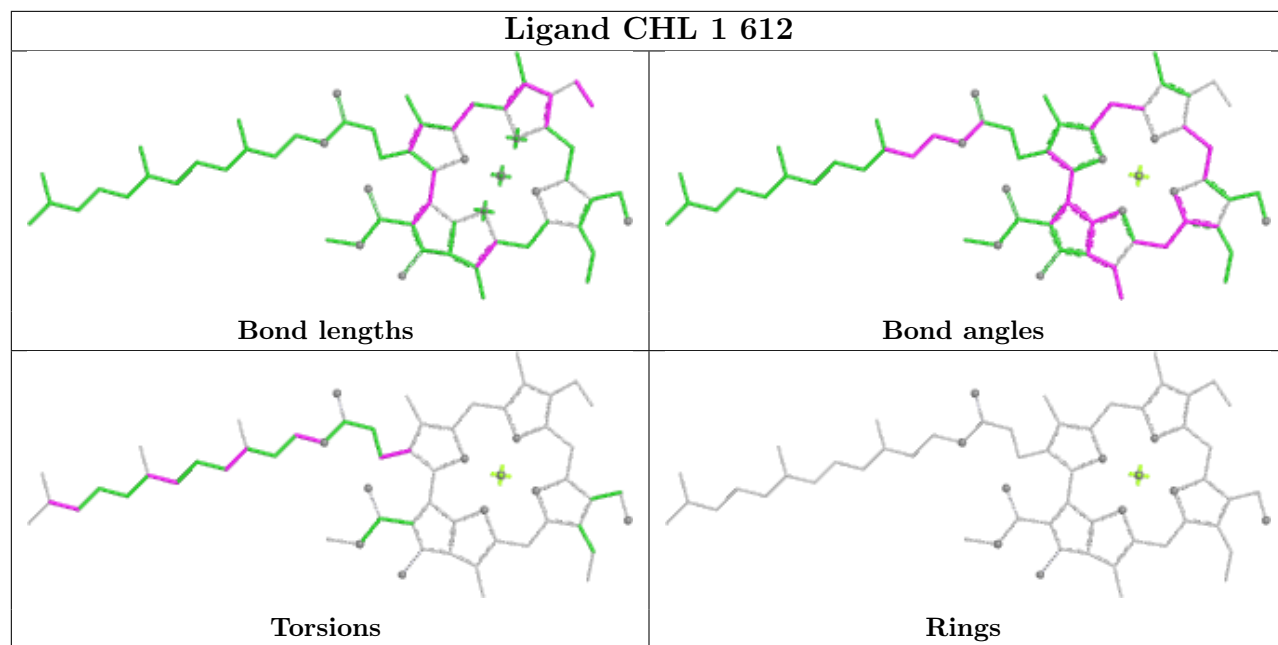
Bond angles

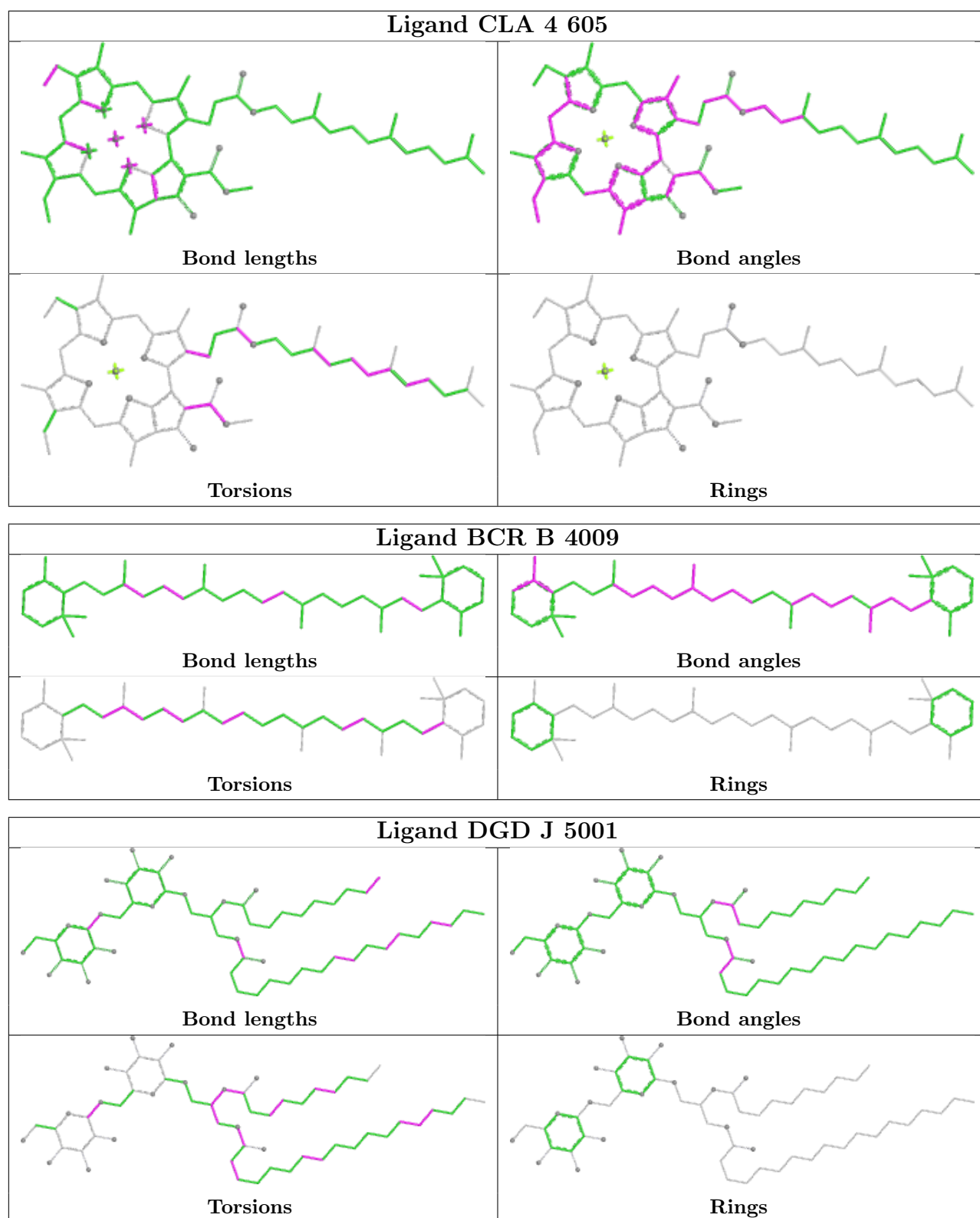


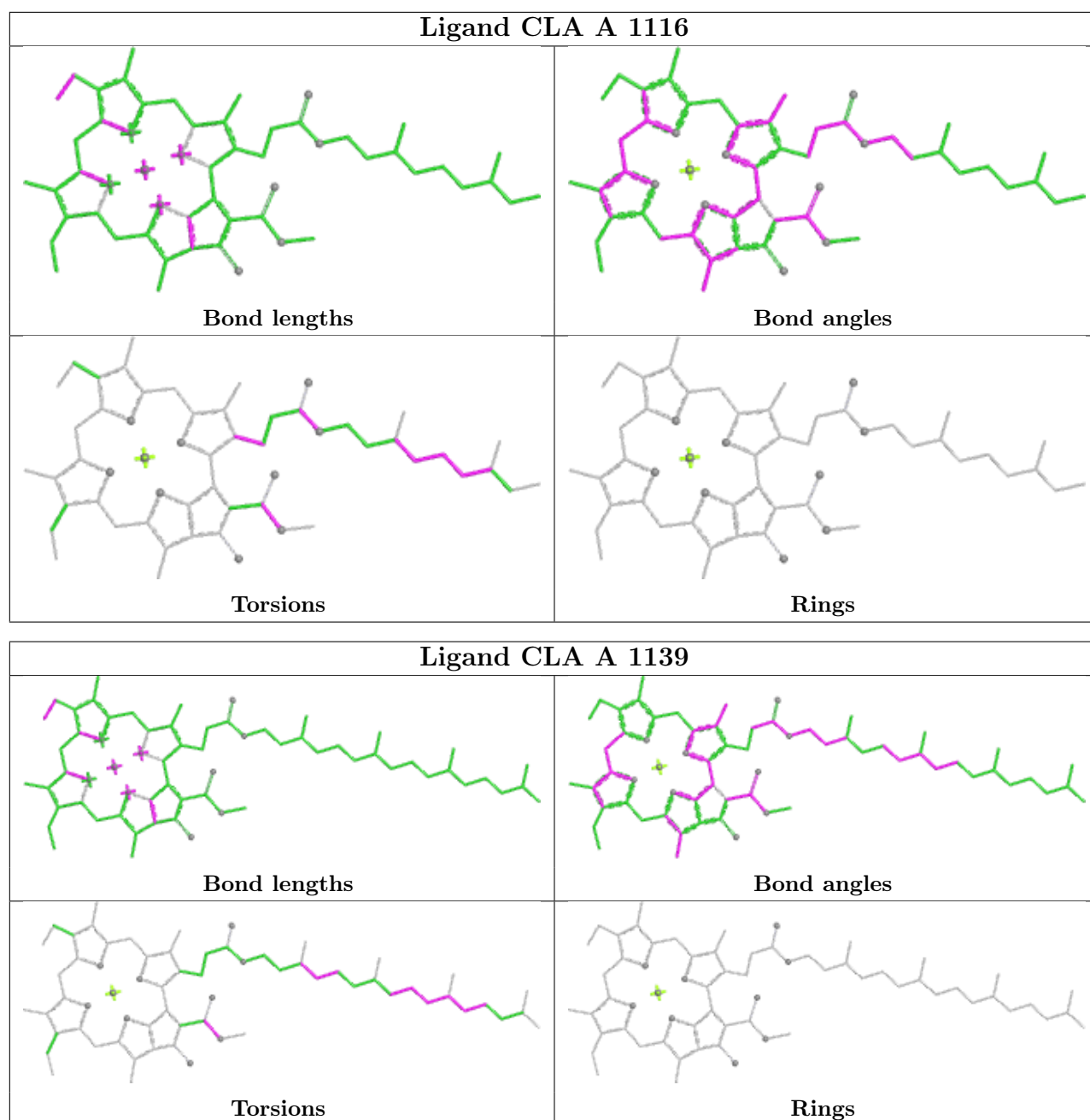
Torsions



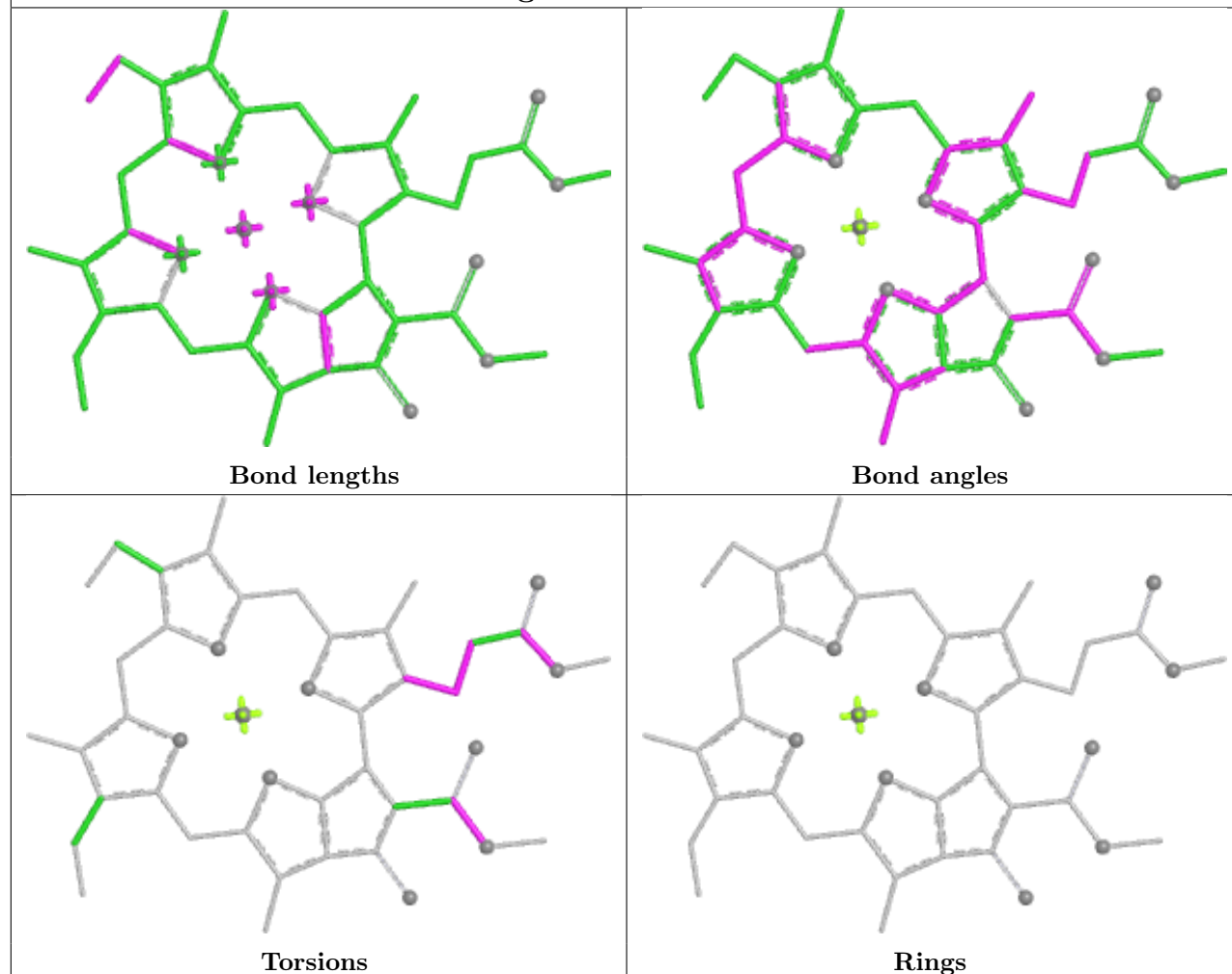
Rings



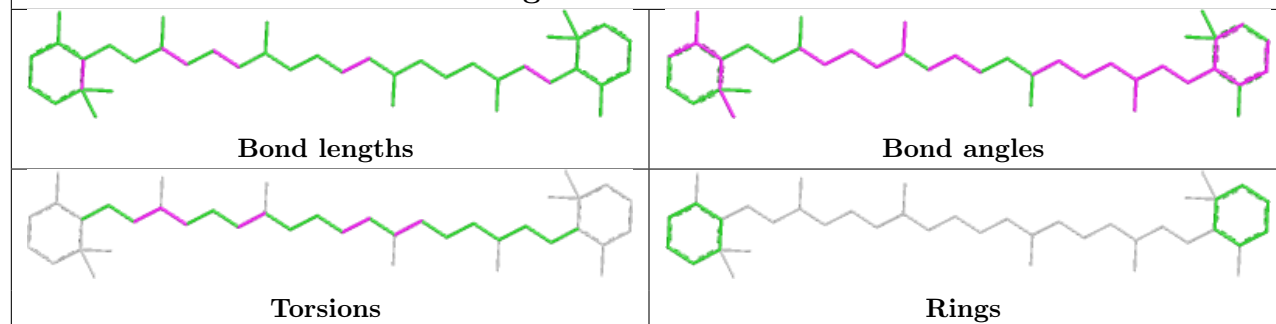




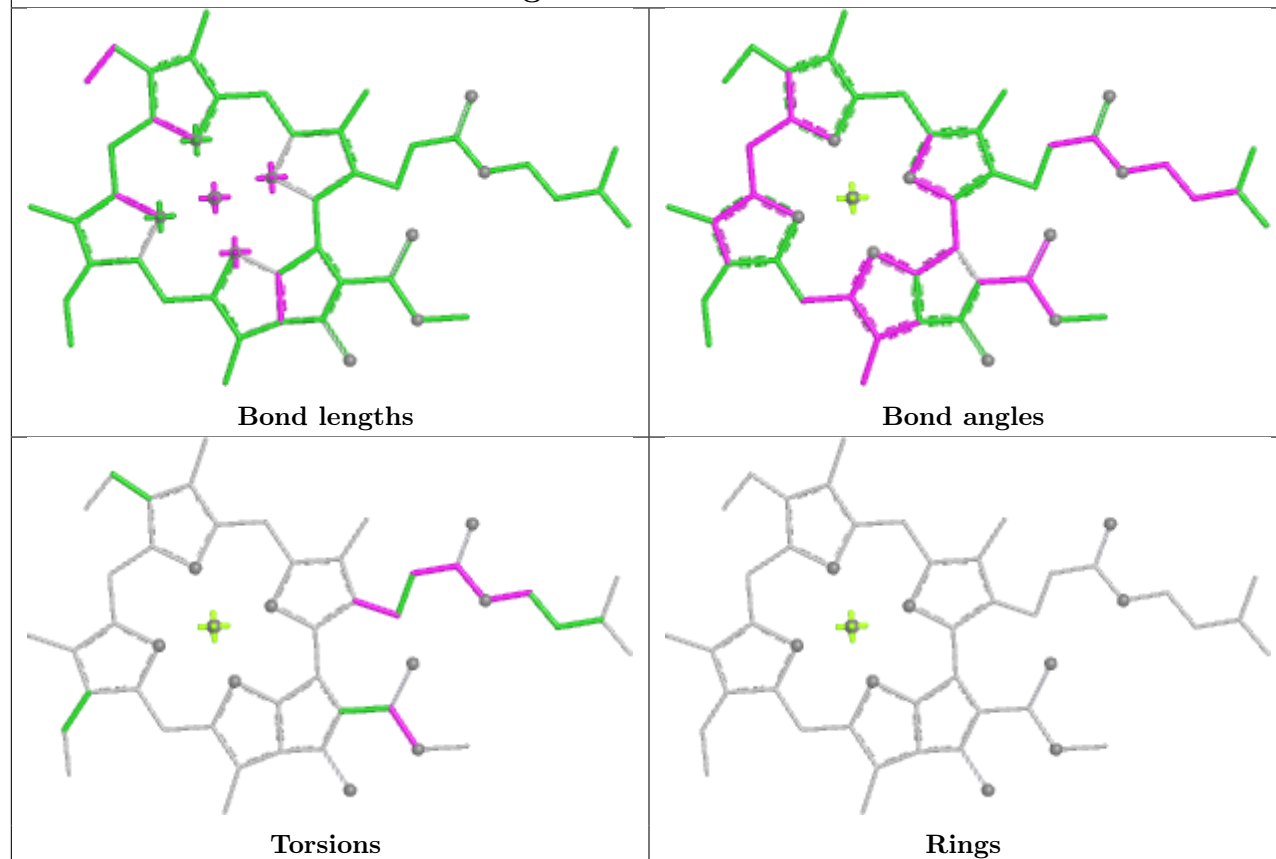
Ligand CLA 3 613



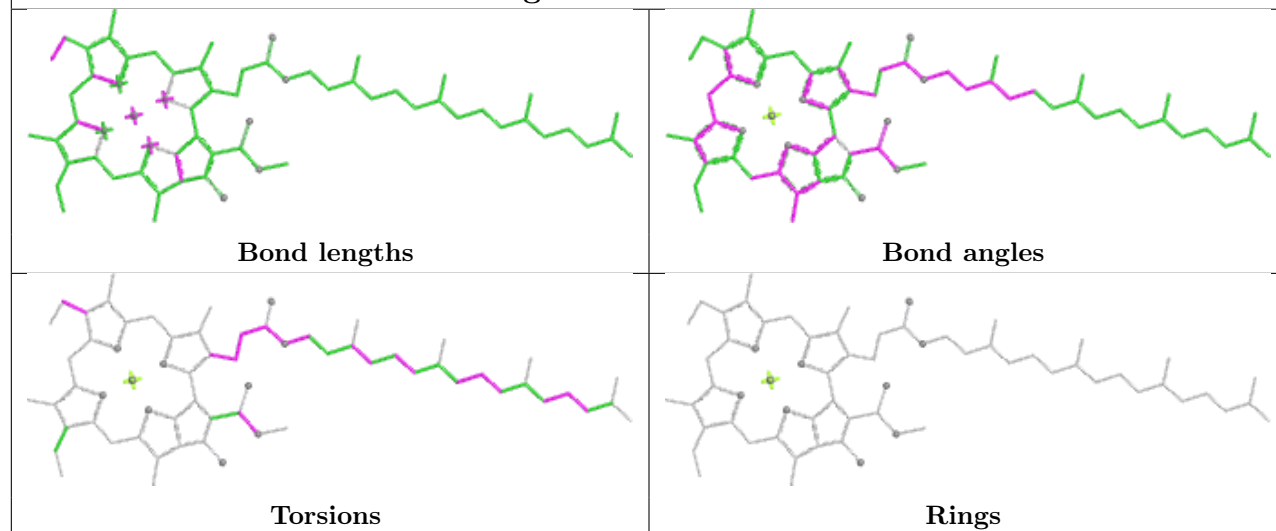
Ligand BCR F 4014

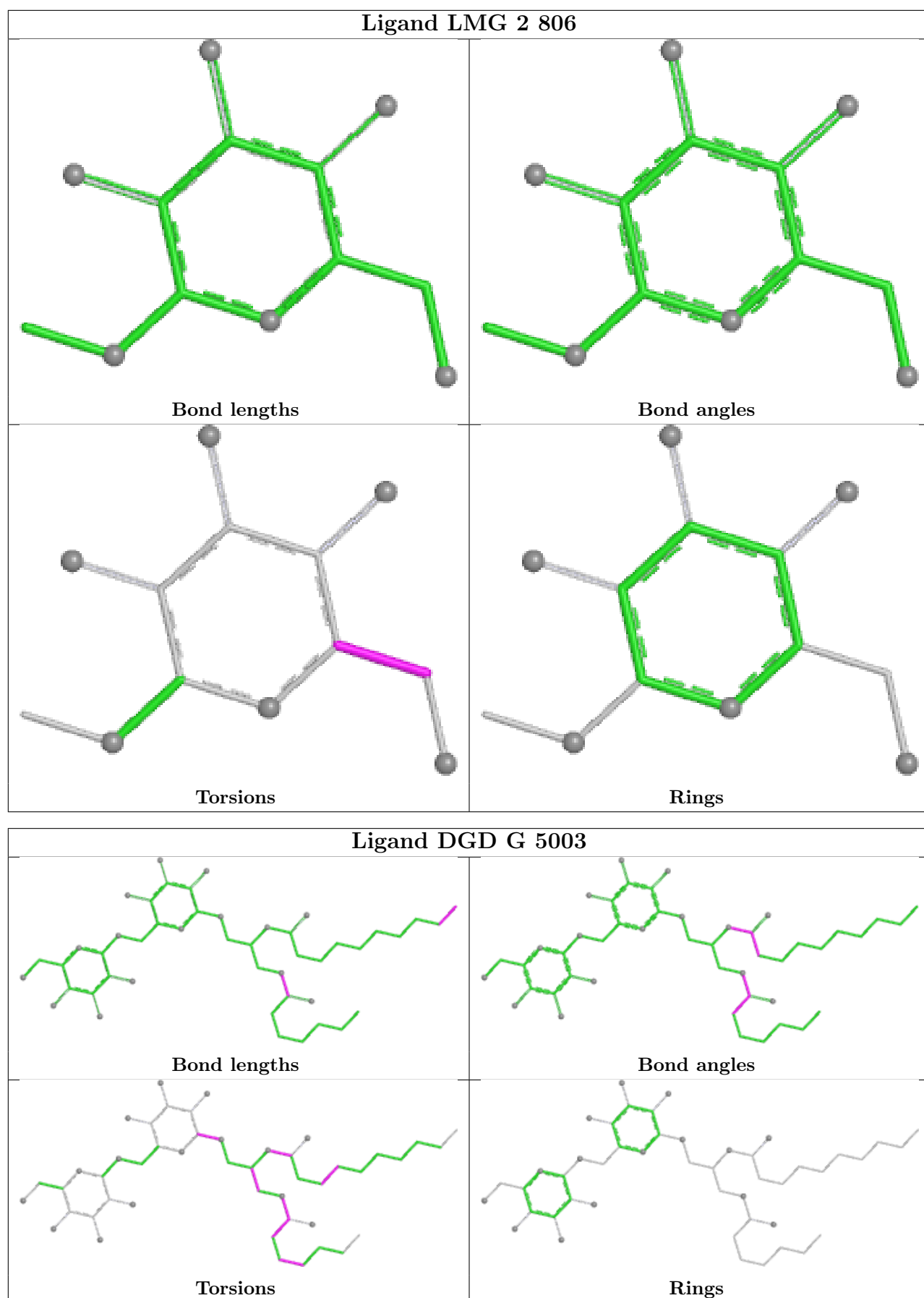


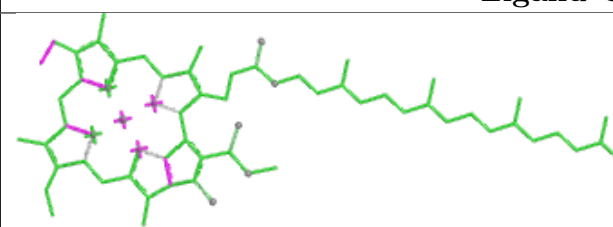
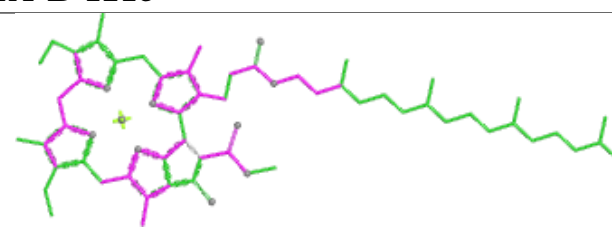
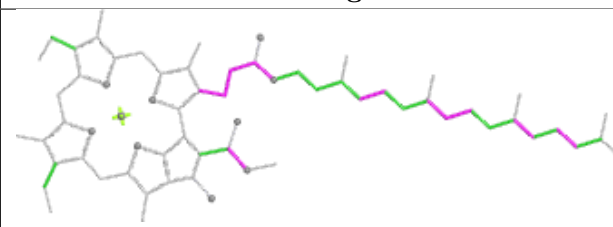
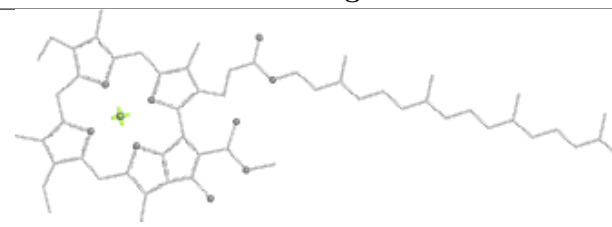
Ligand CLA B 1236

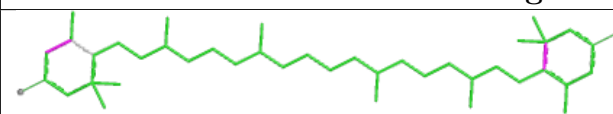
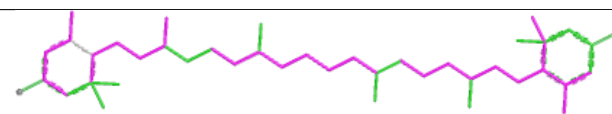
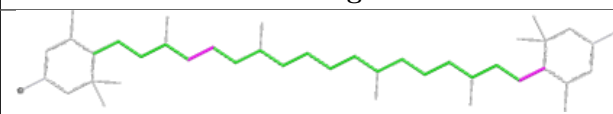
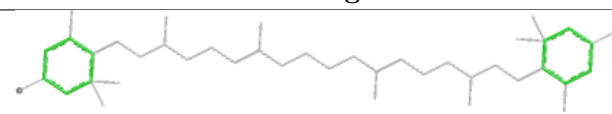


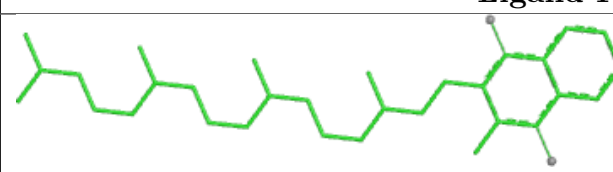
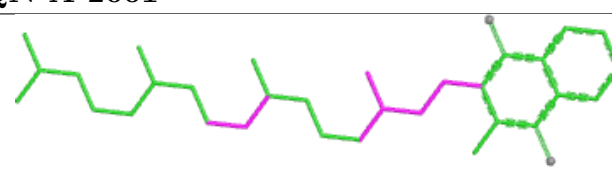
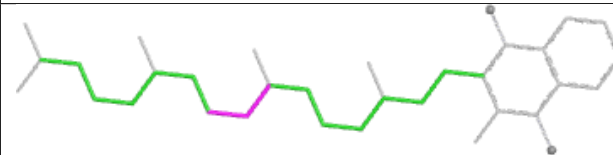
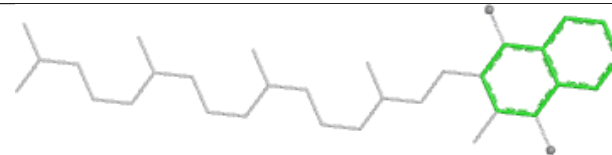
Ligand CLA B 1224

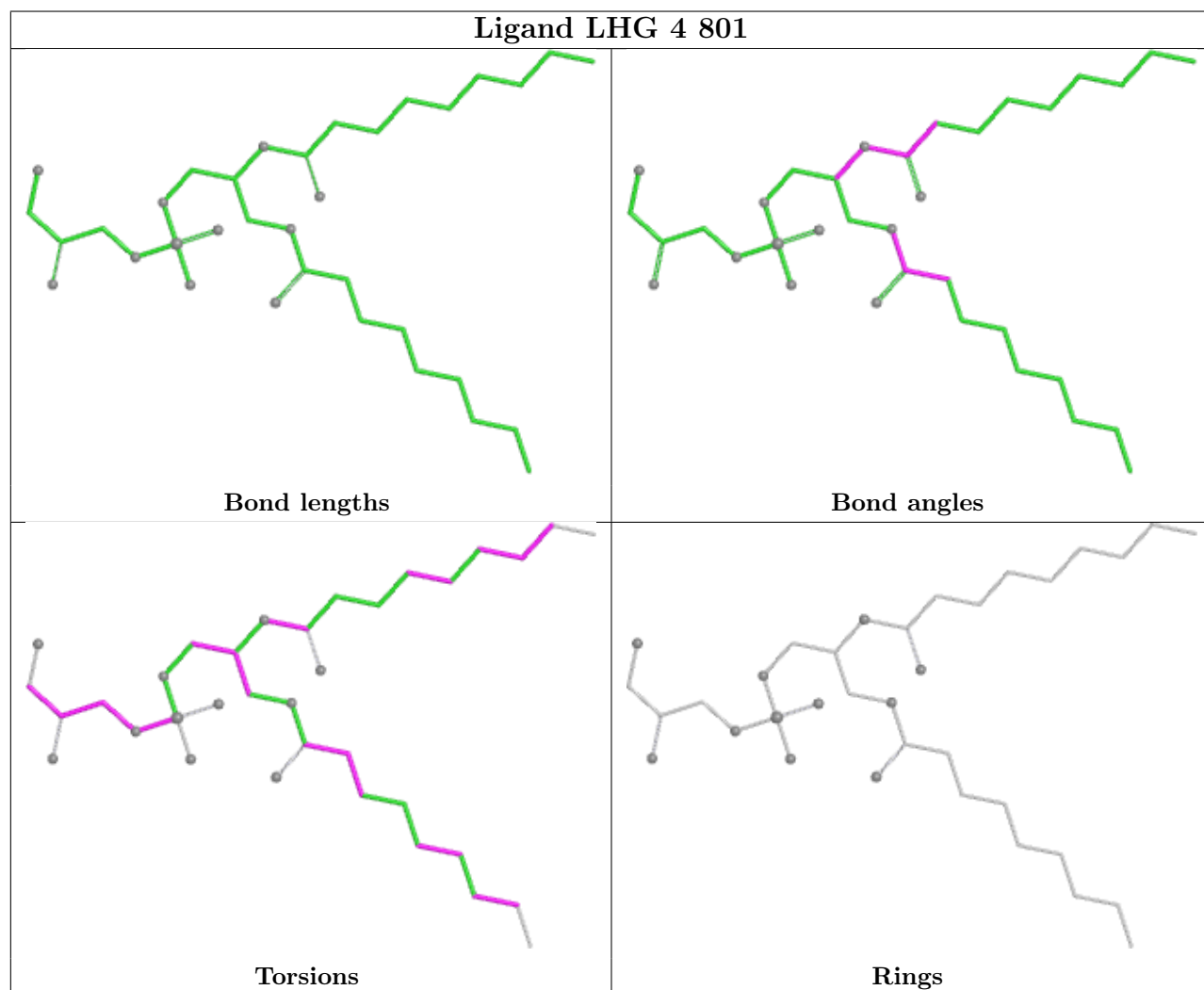
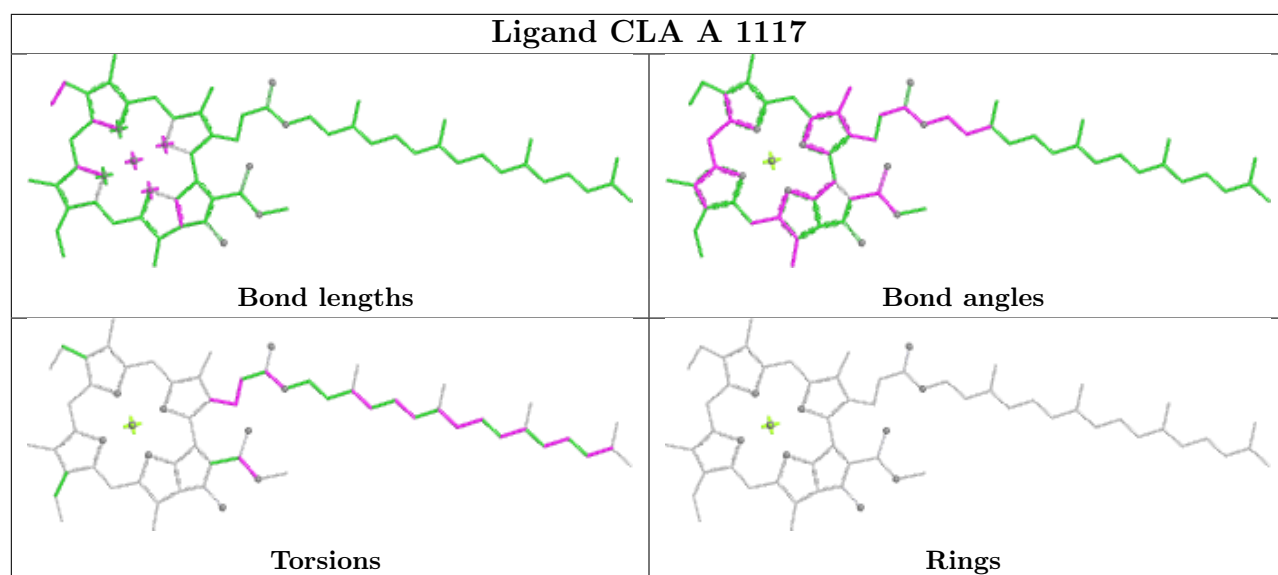


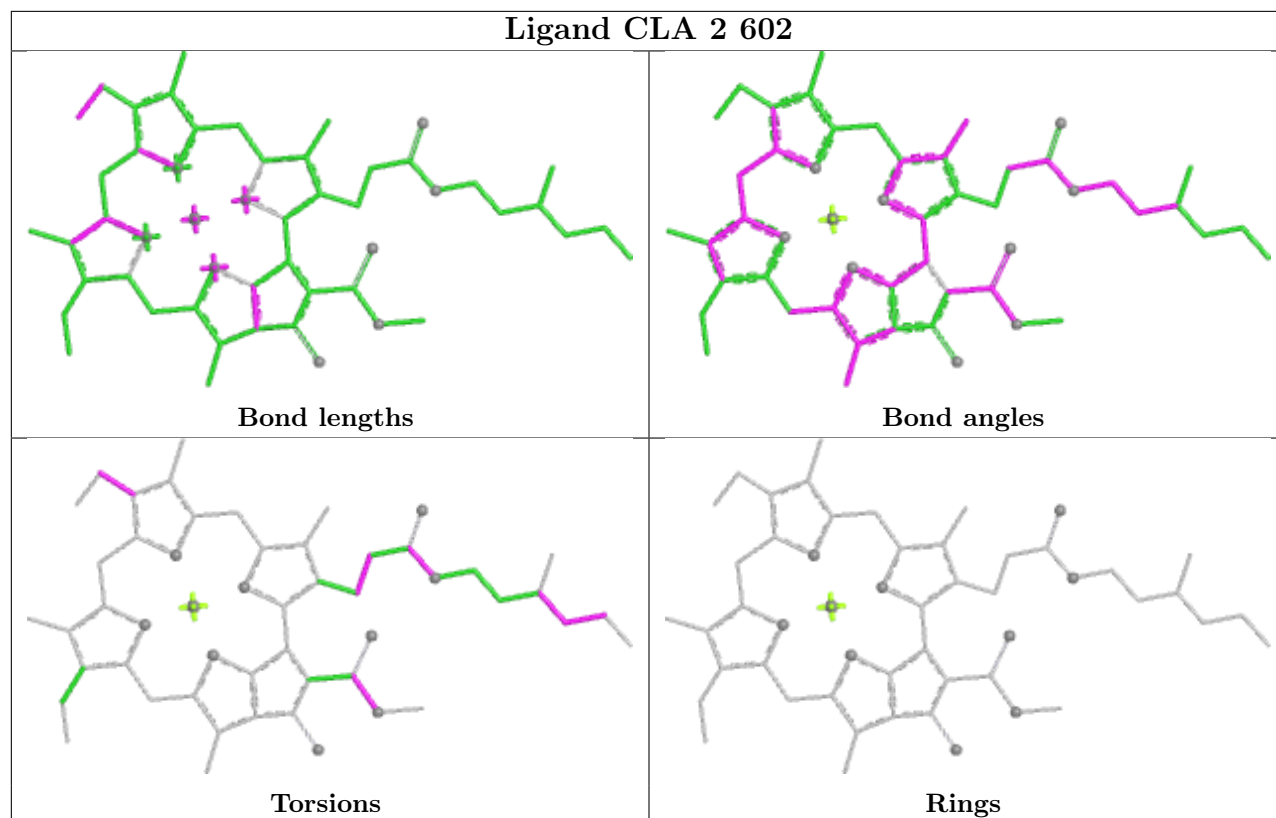
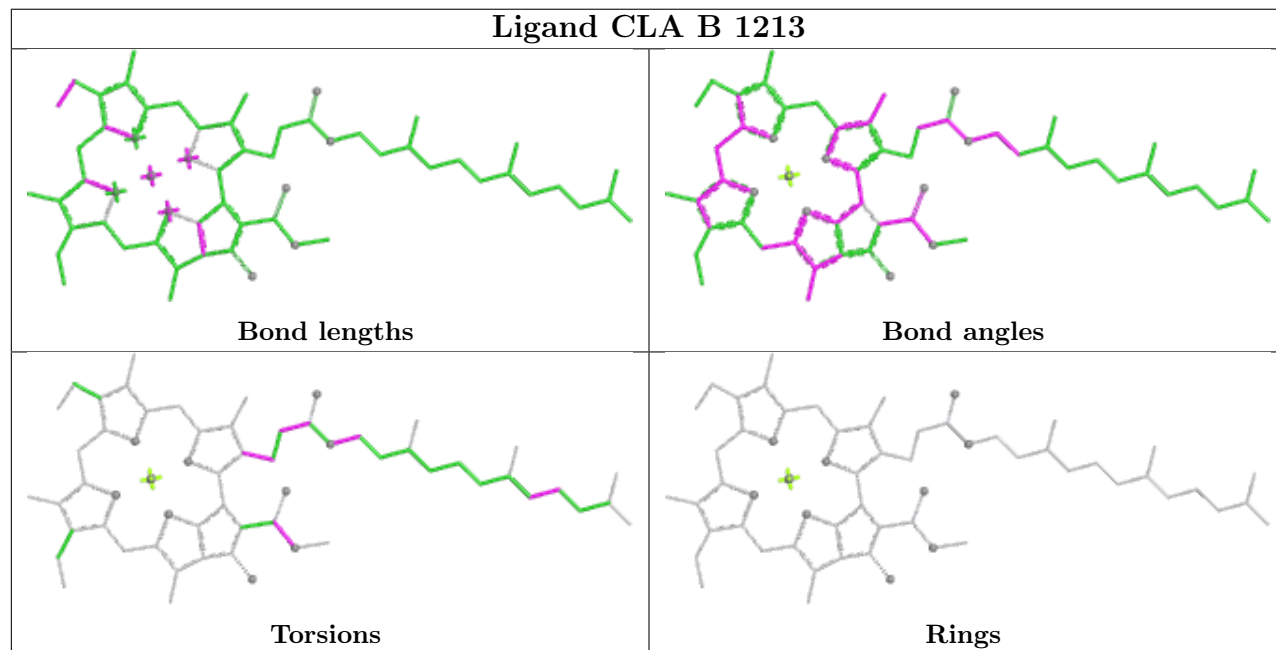


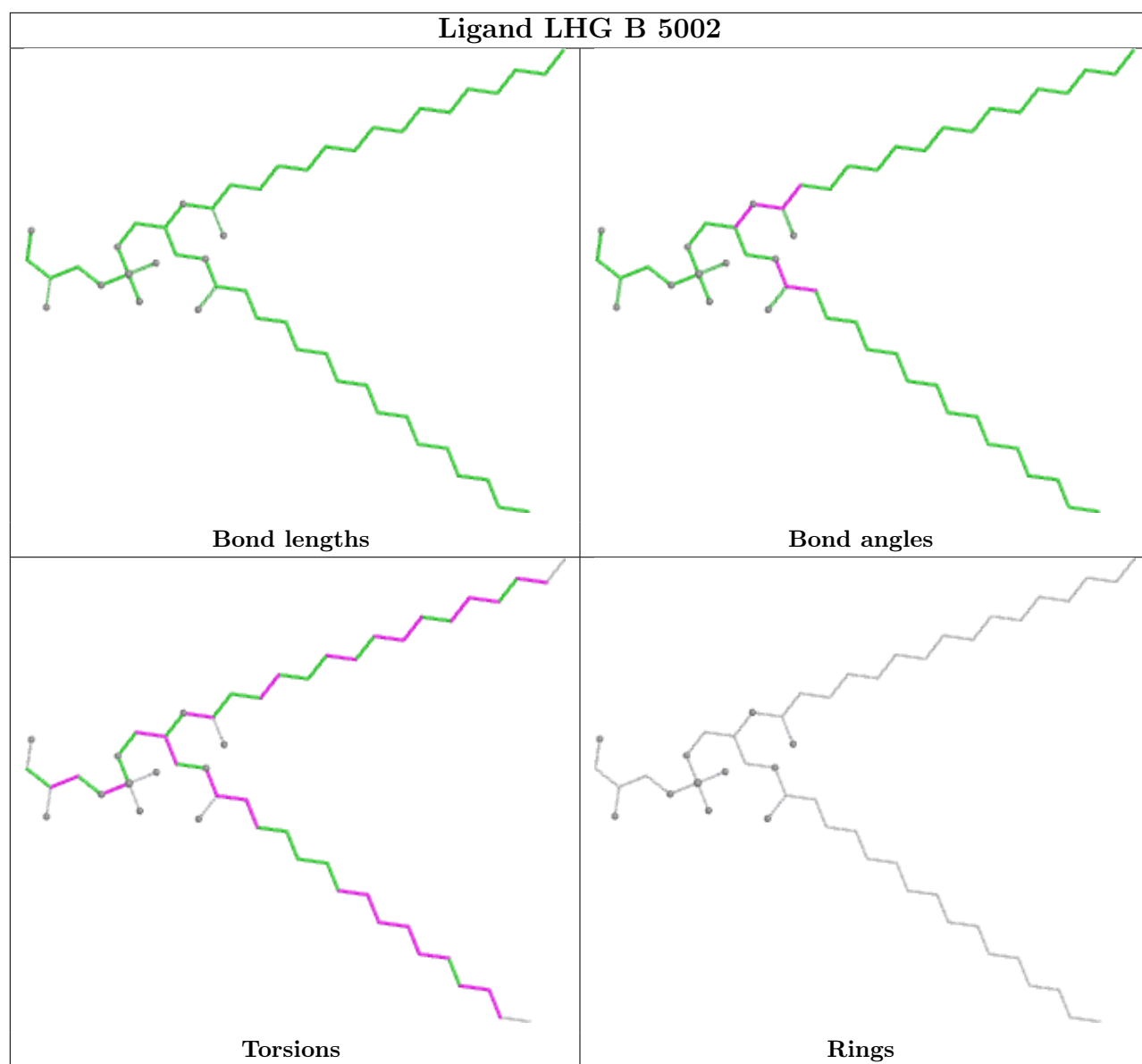
Ligand CLA B 1215	
	
Bond lengths	Bond angles
	
Torsions	Rings

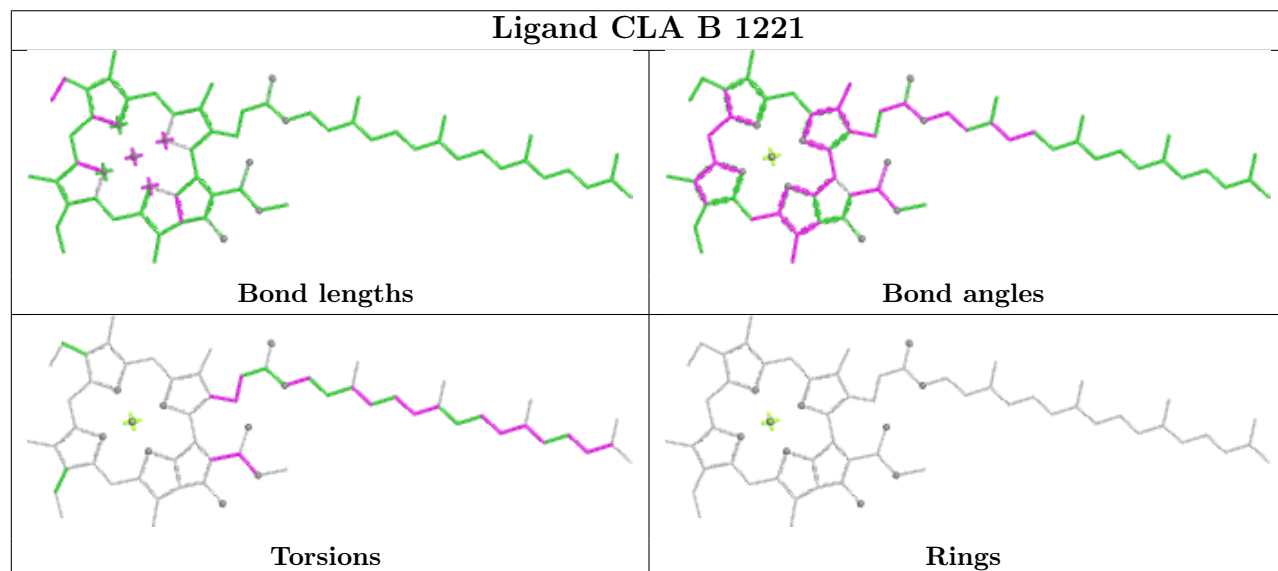
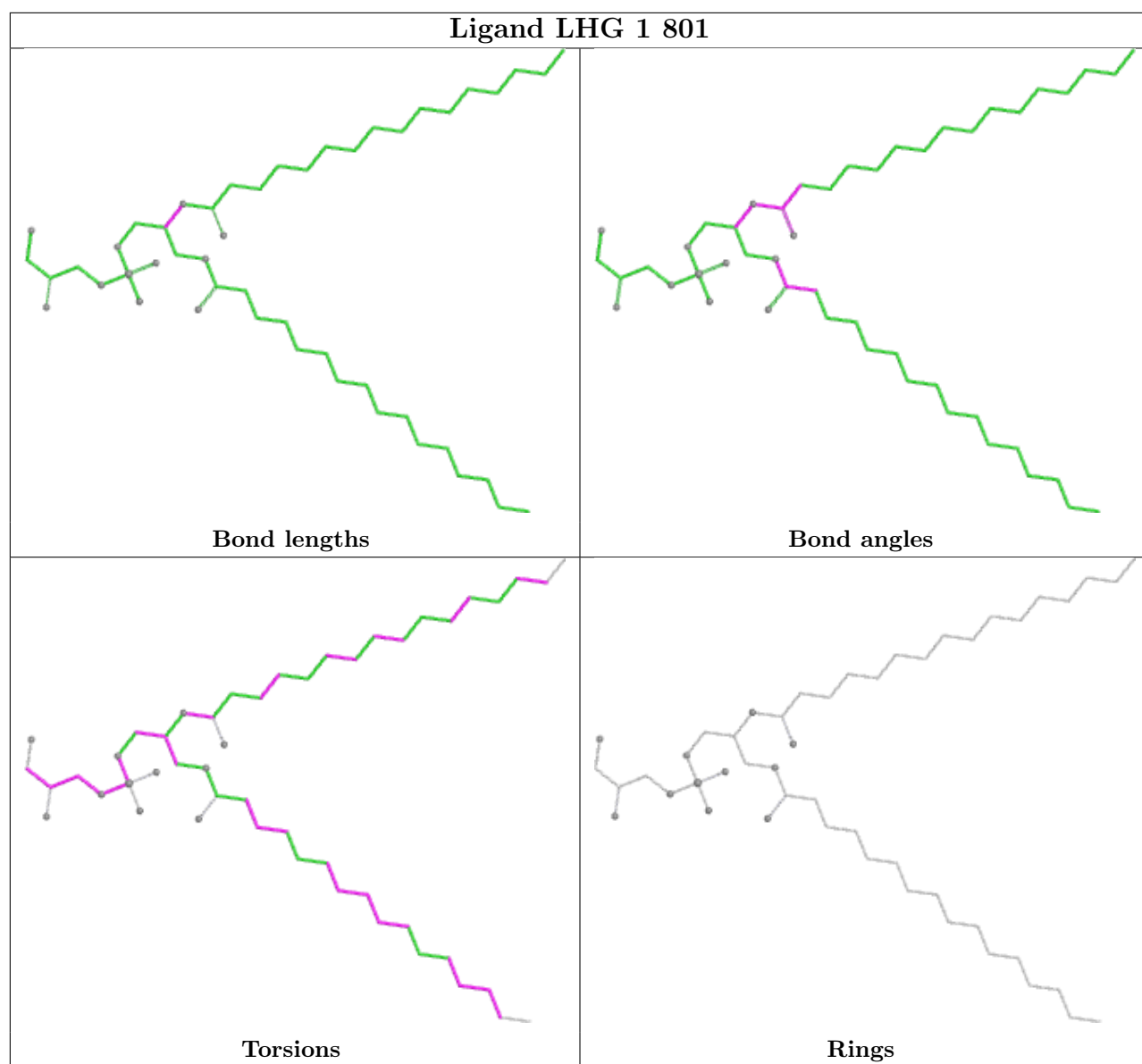
Ligand LUT 1 501	
	
Bond lengths	Bond angles
	
Torsions	Rings

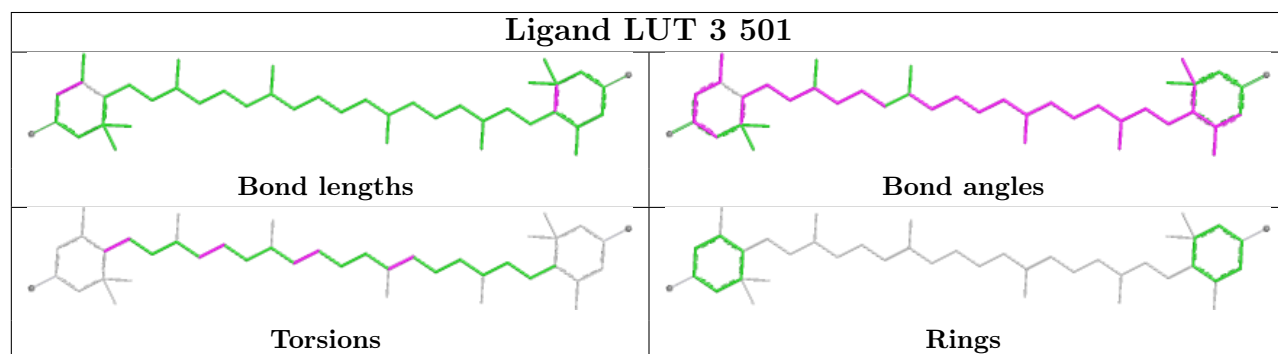
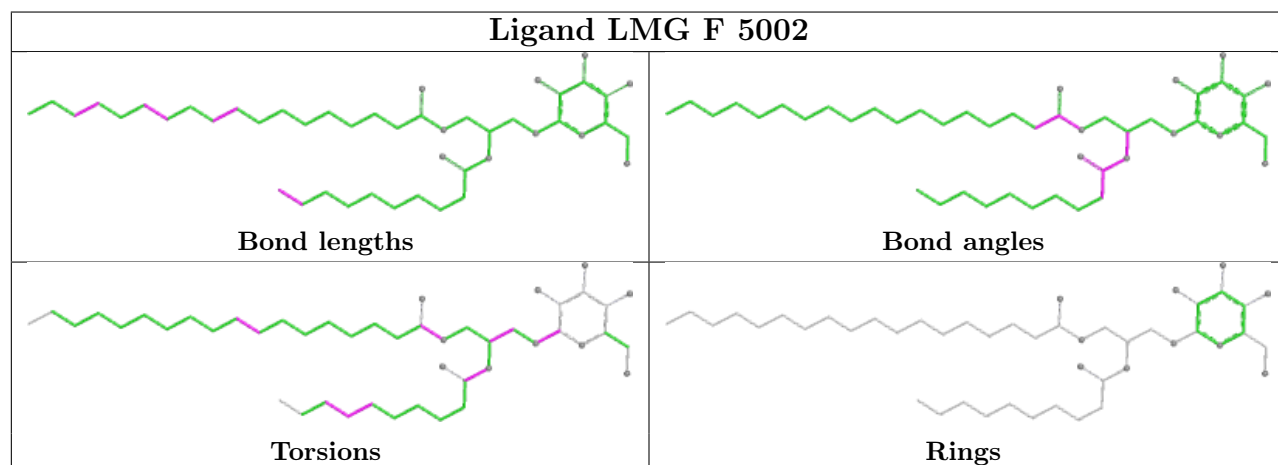
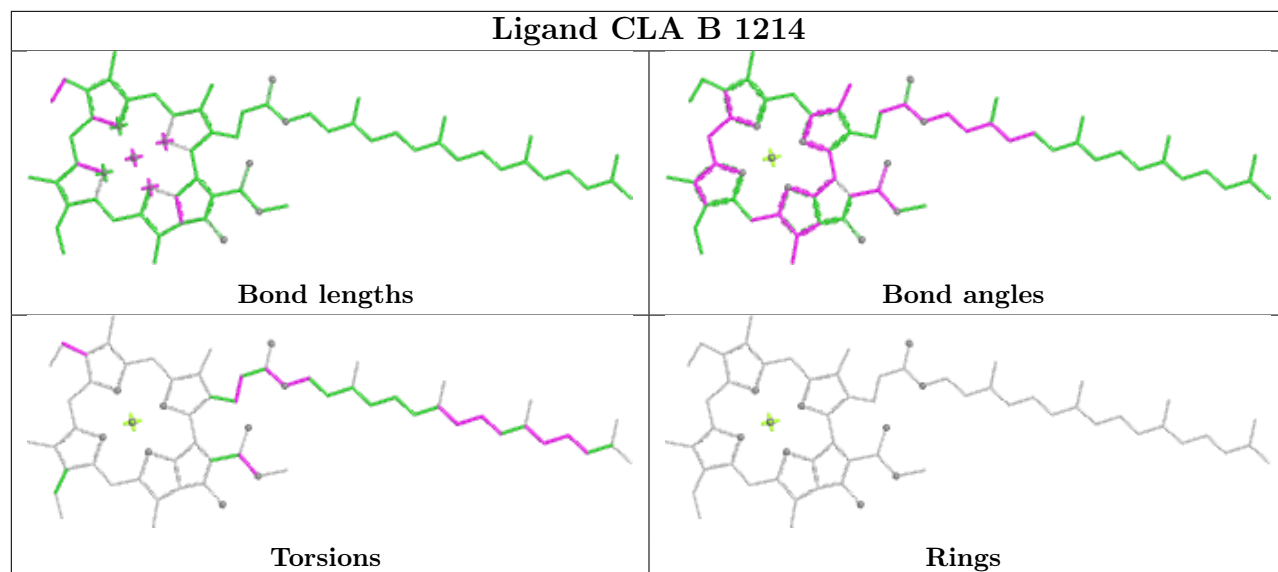
Ligand PQN A 2001	
	
Bond lengths	Bond angles
	
Torsions	Rings

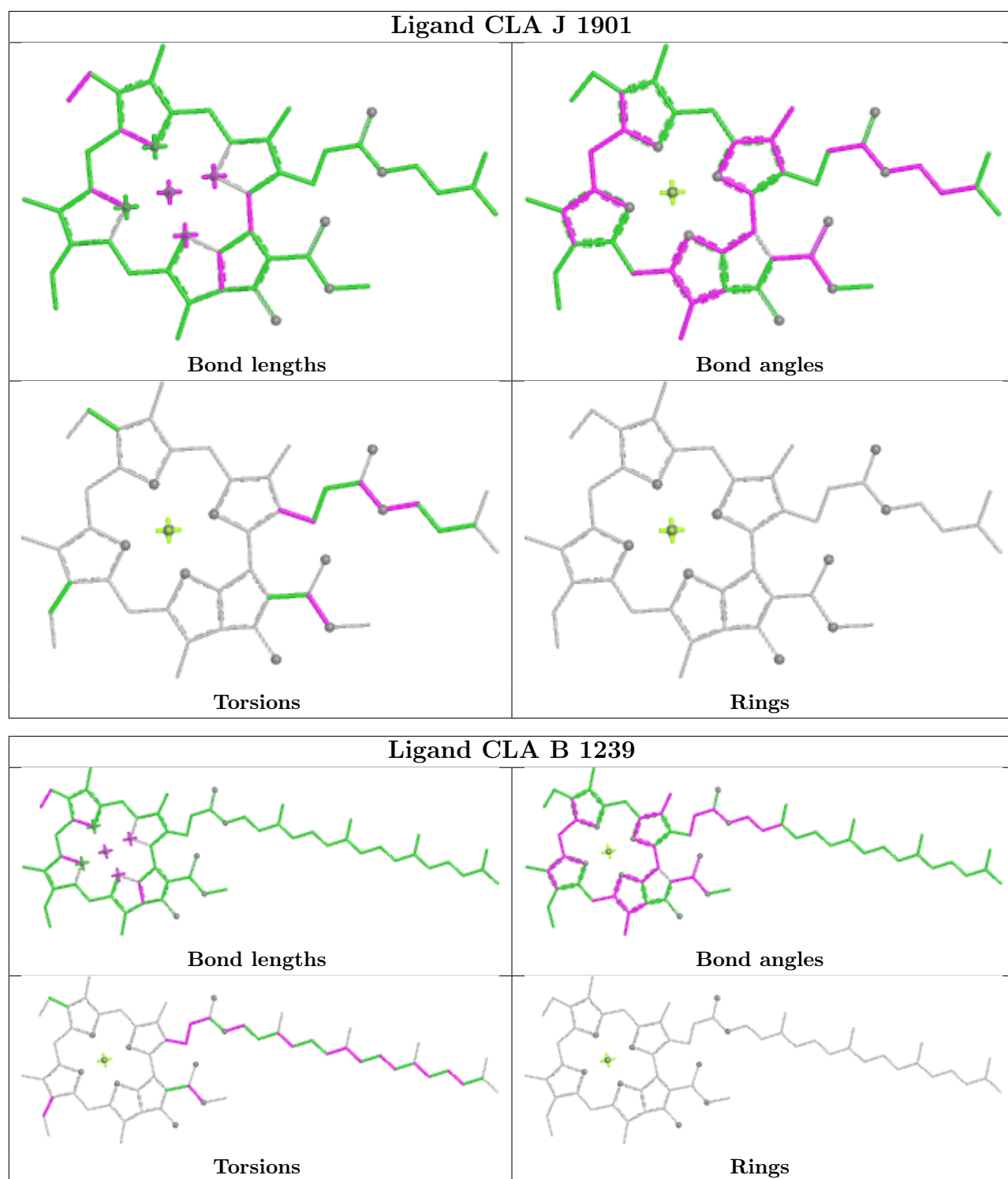


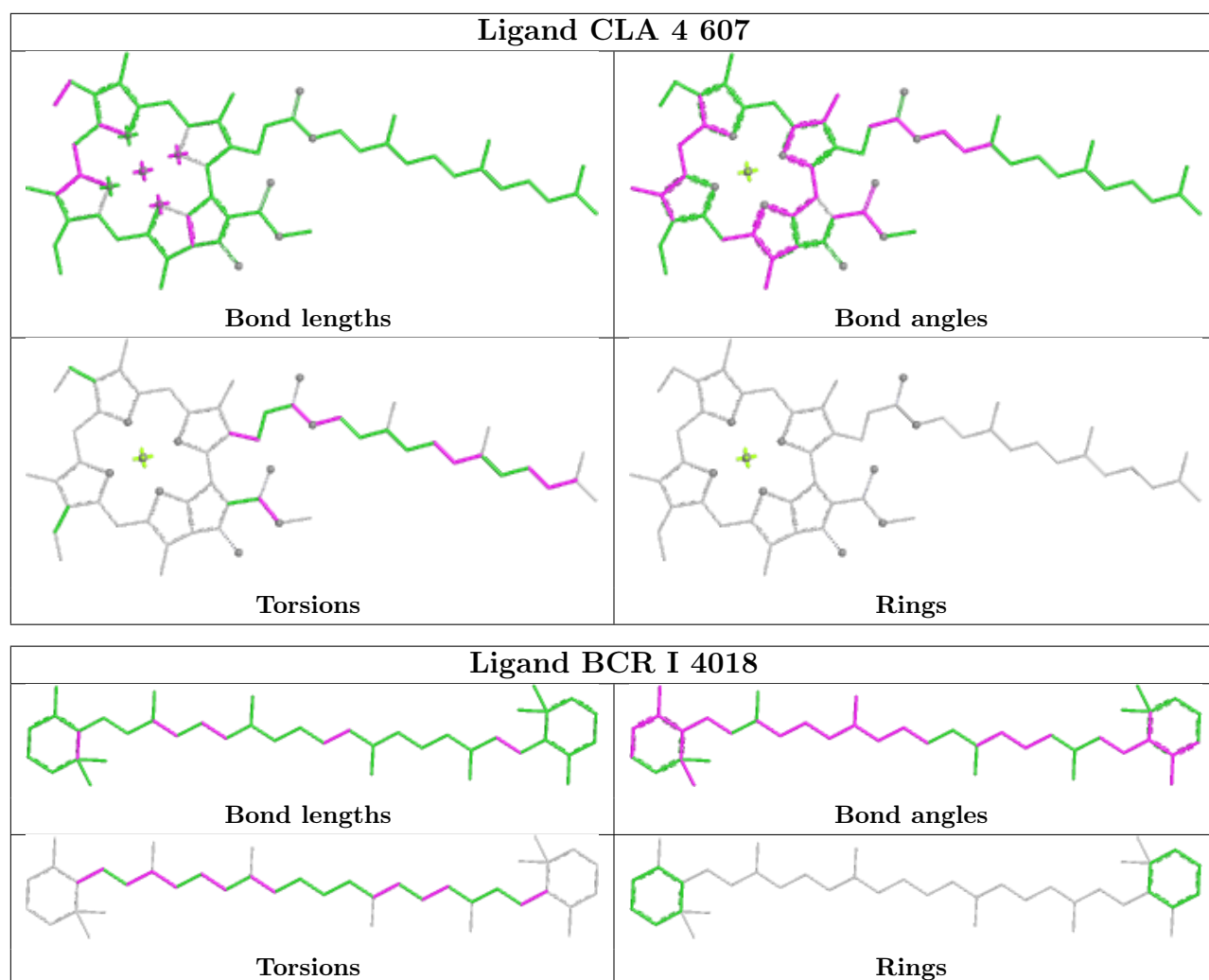
Ligand CLA 2 602**Ligand CLA B 1213**

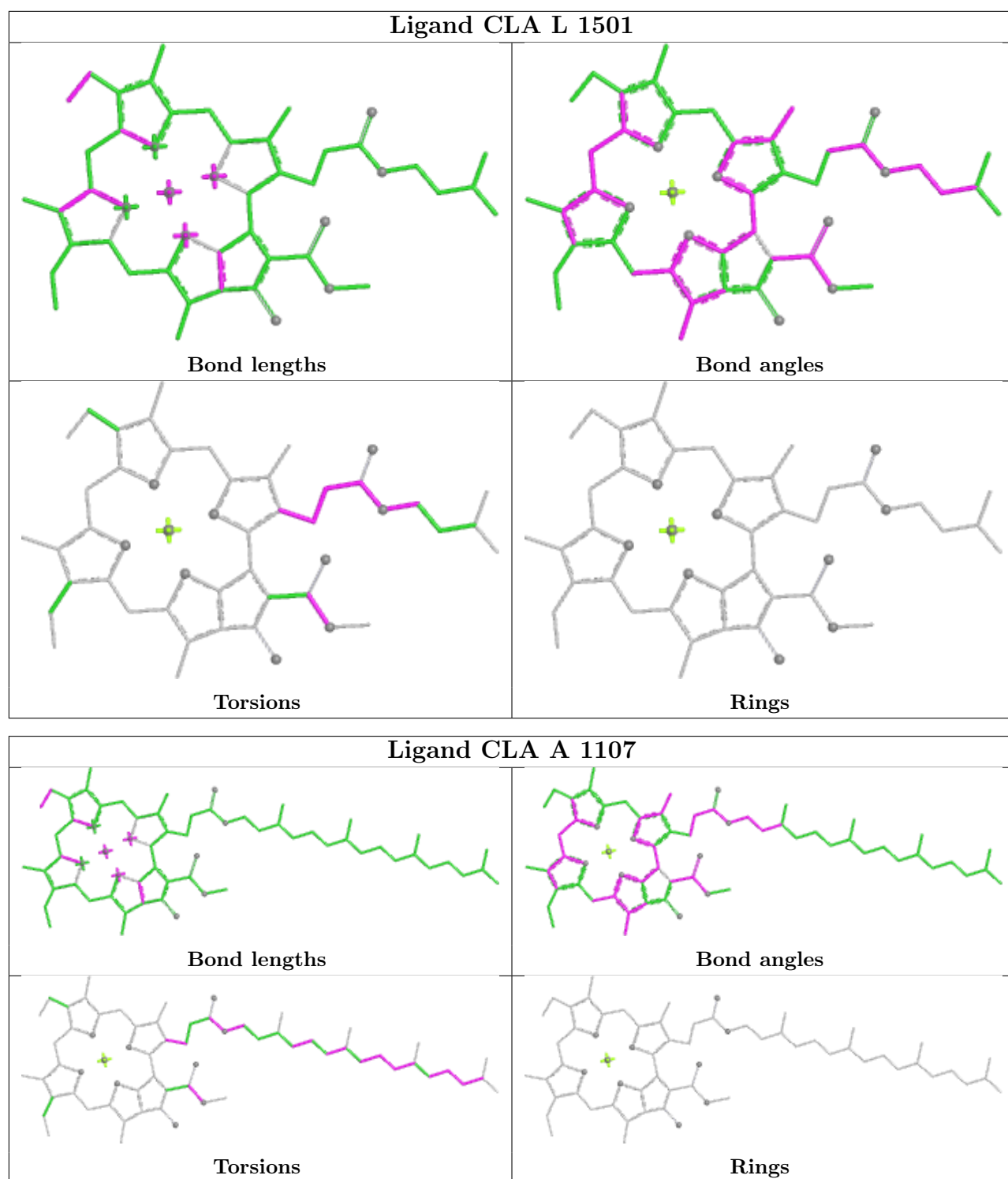


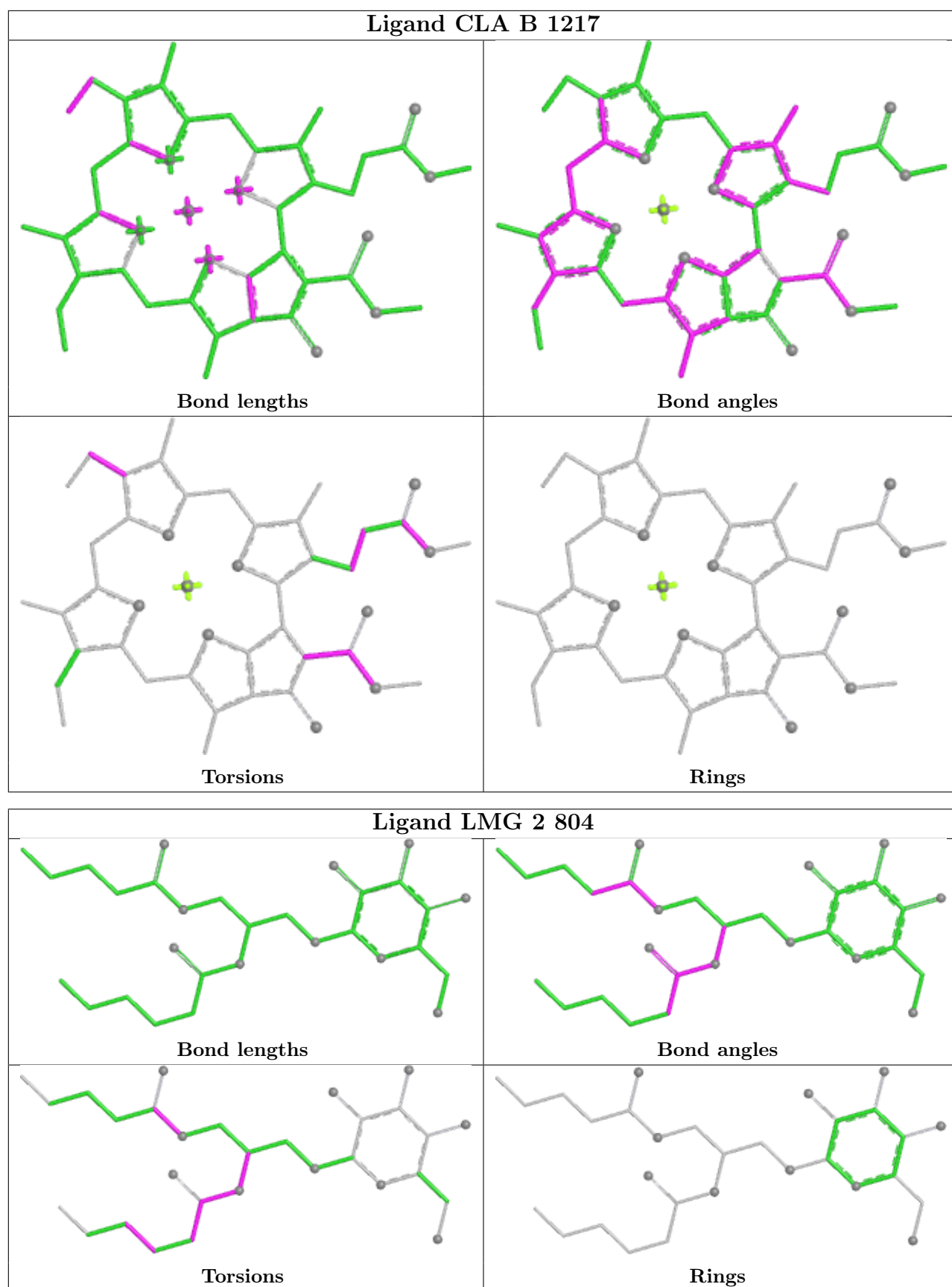


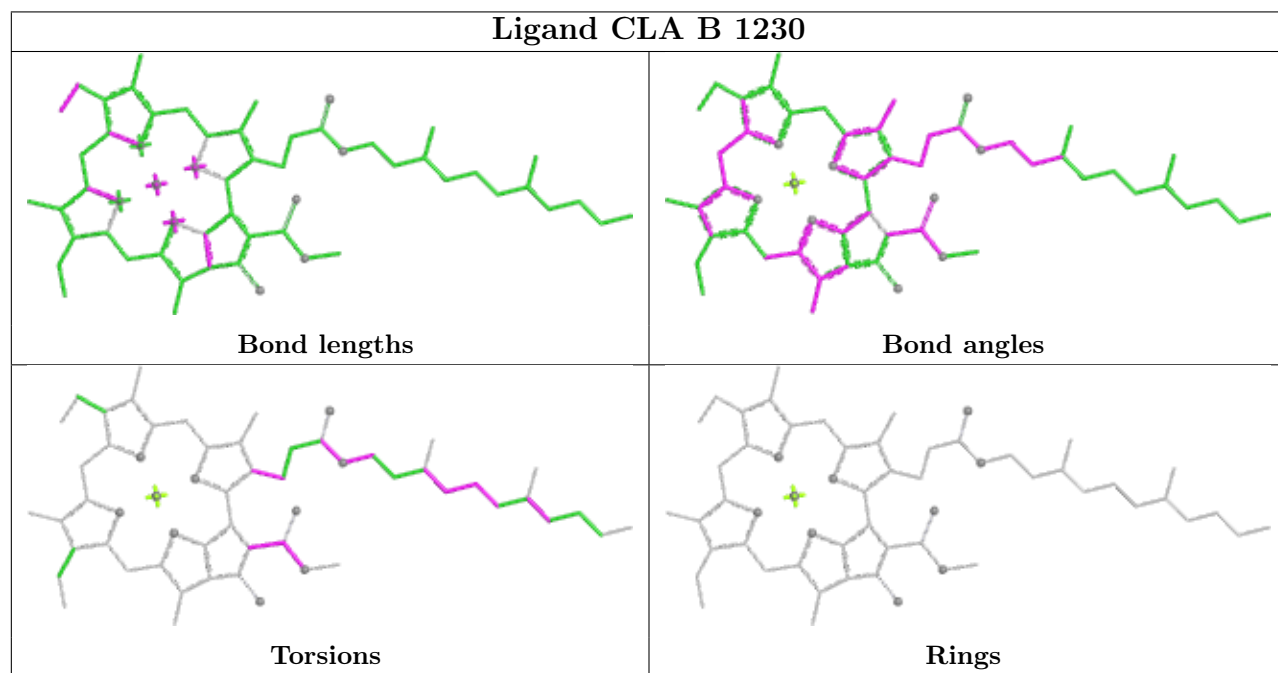
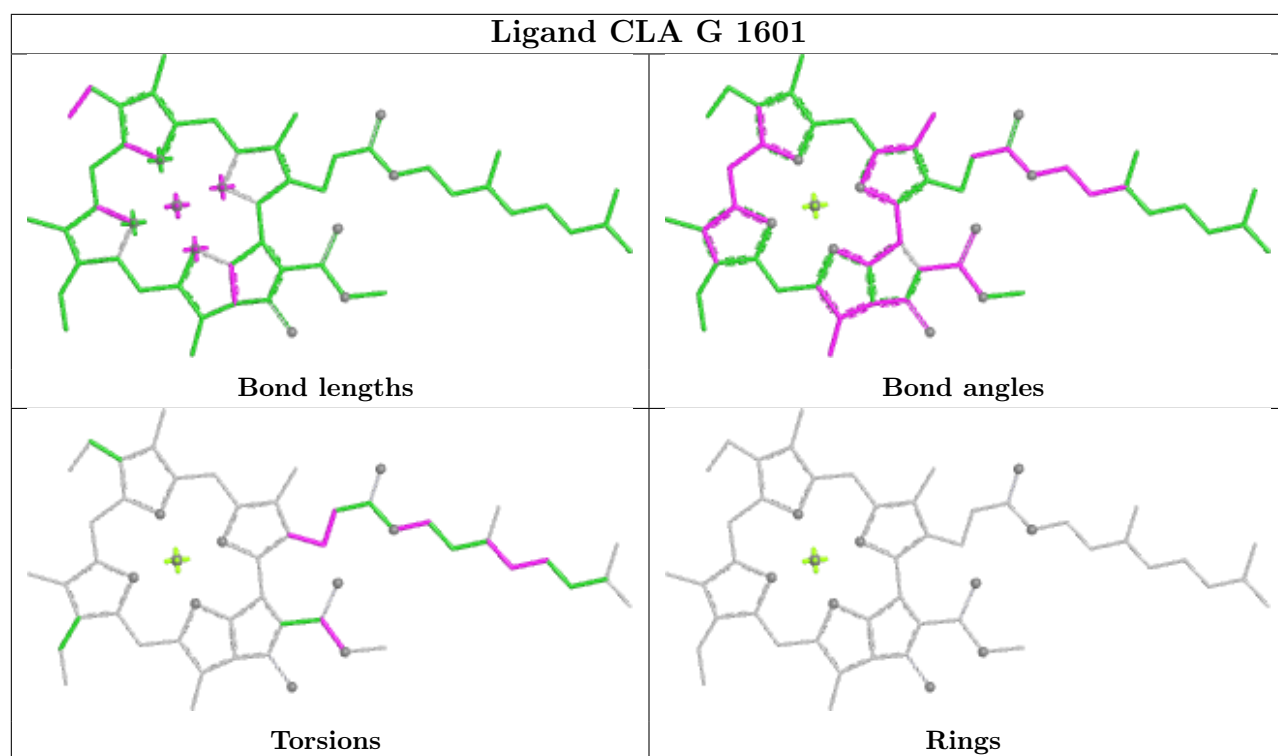


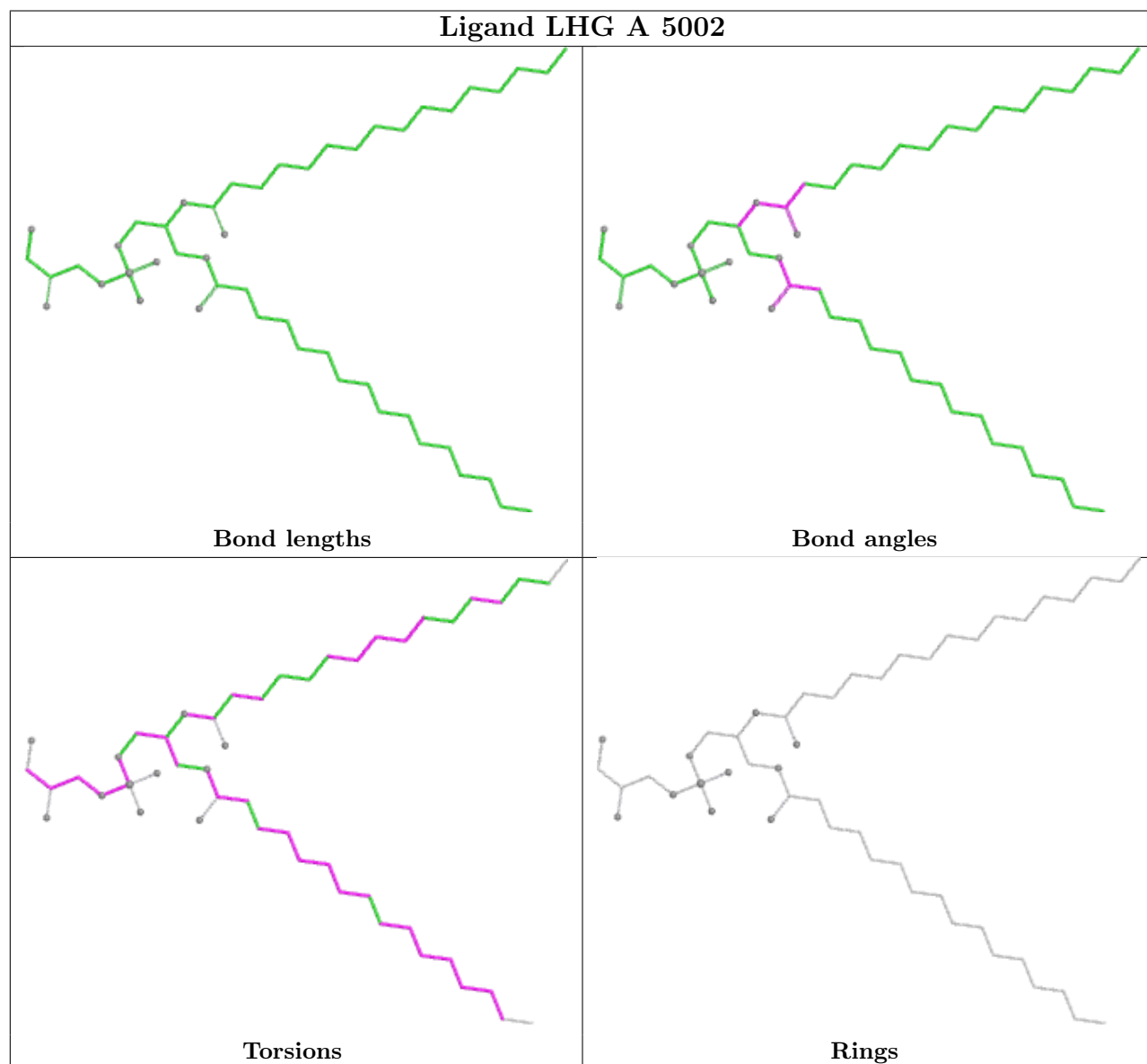
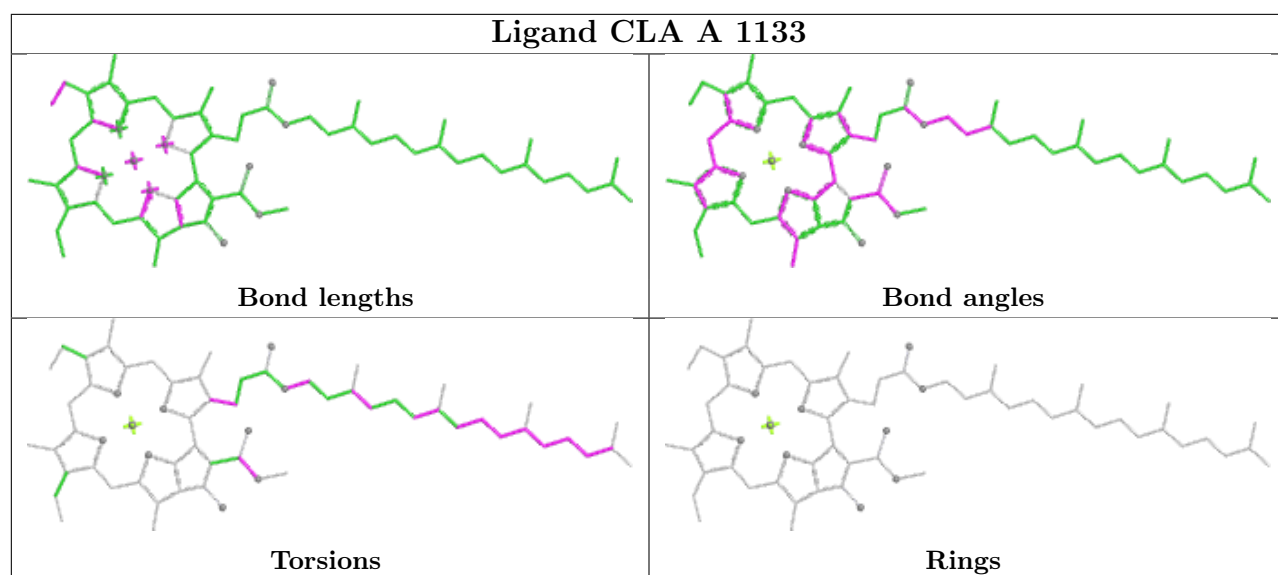


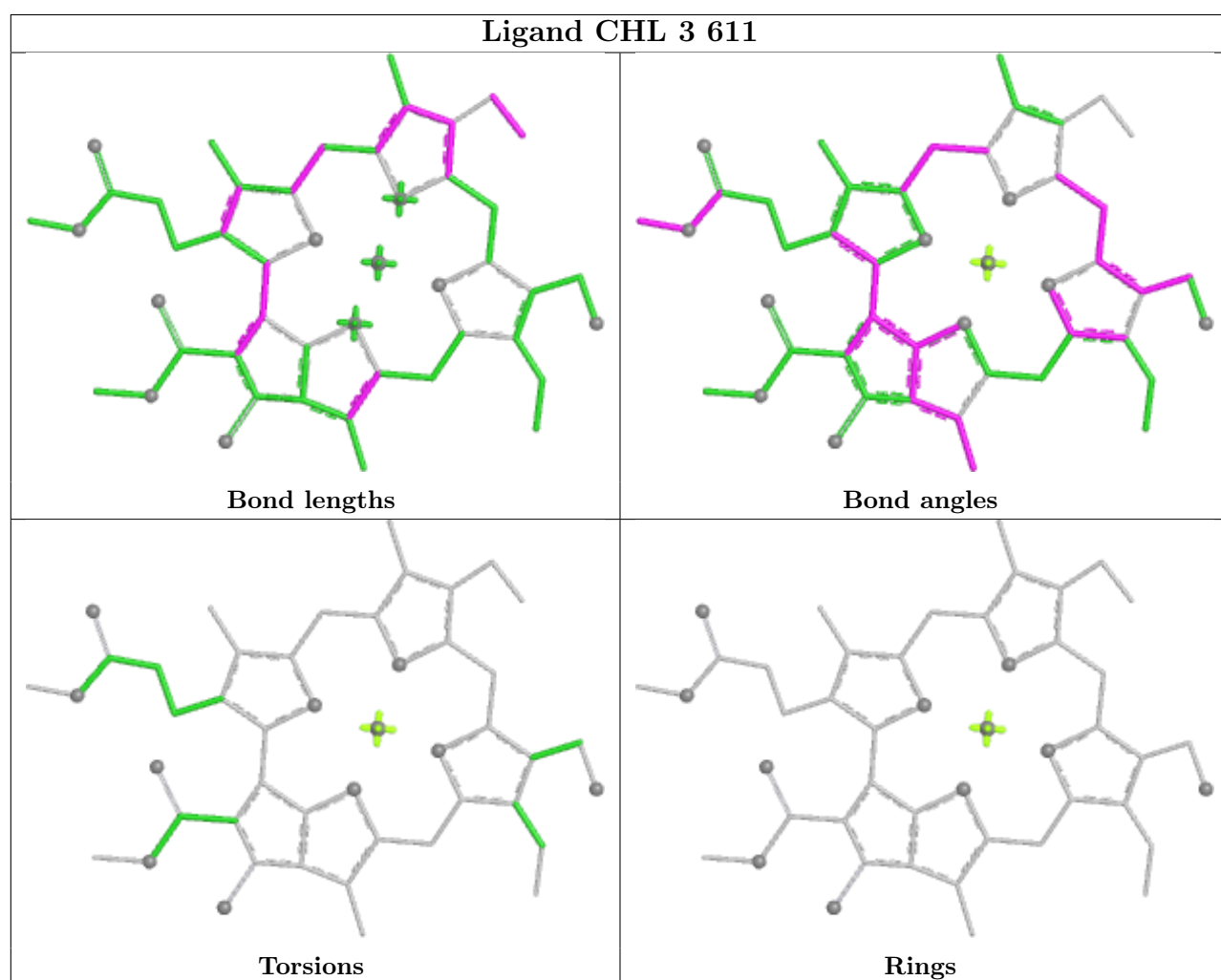
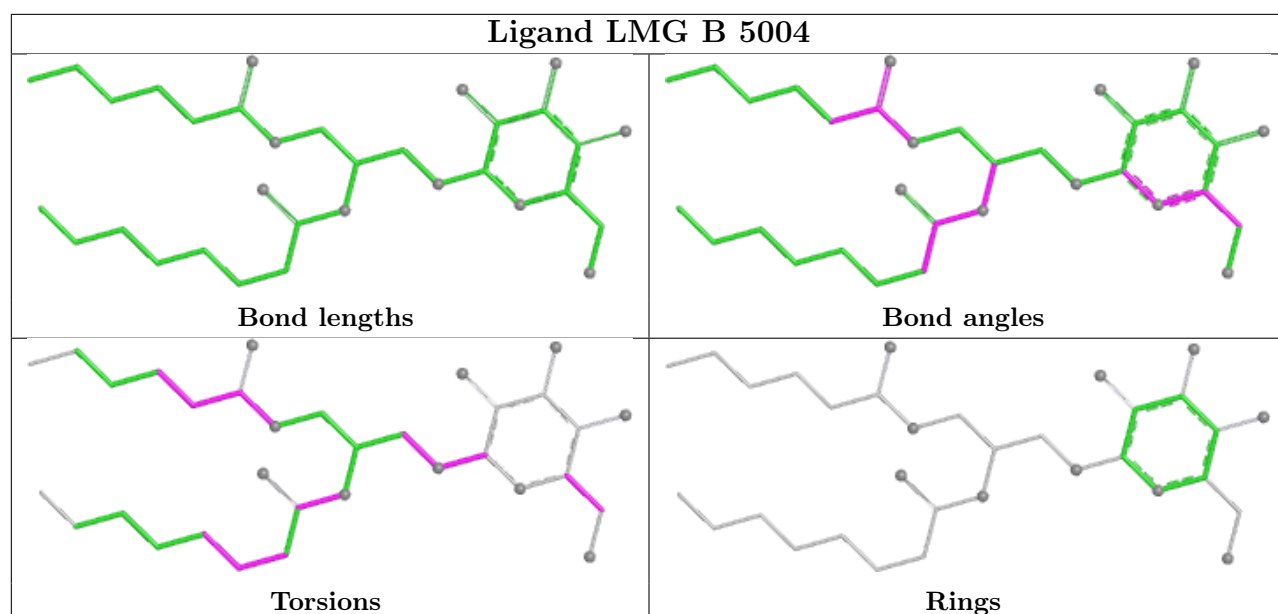


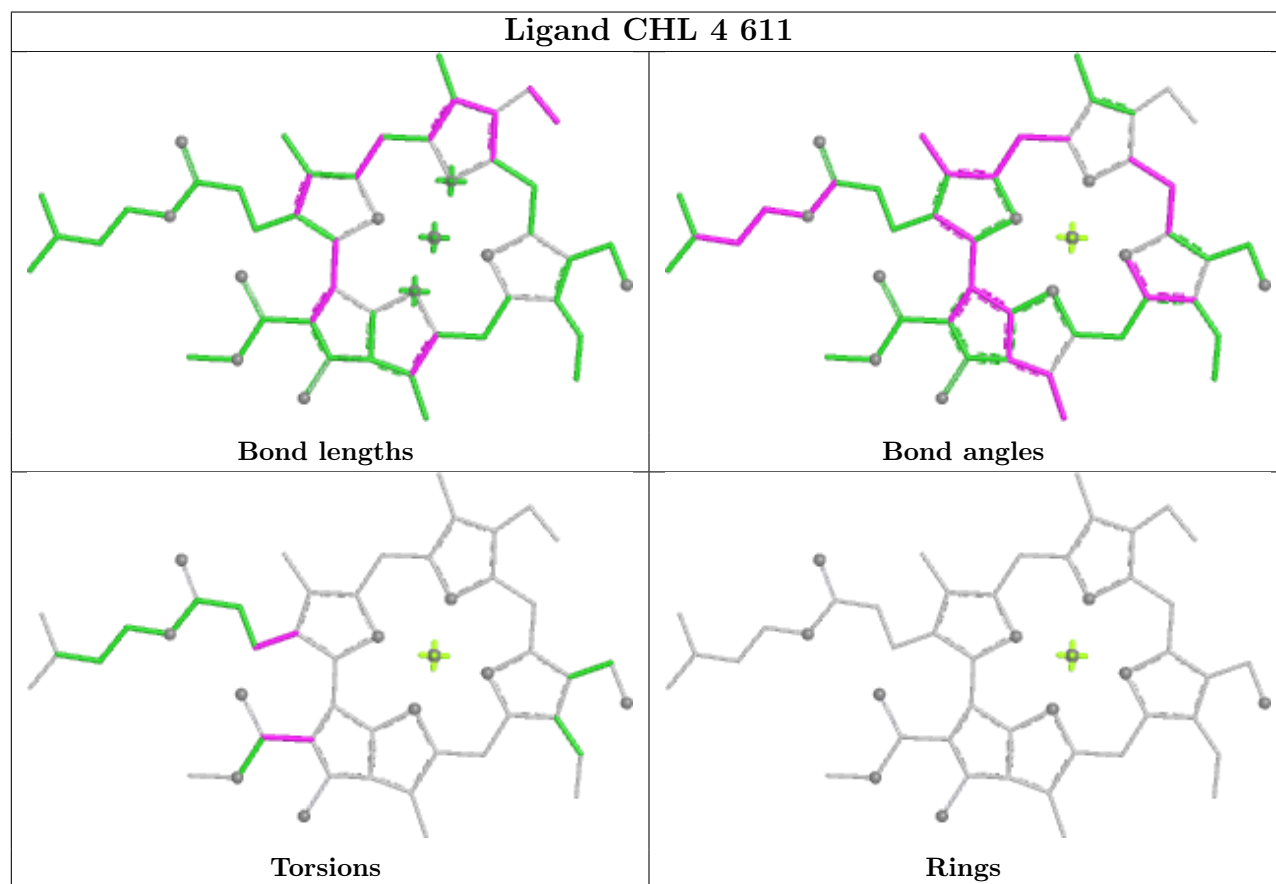
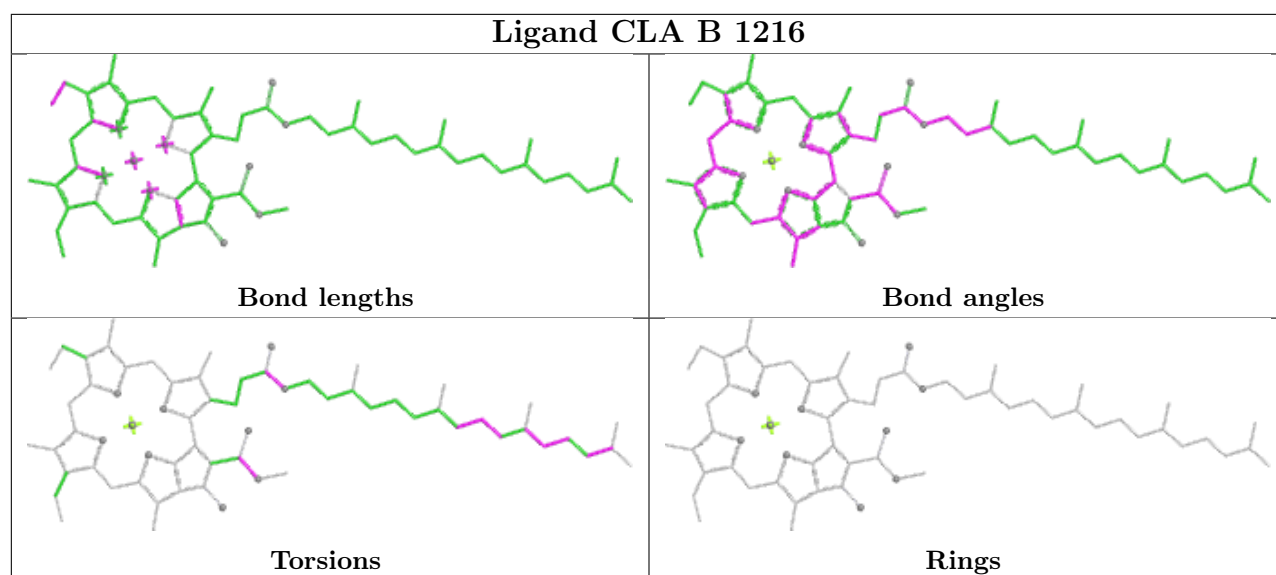


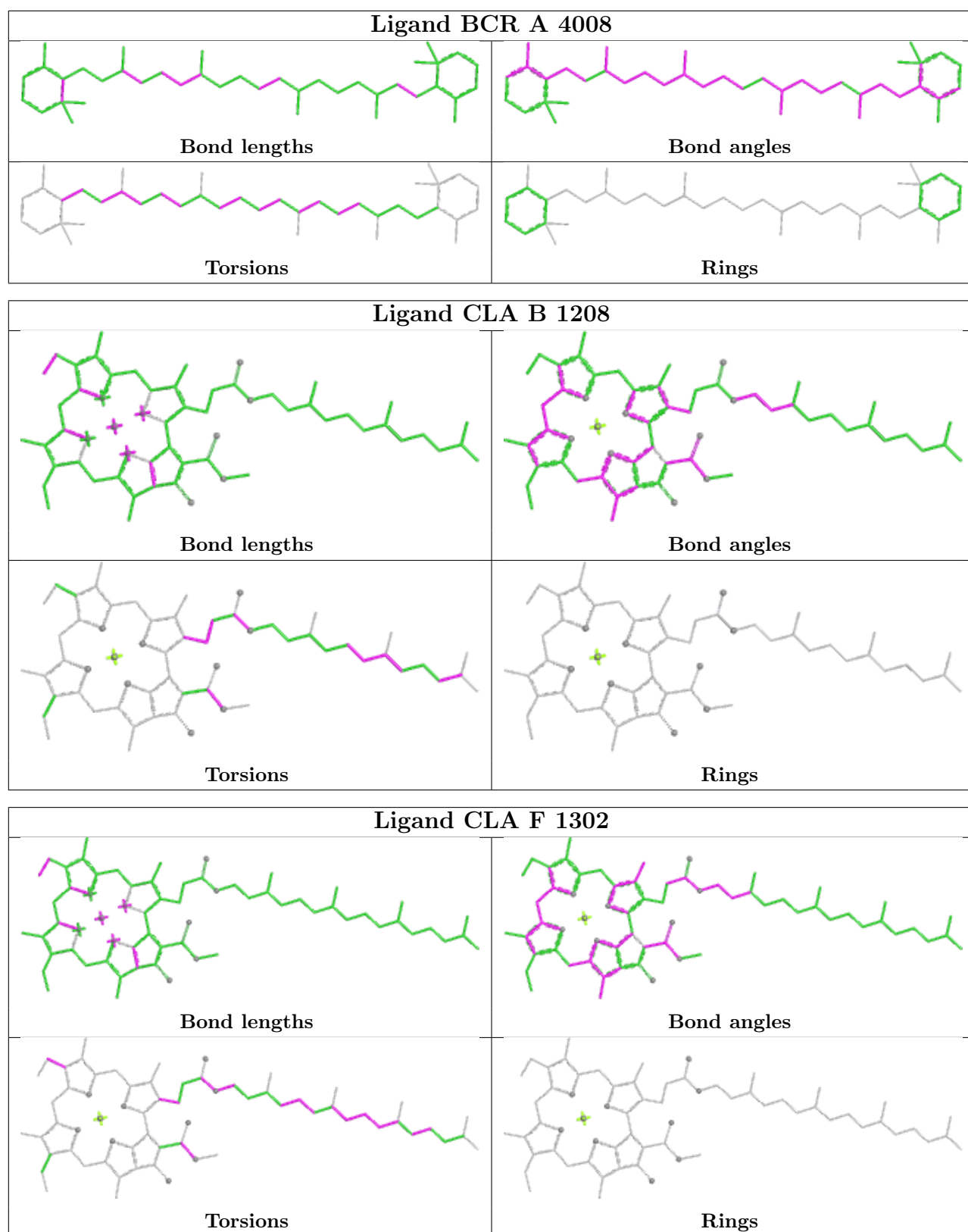


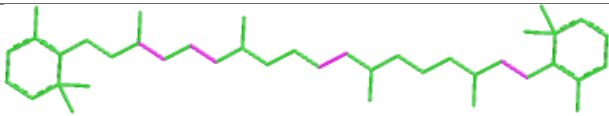
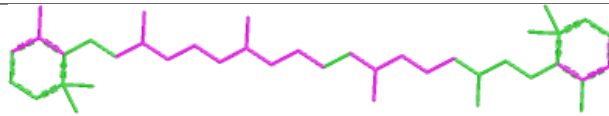
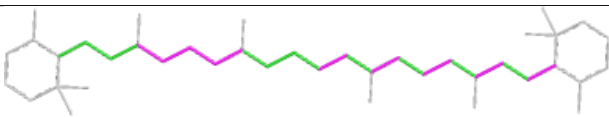
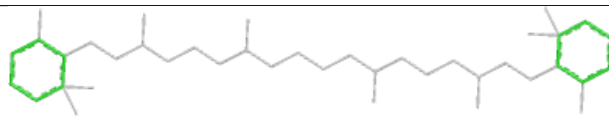


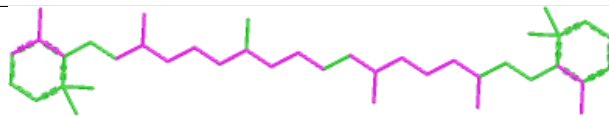
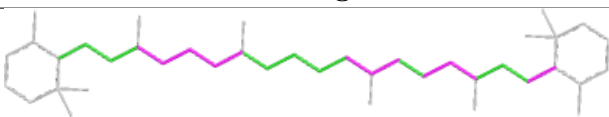
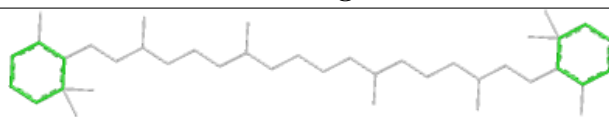


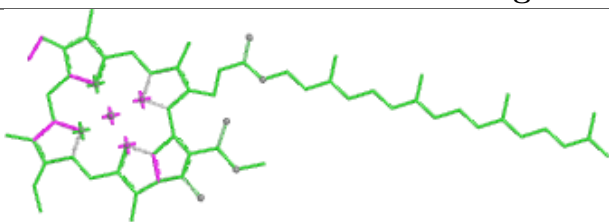
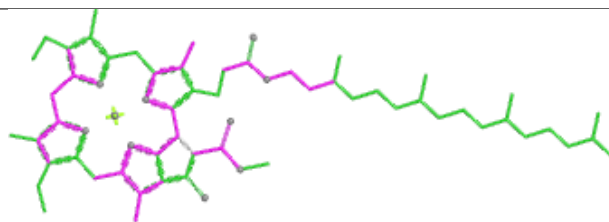
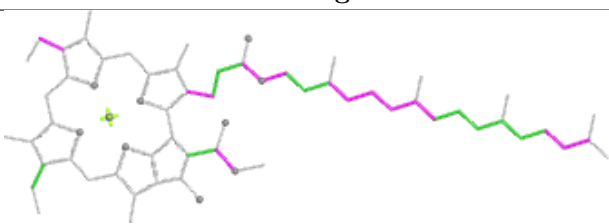
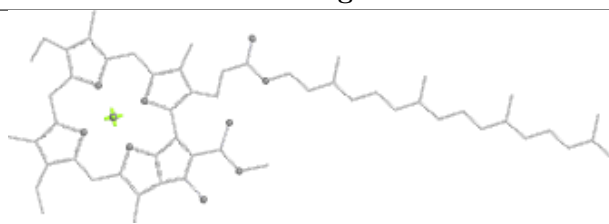


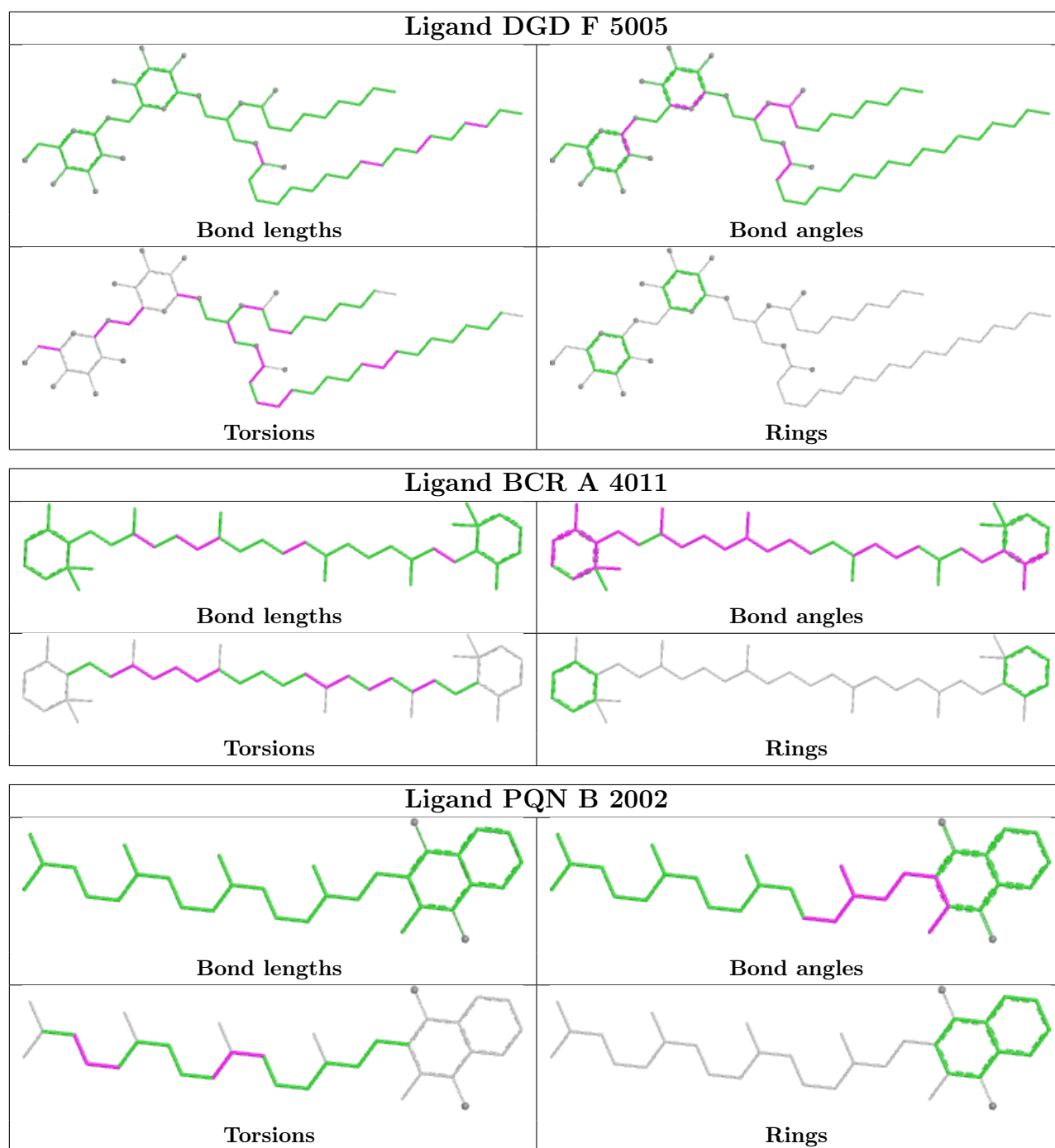


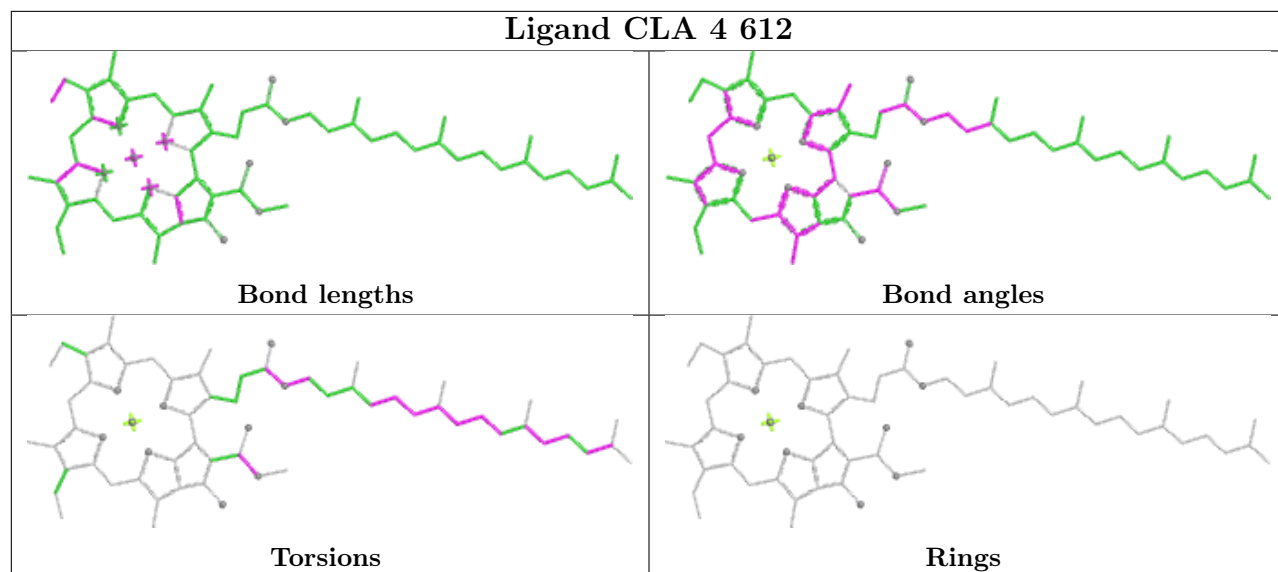
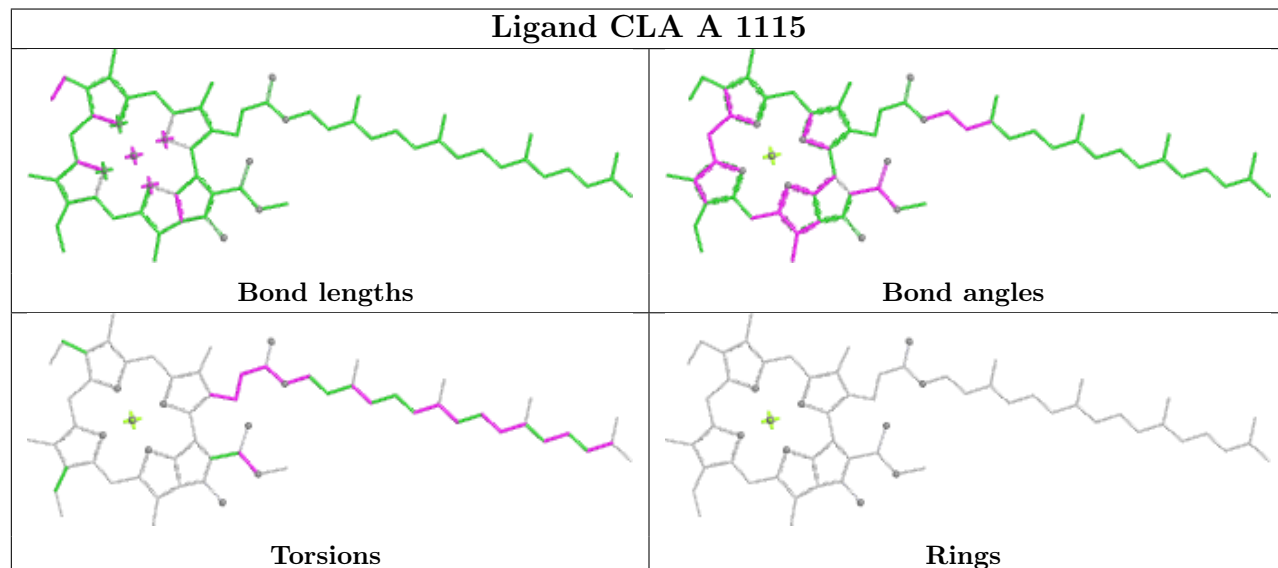


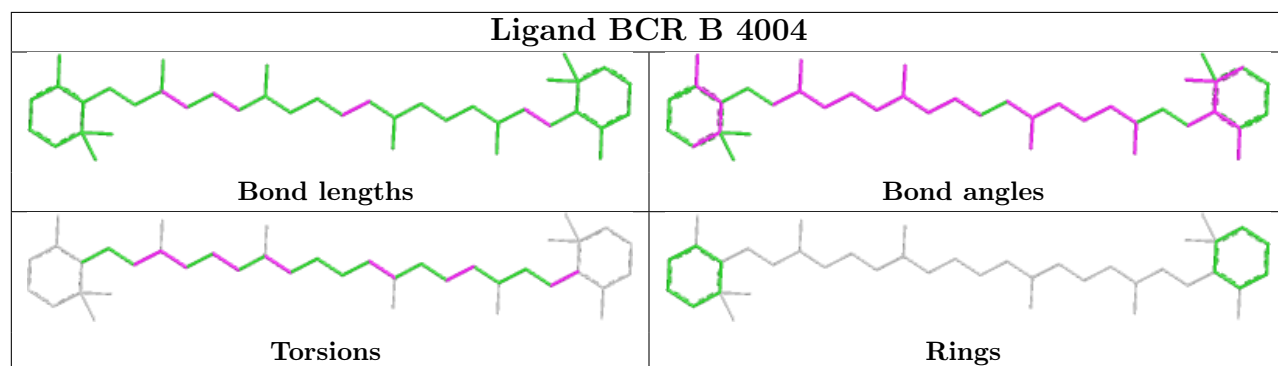
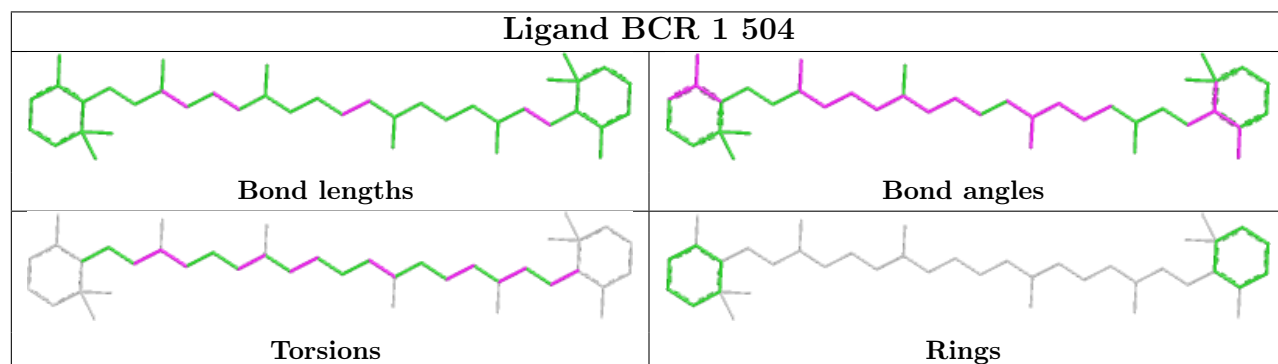
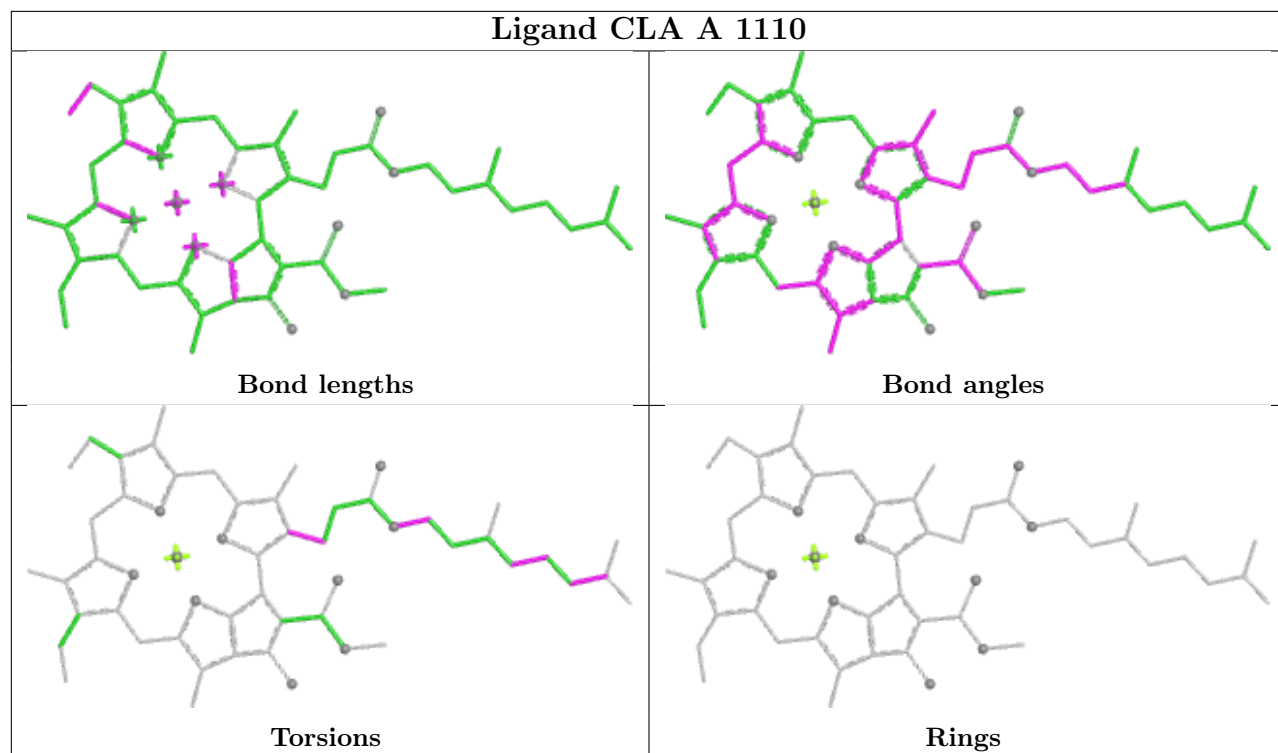
Ligand BCR L 4019	
	
Bond lengths	Bond angles
	
Torsions	Rings

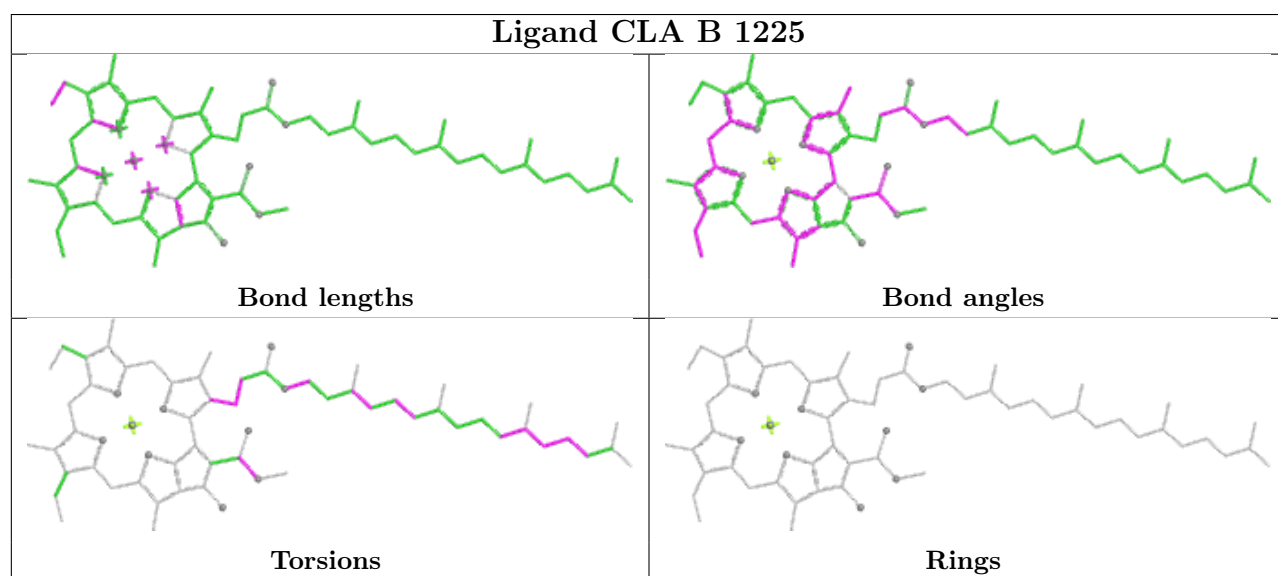
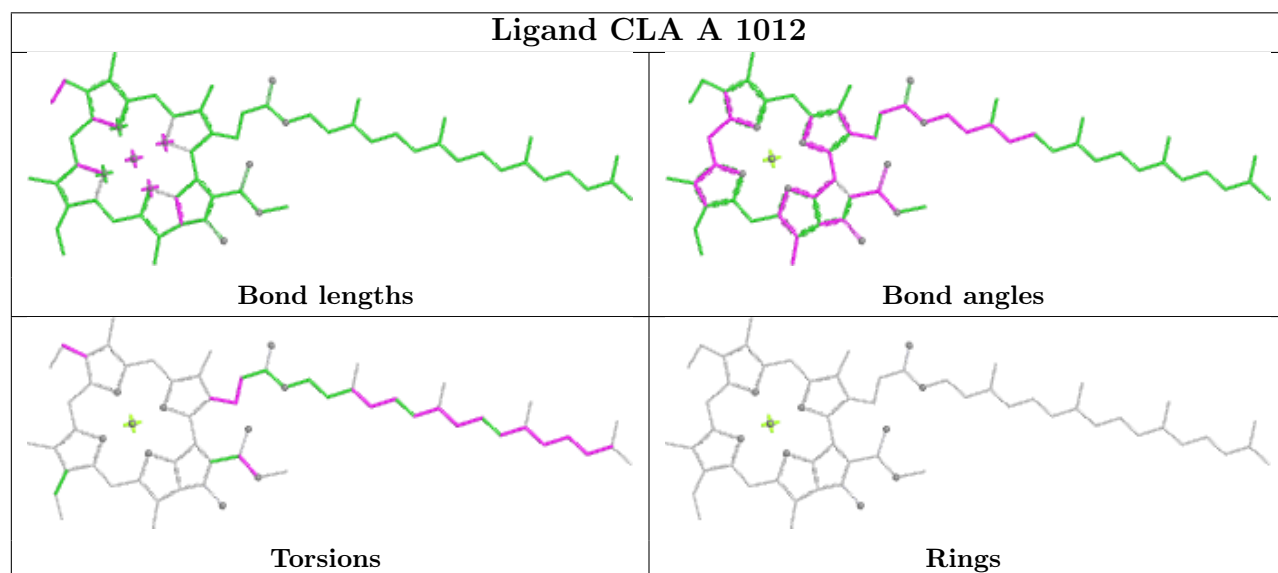
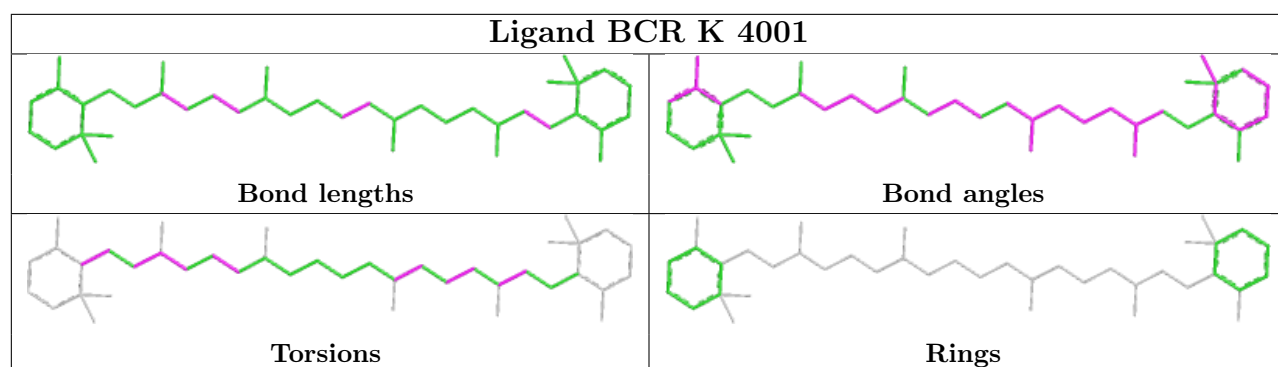
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Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA 2 604	
	
Bond lengths	Bond angles
	
Torsions	Rings

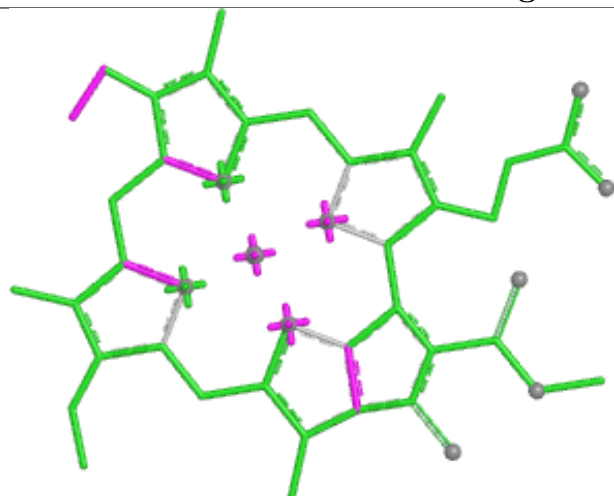


Ligand CLA 4 612**Ligand CLA A 1115**

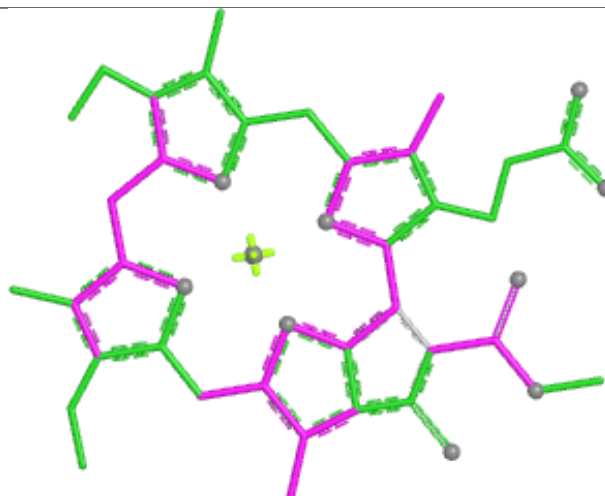




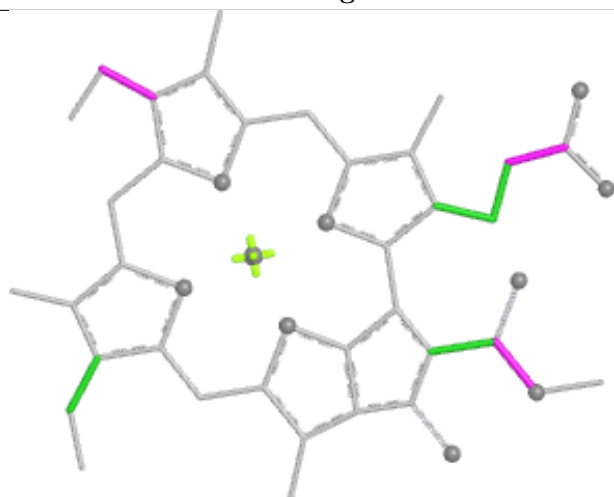
Ligand CLA 1 613



Bond lengths



Bond angles

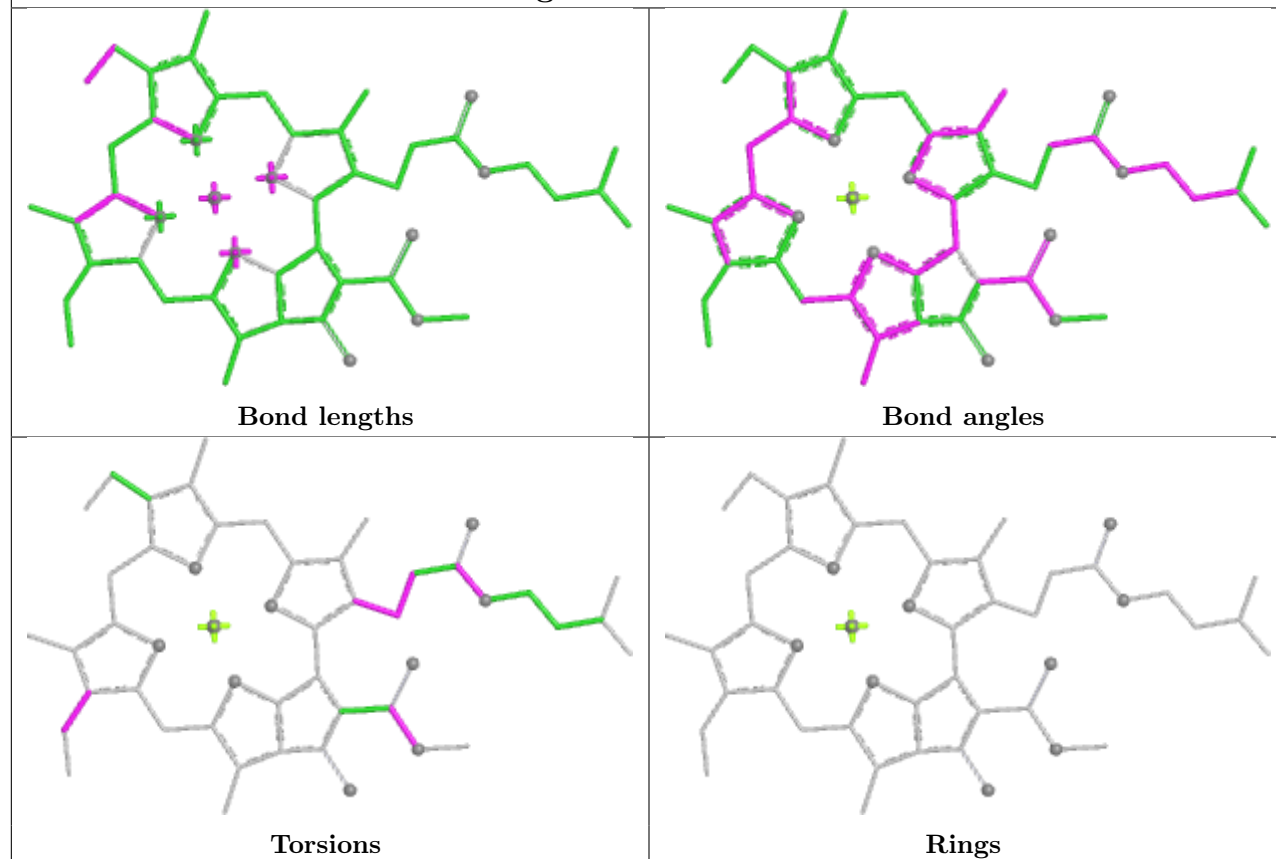


Torsions

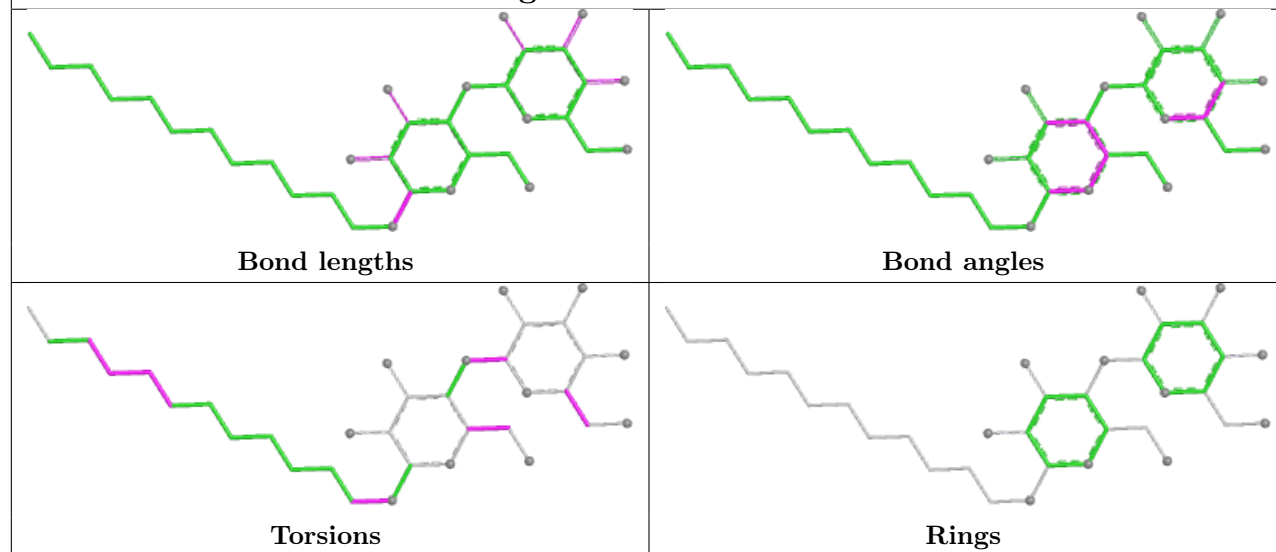


Rings

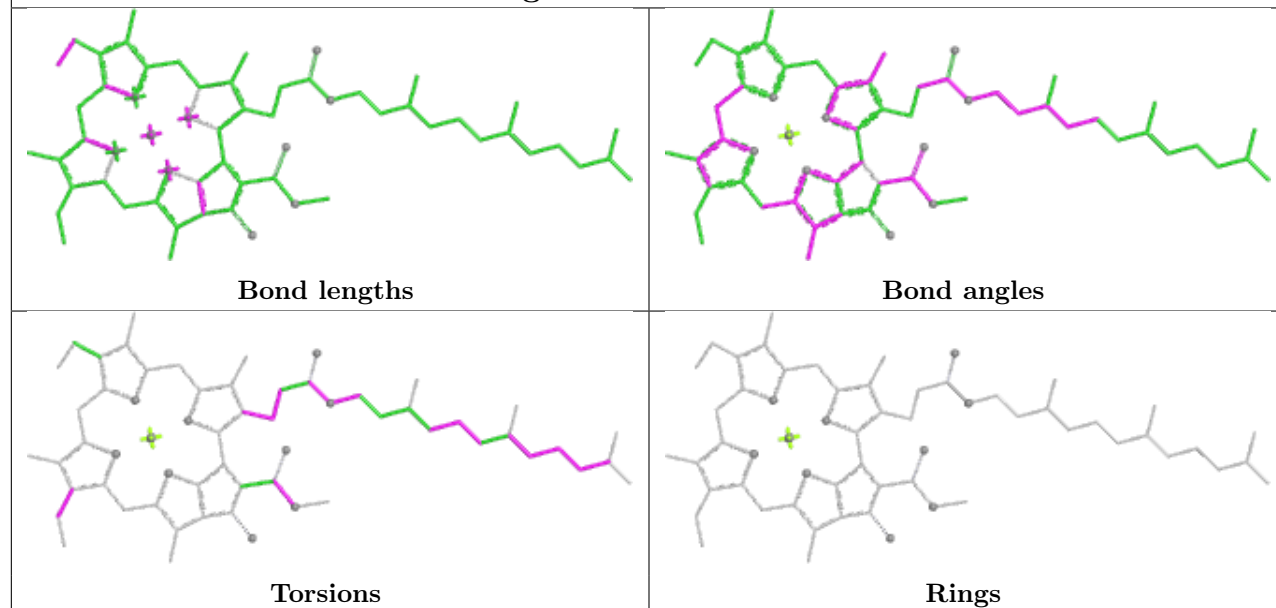
Ligand CLA 2 606



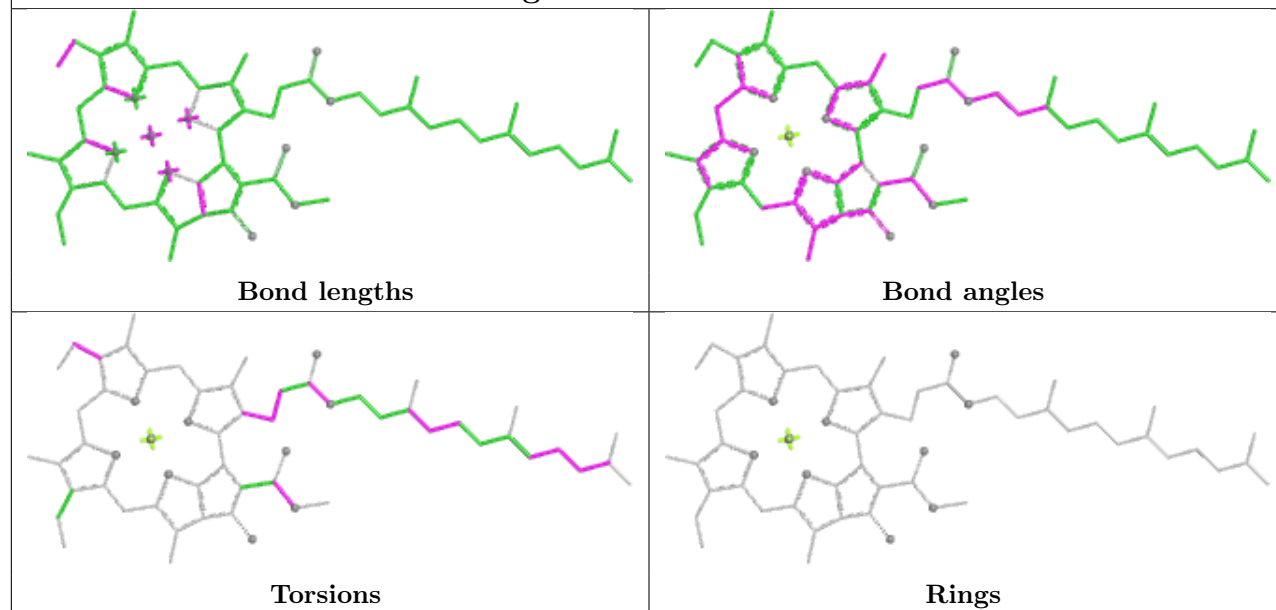
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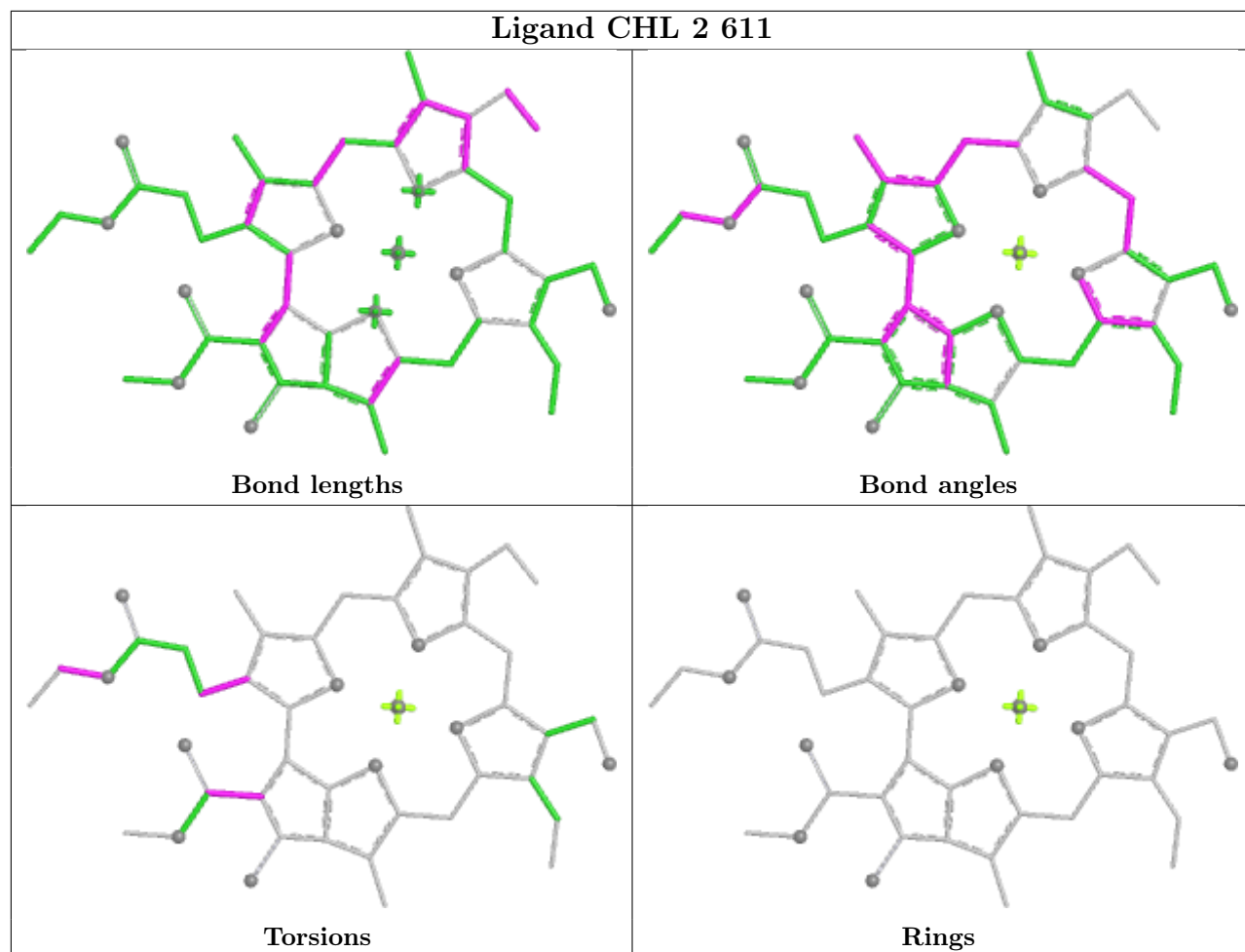
Ligand CLA A 1105



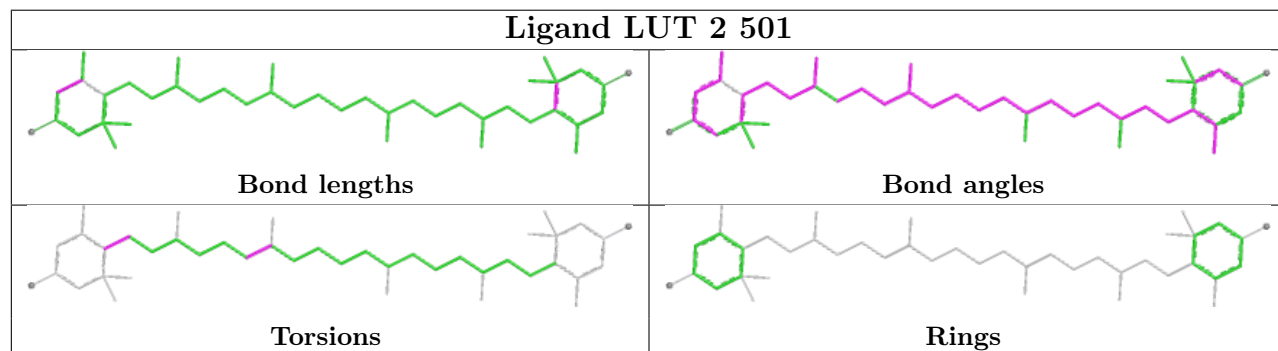
Ligand CLA A 1121

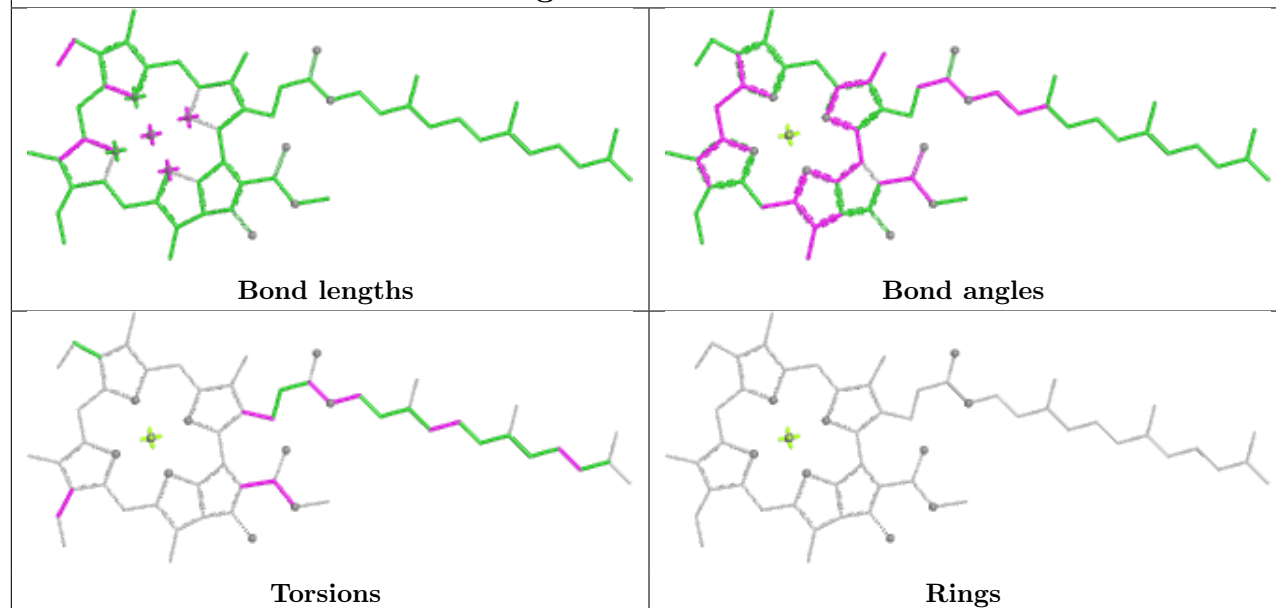
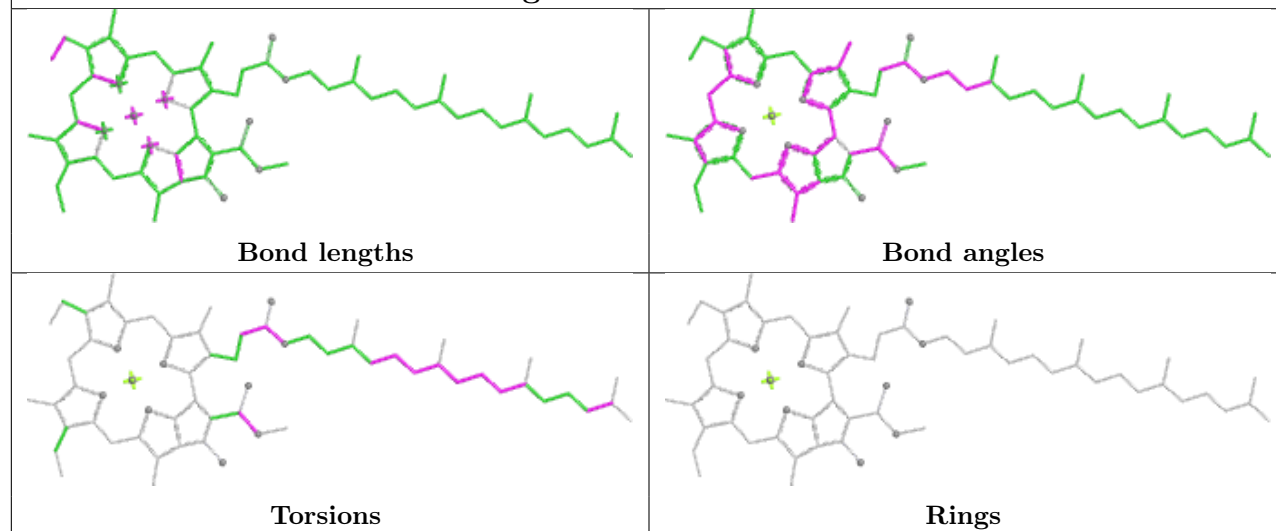


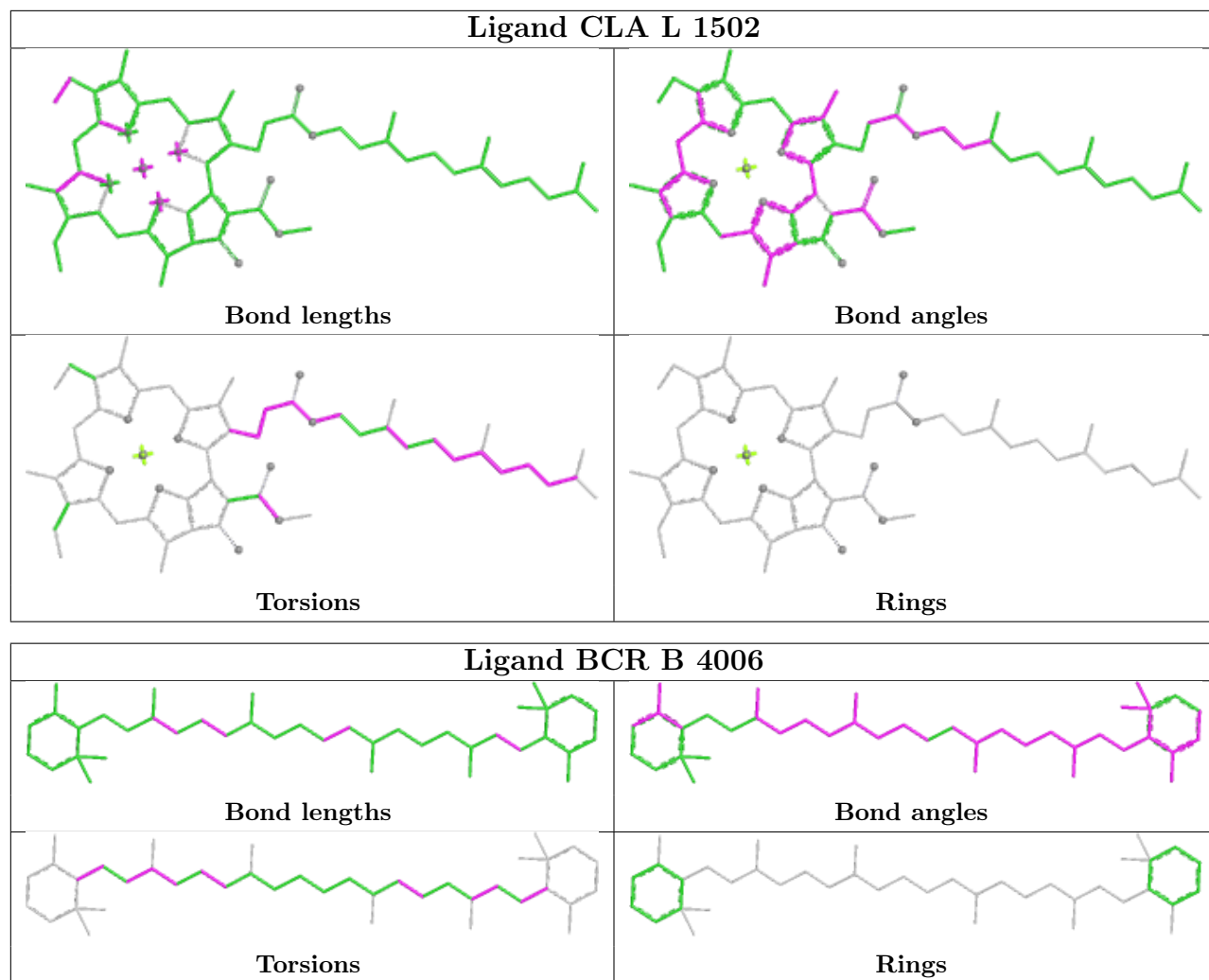
Ligand CHL 2 611

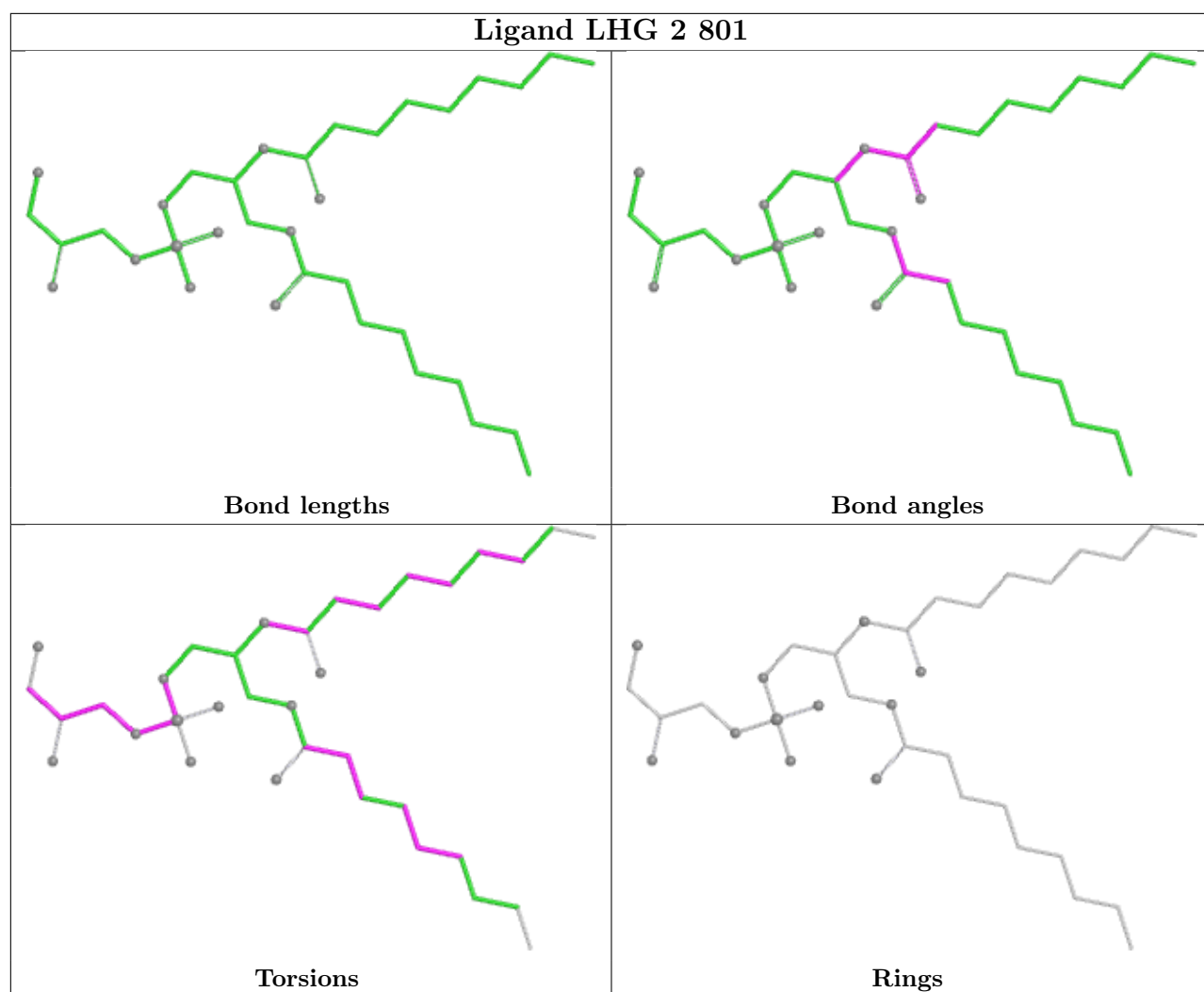


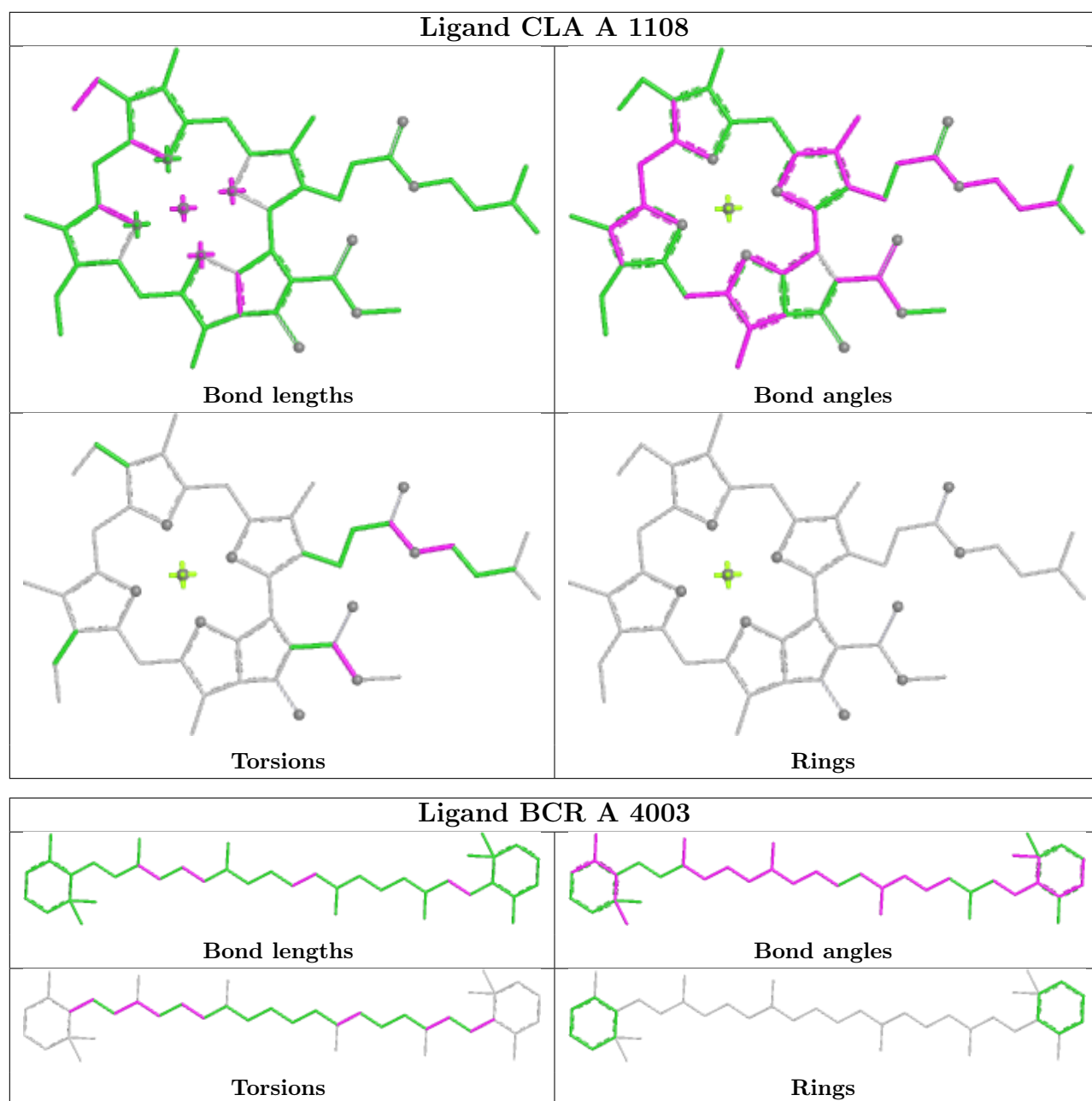
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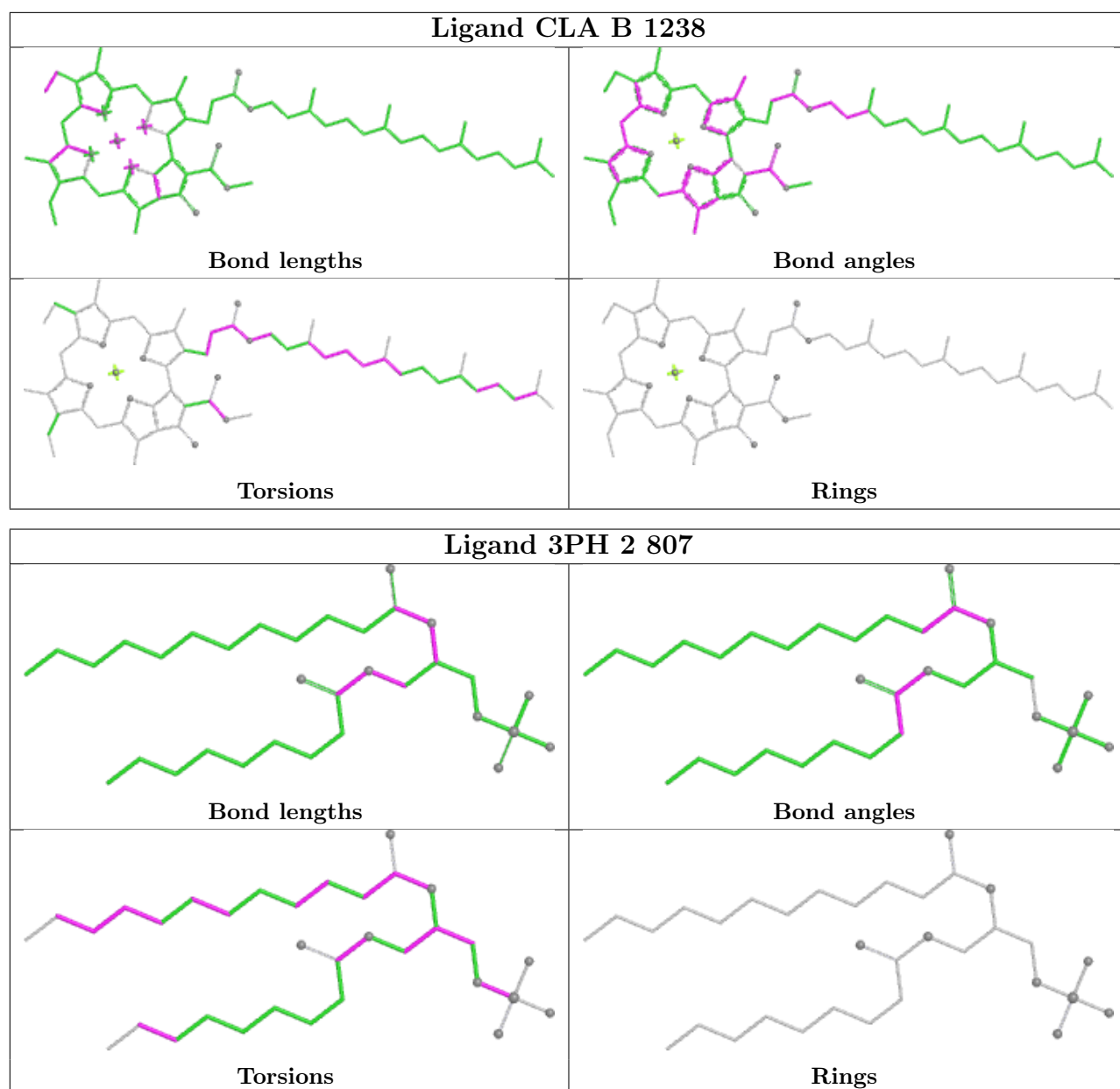


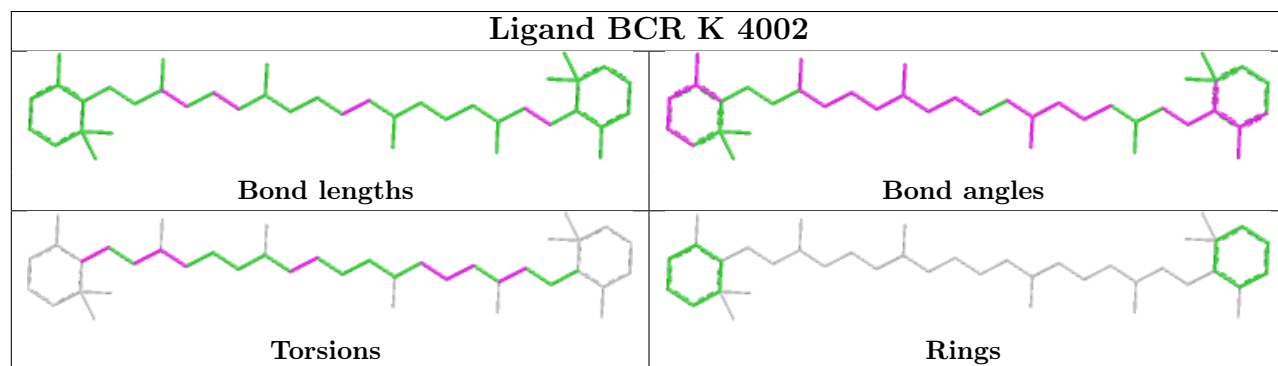
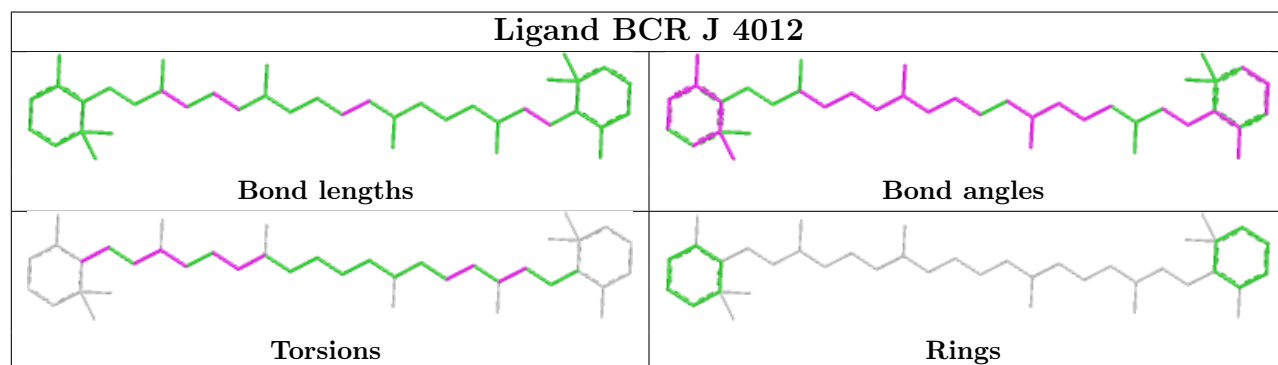
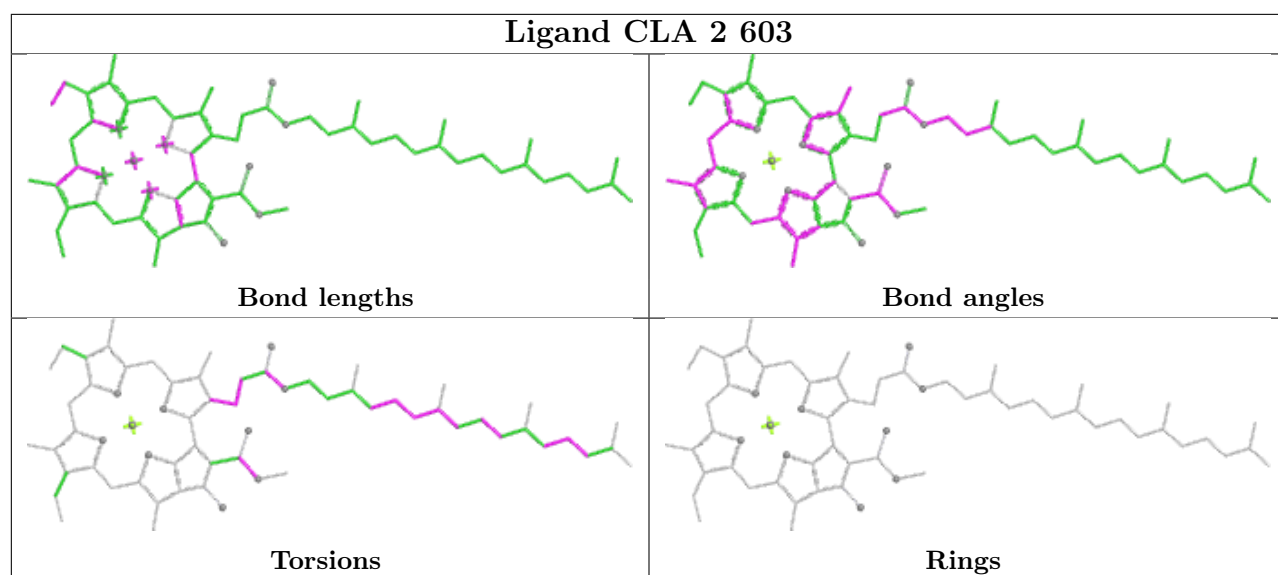
Ligand CLA 3 617**Ligand CLA A 1125**

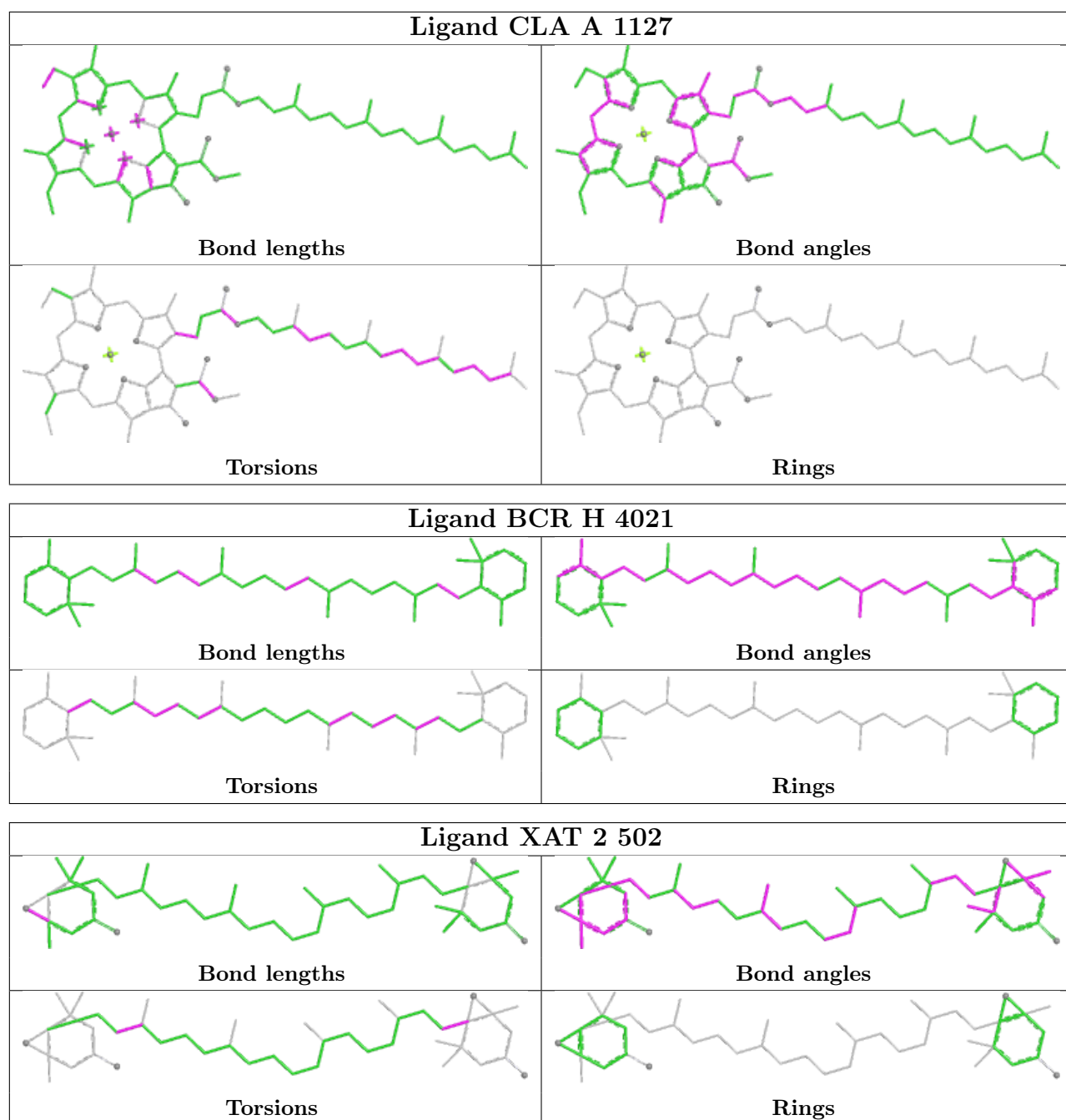




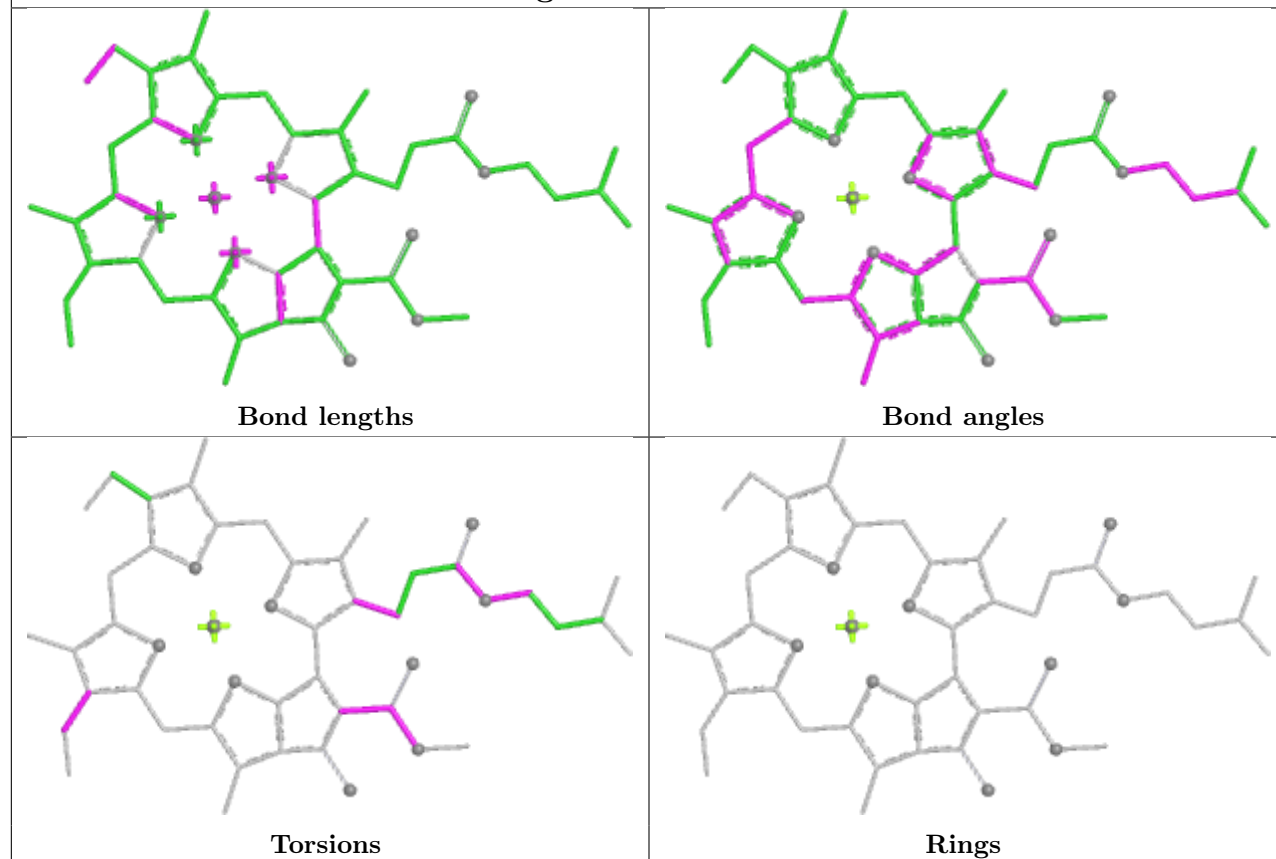




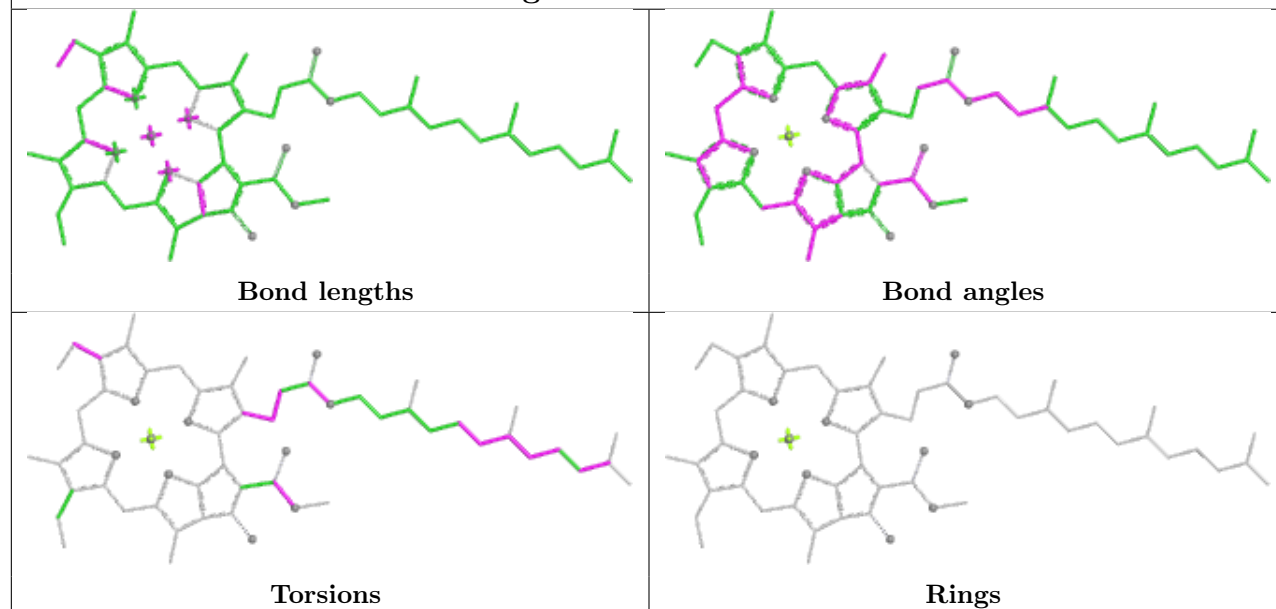


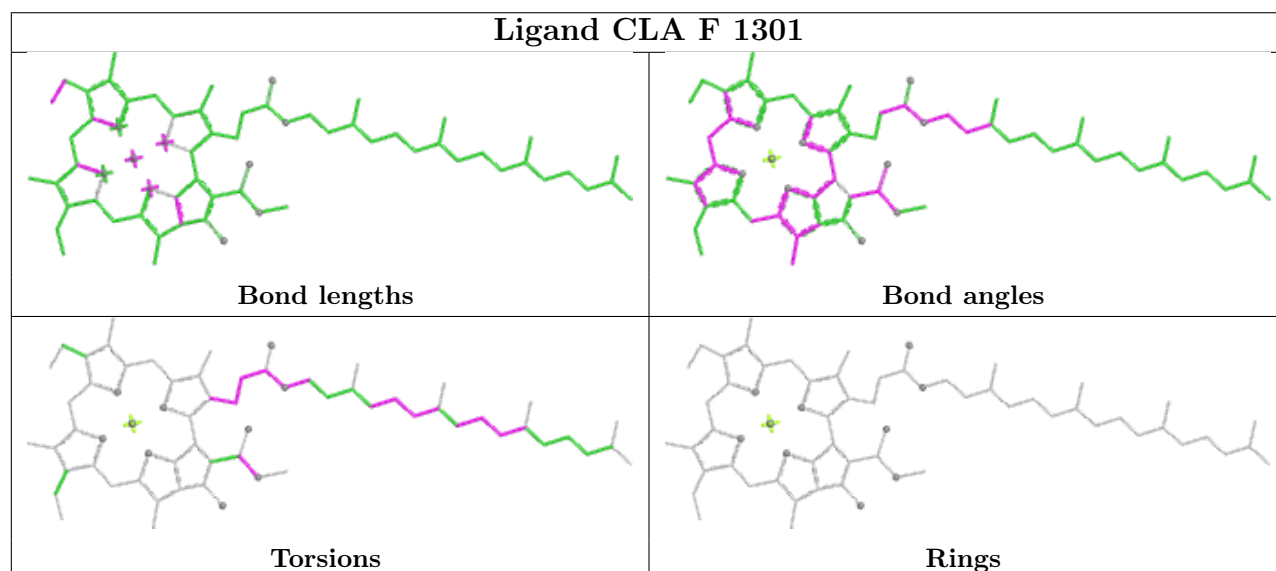
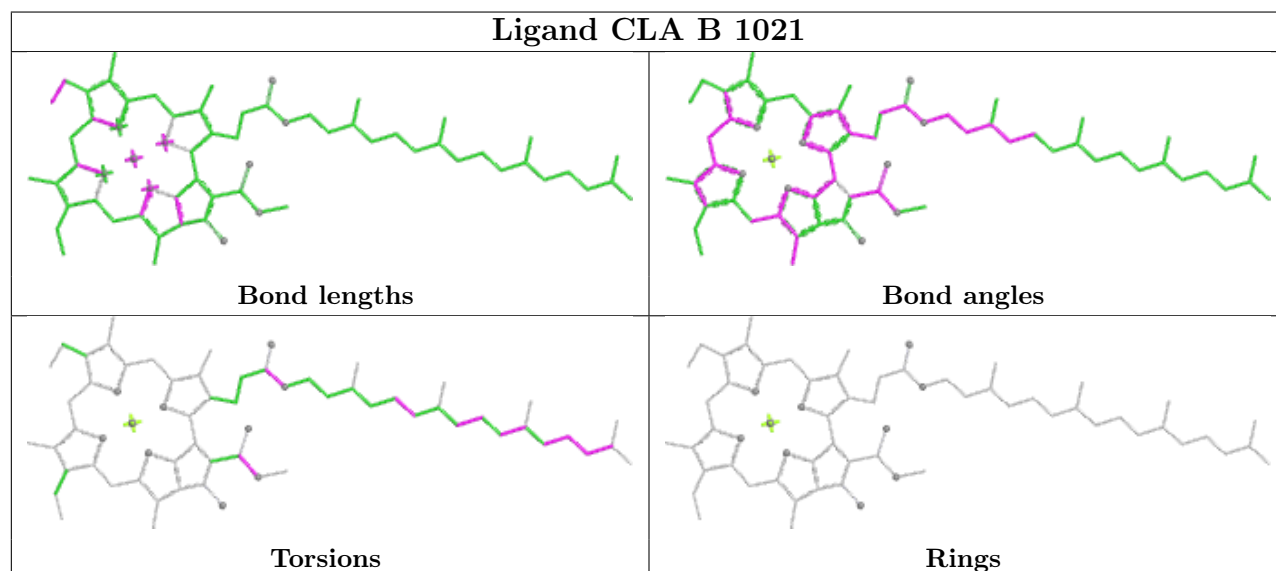
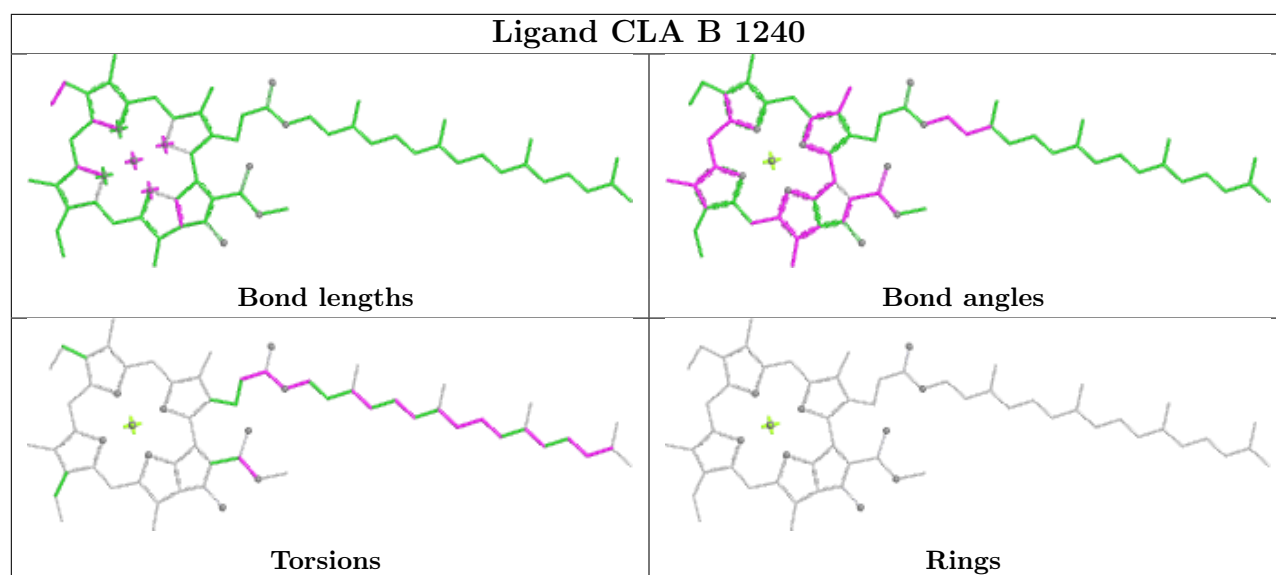


Ligand CLA 3 606

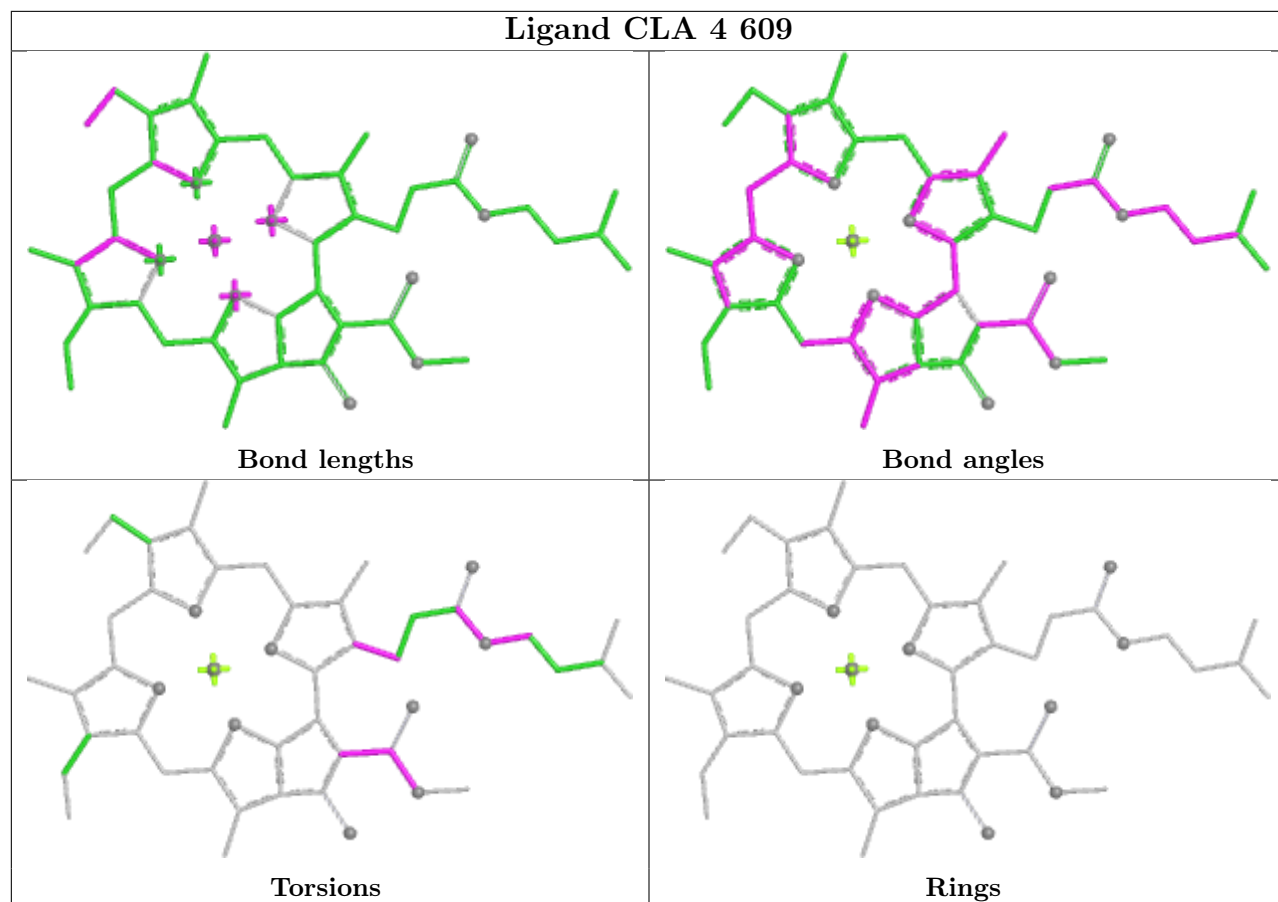


Ligand CLA A 1141

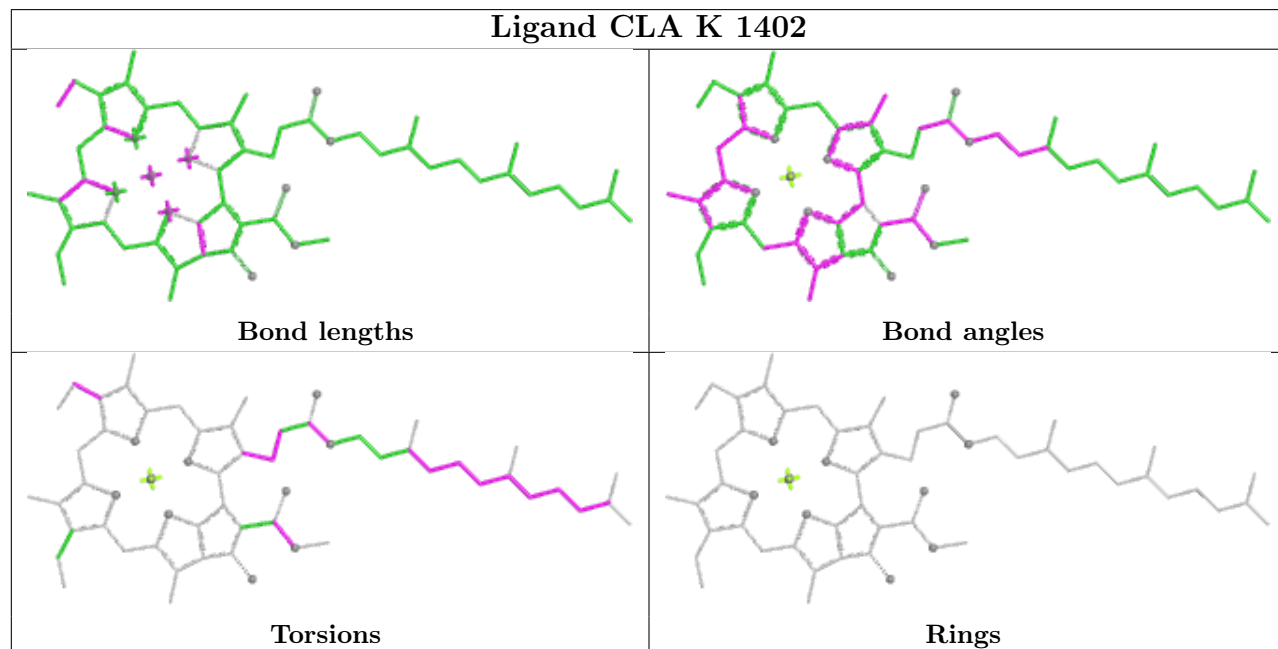


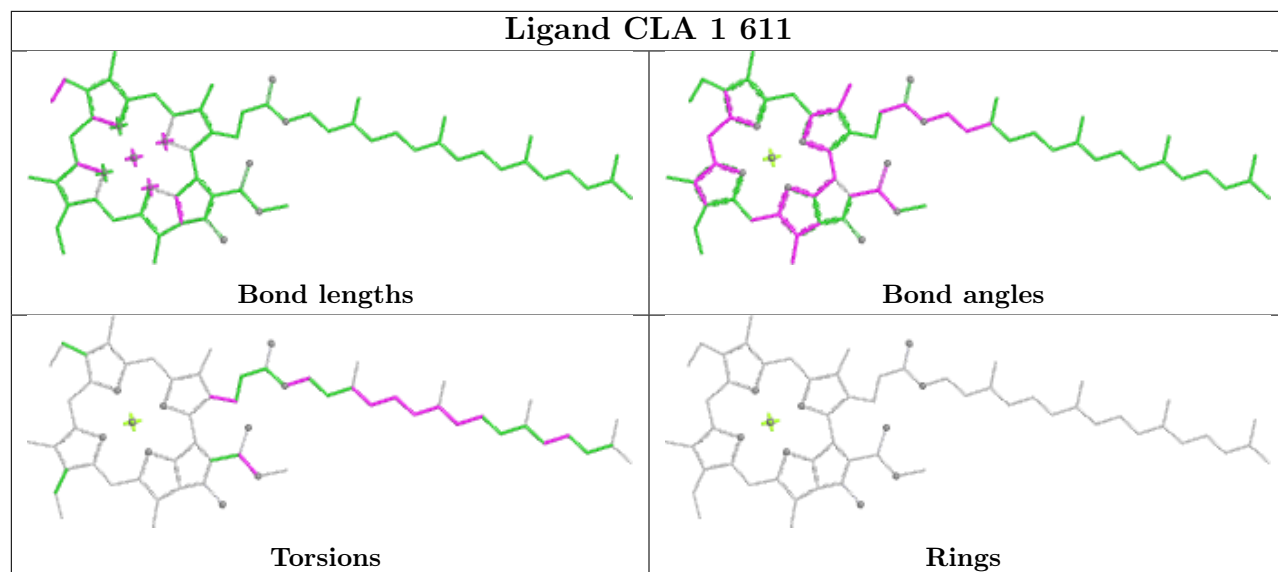
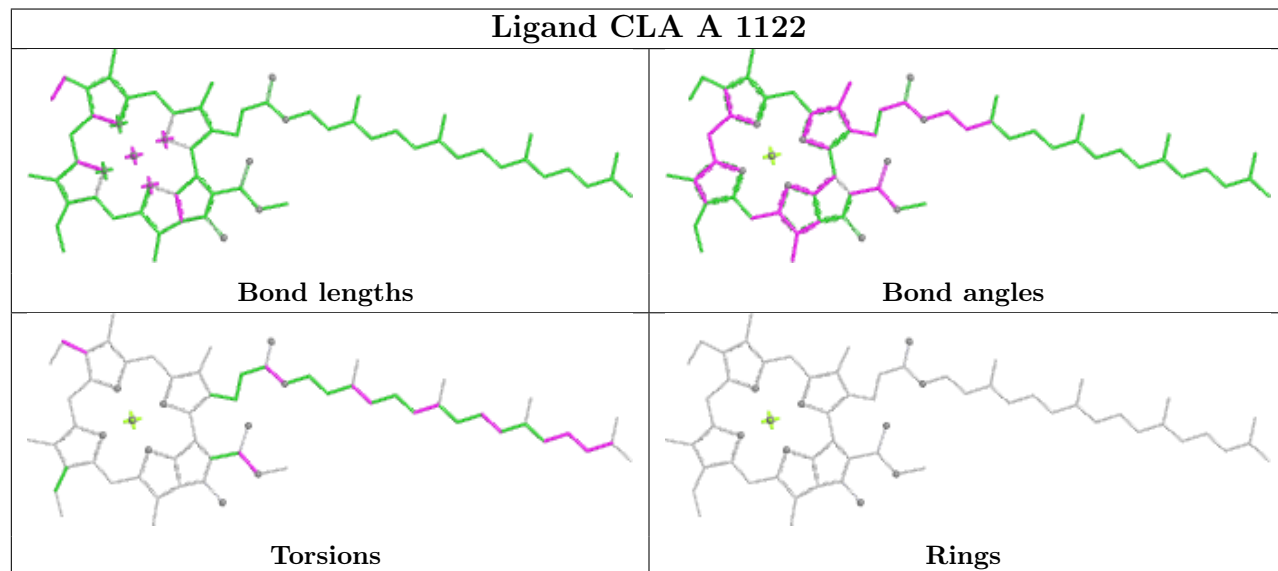


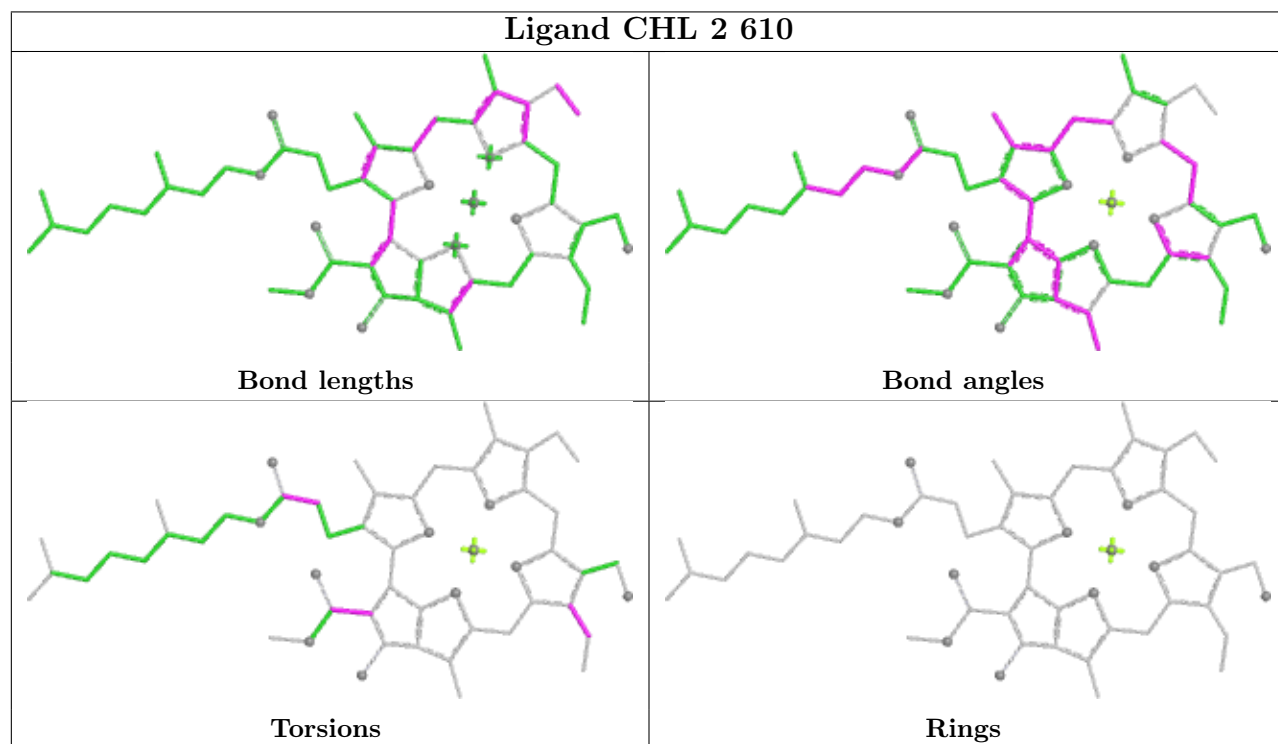
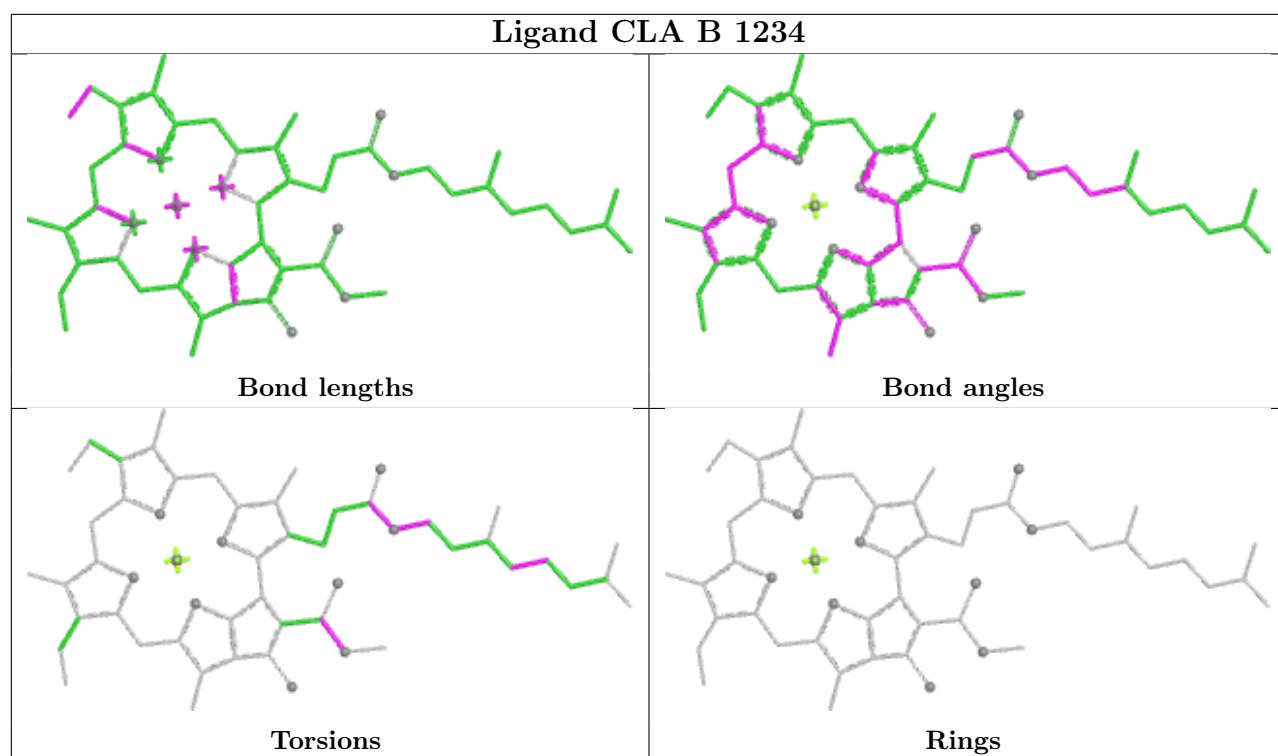
Ligand CLA 4 609

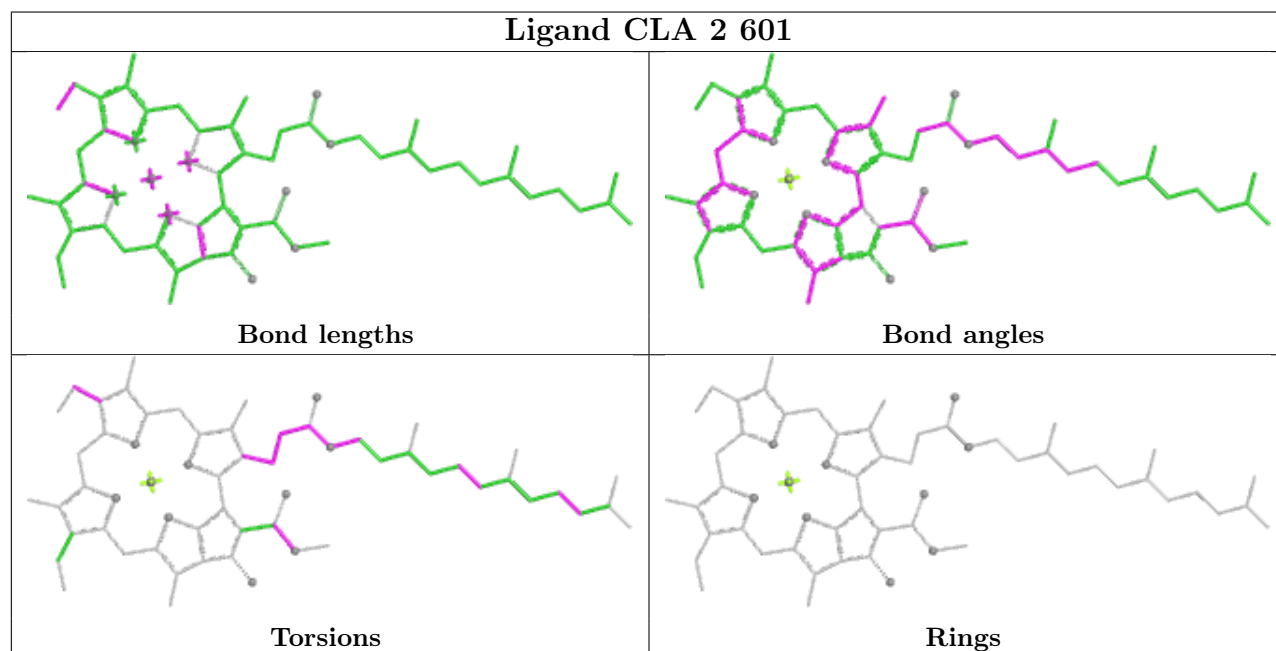
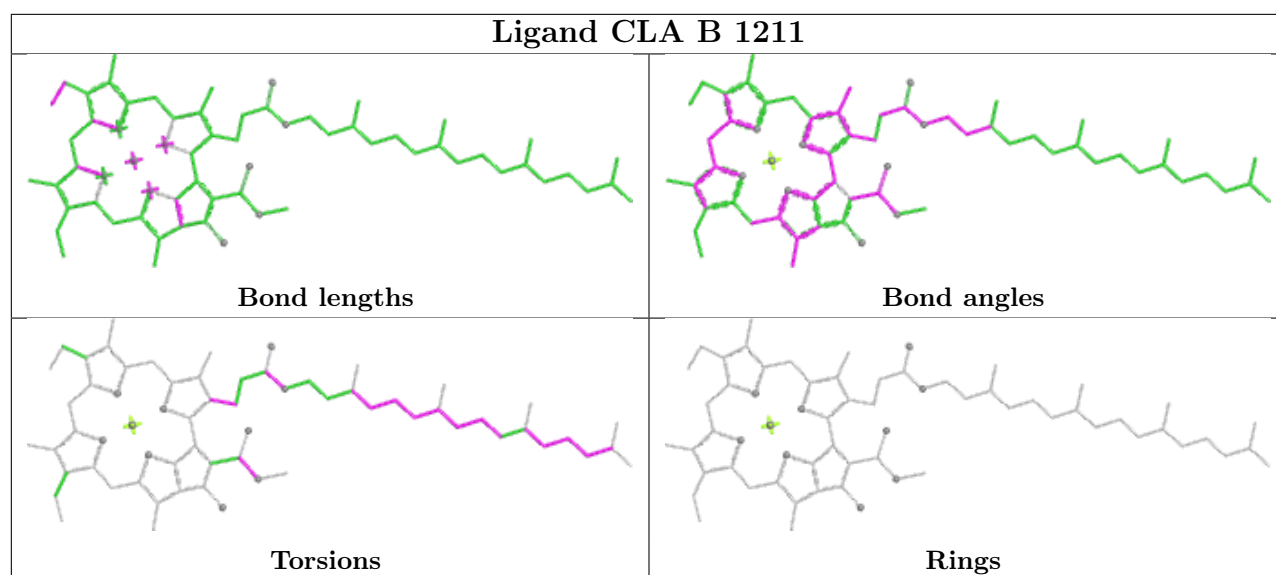


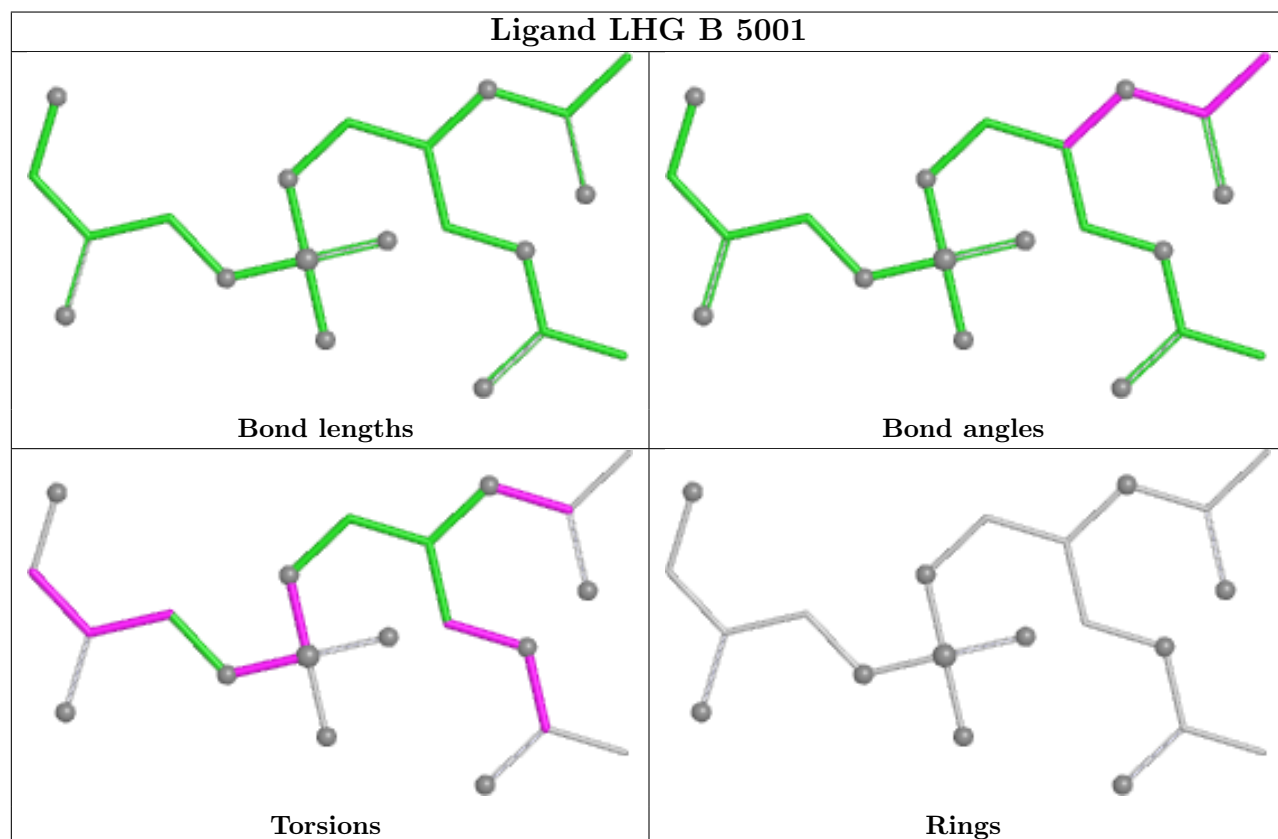
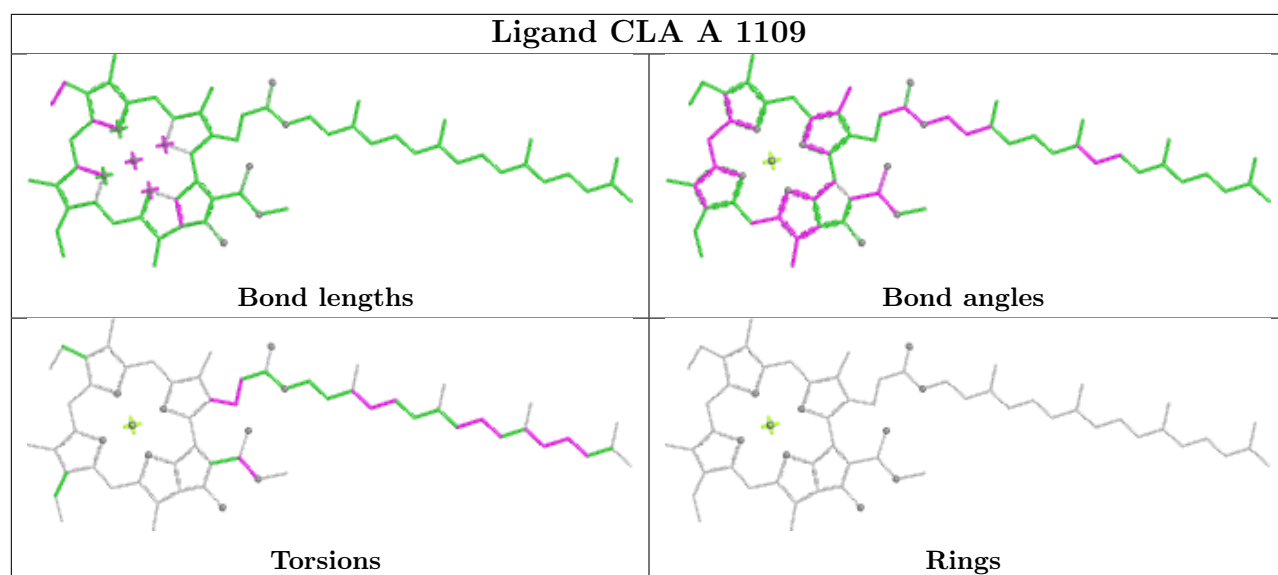
Ligand CLA K 1402

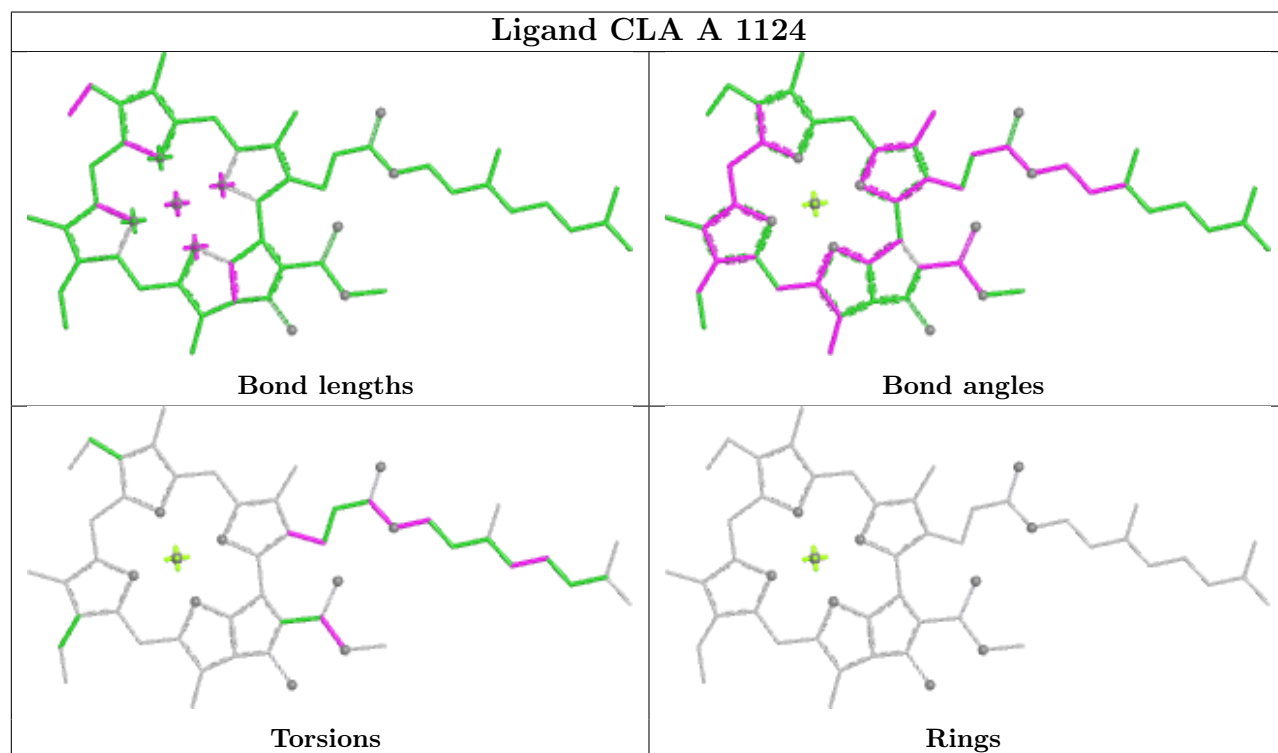
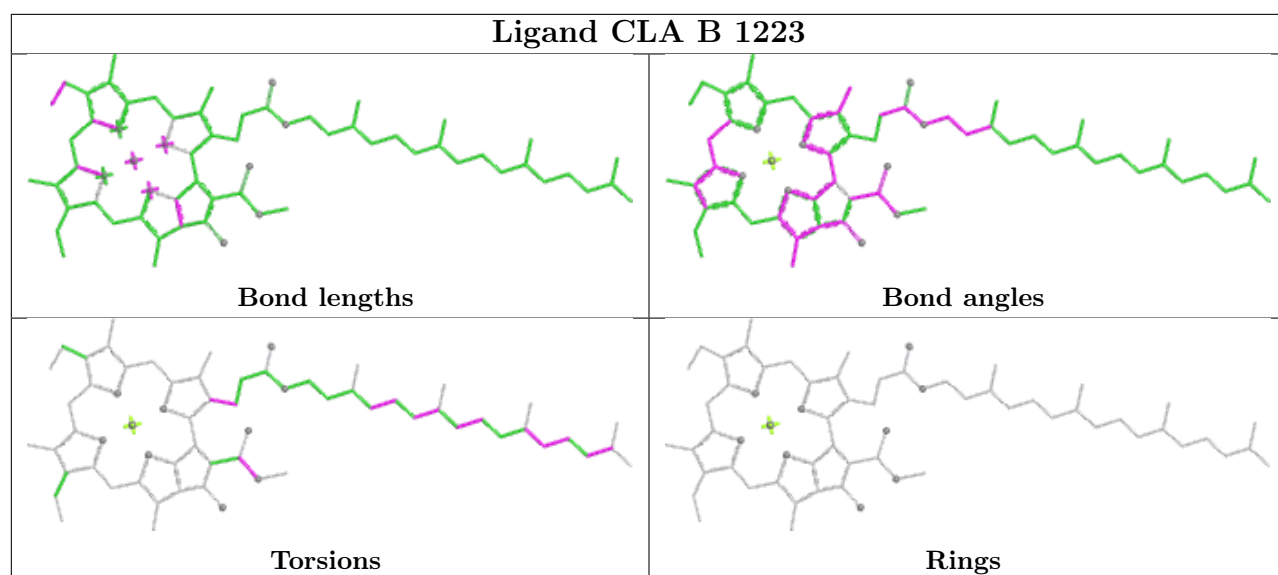


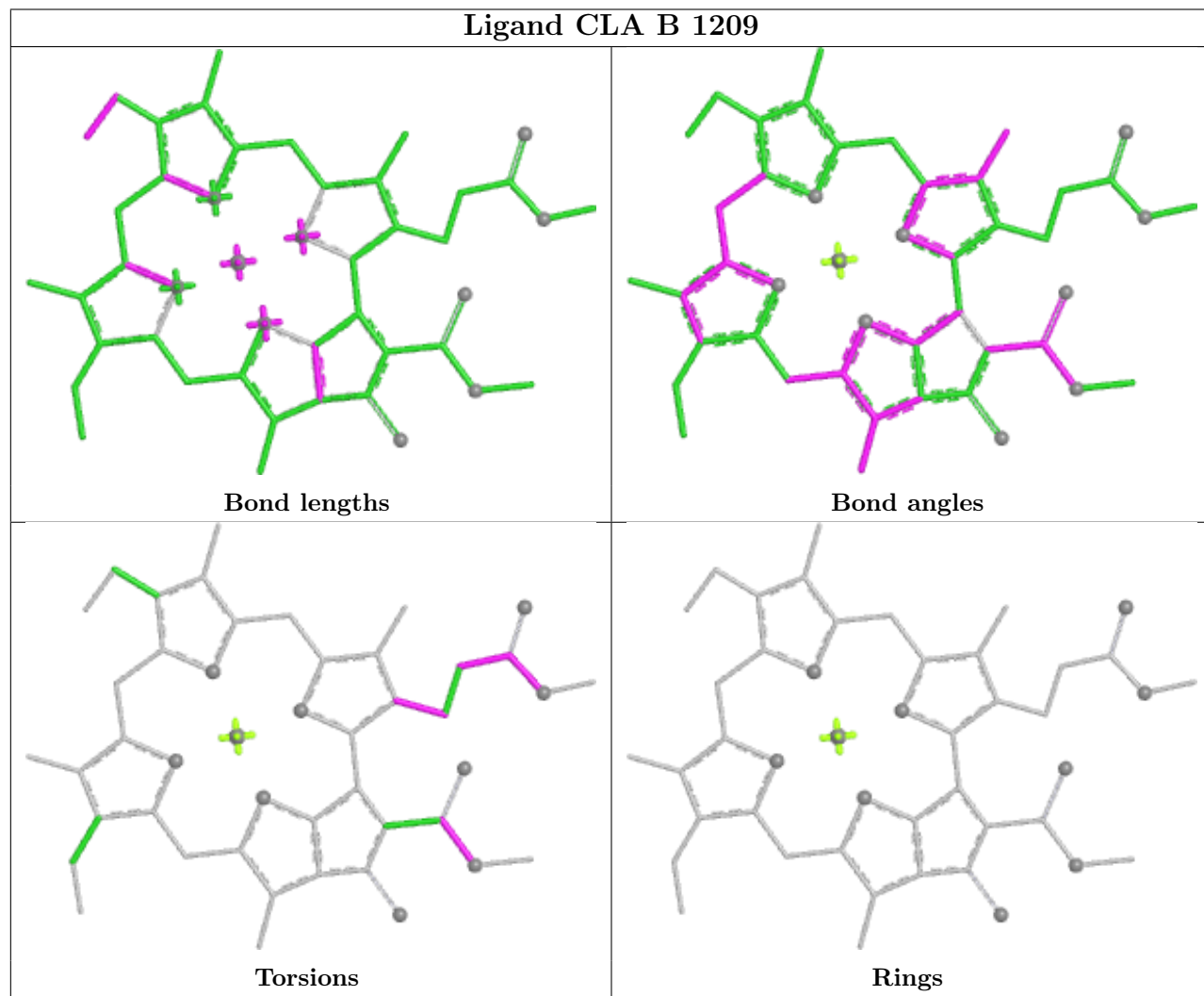
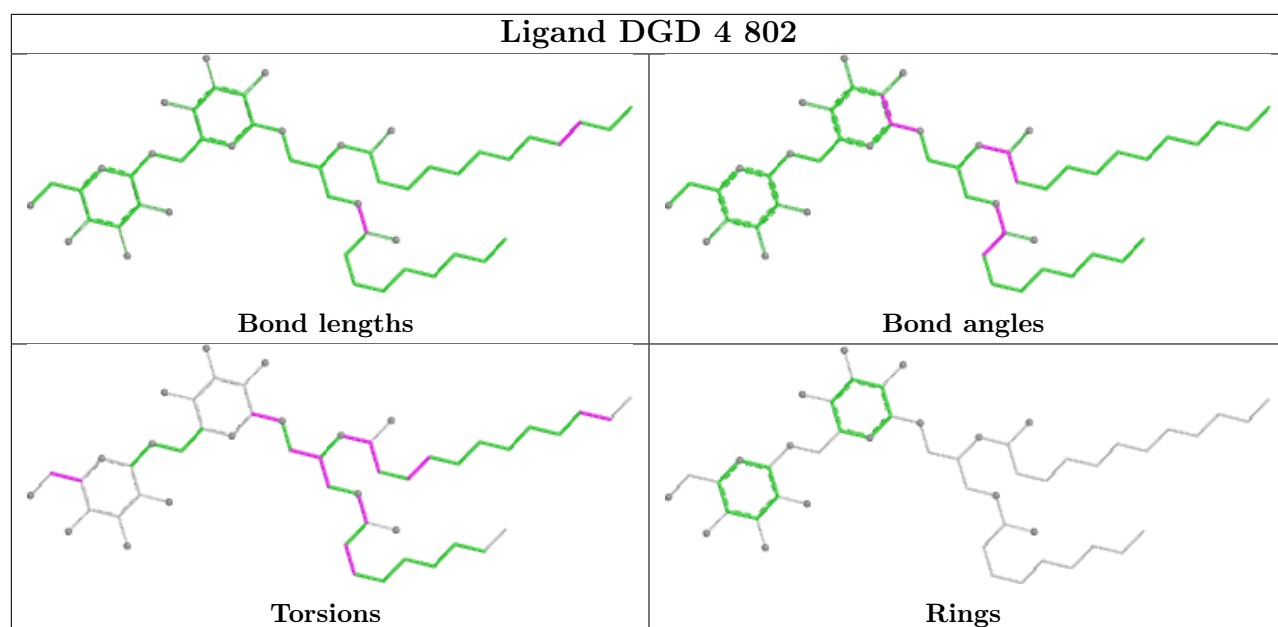
Ligand CLA 1 611**Ligand CLA A 1122**

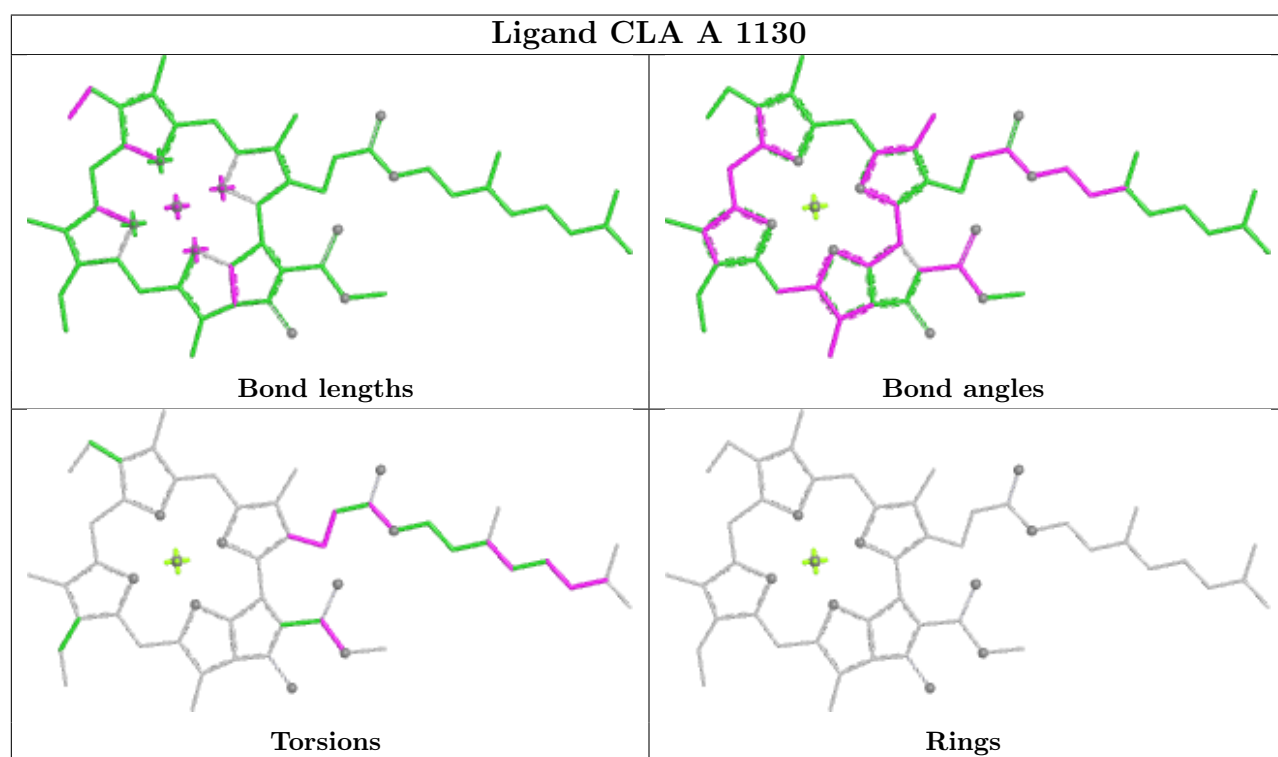
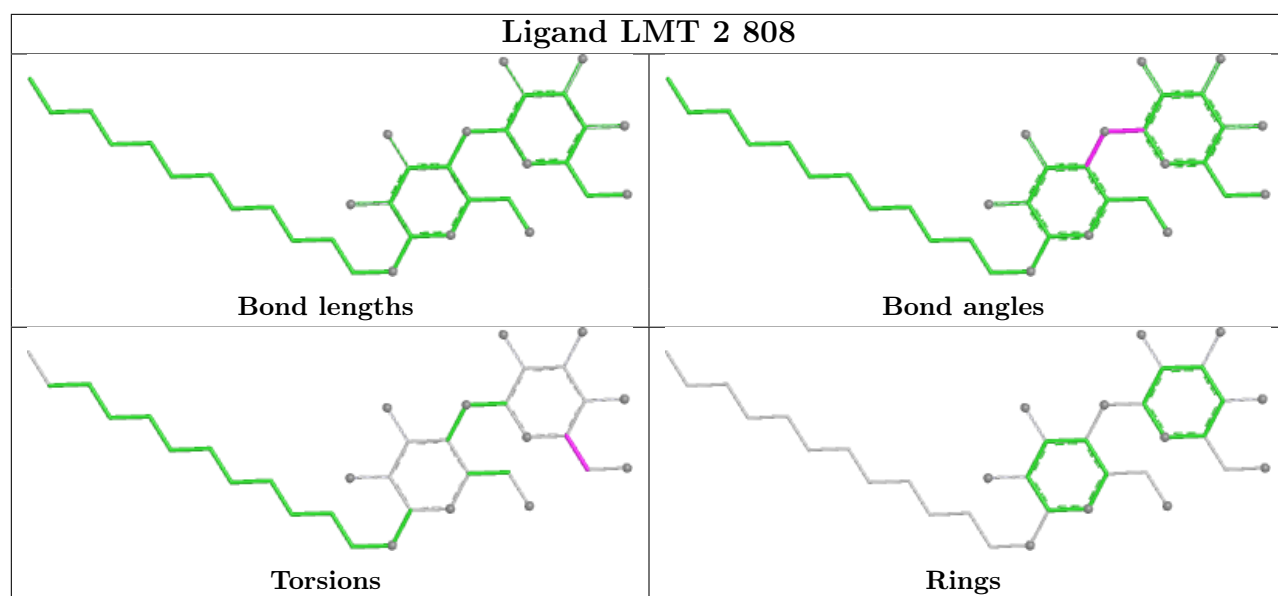




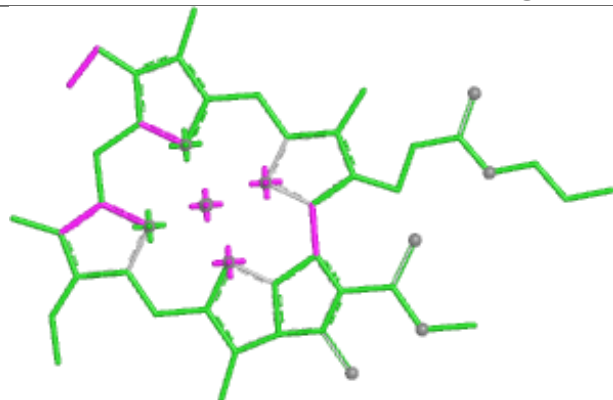




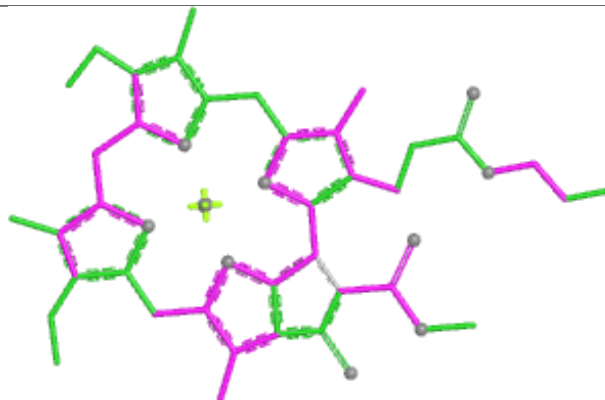




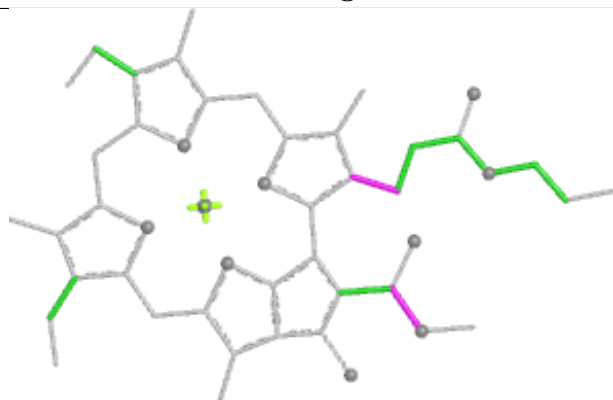
Ligand CLA 3 608



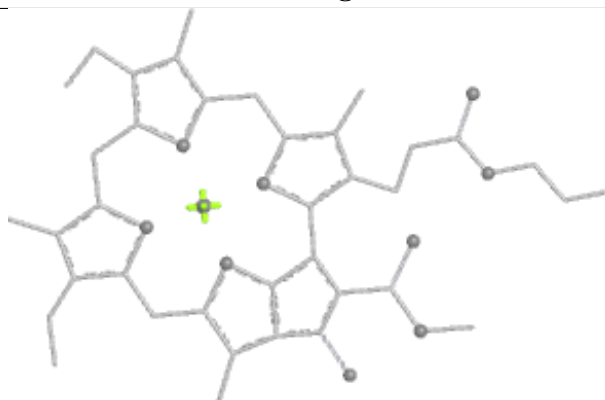
Bond lengths



Bond angles

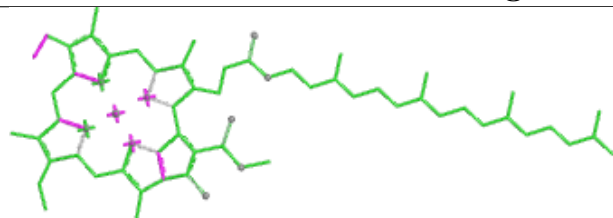


Torsions

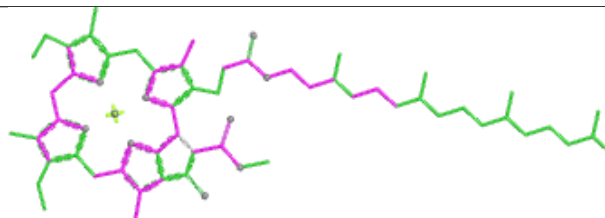


Rings

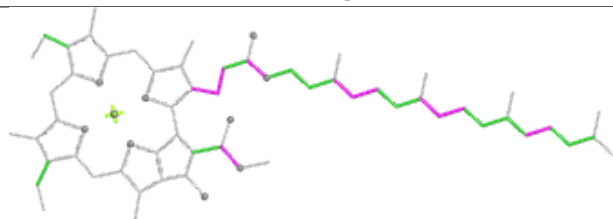
Ligand CLA A 1119



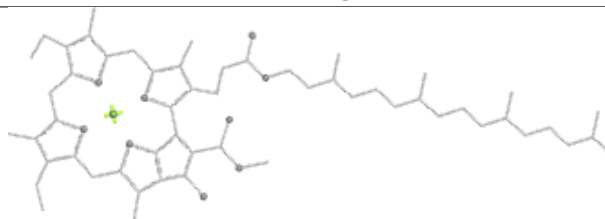
Bond lengths



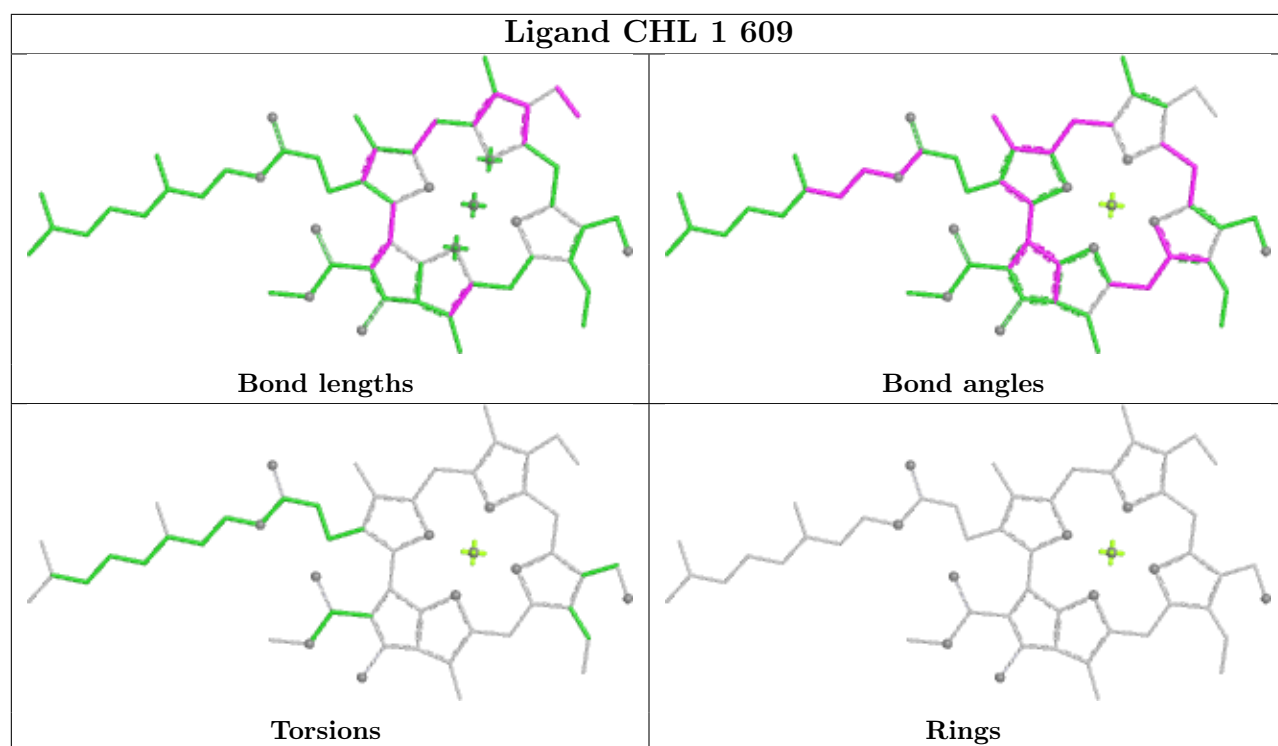
Bond angles



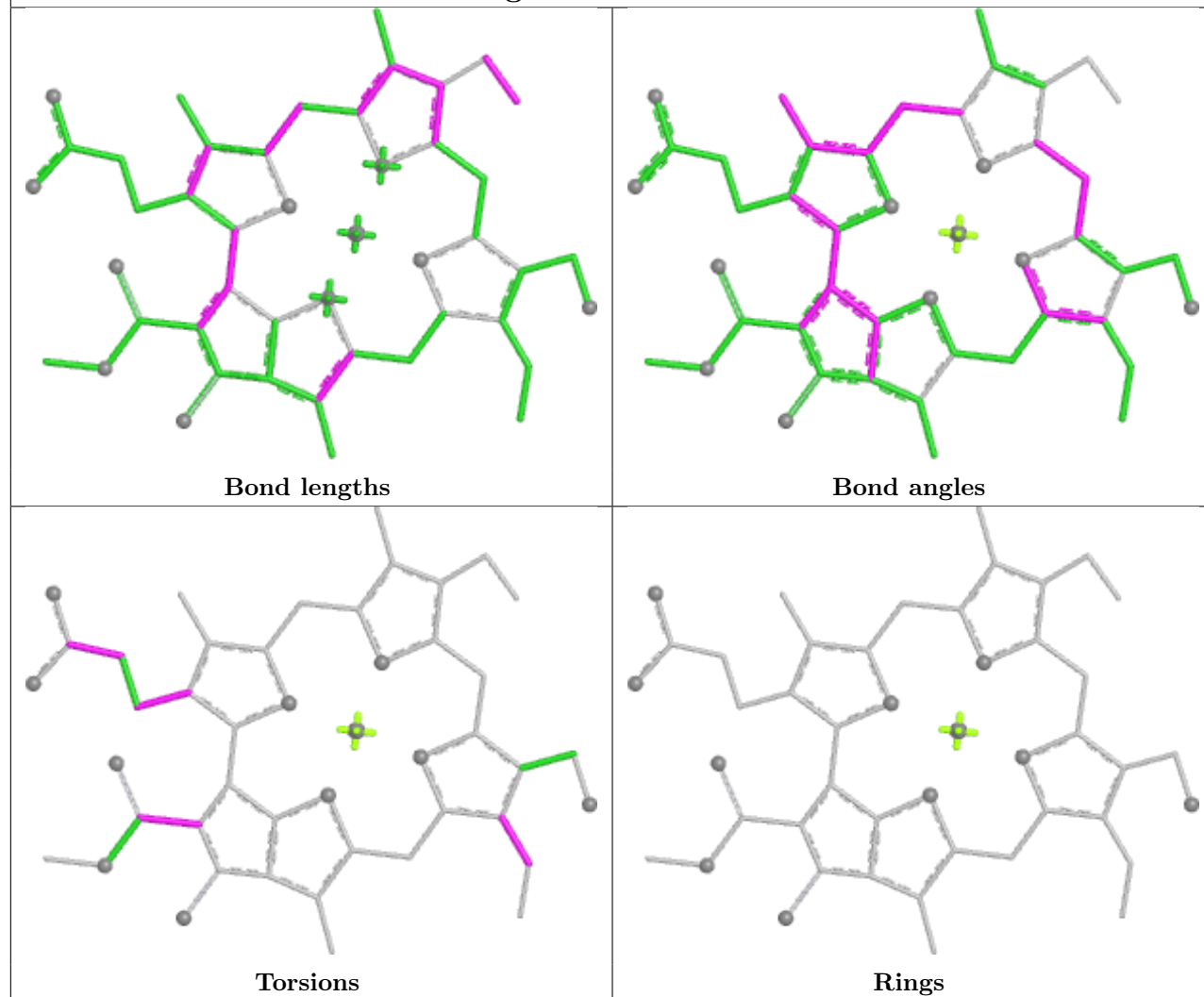
Torsions



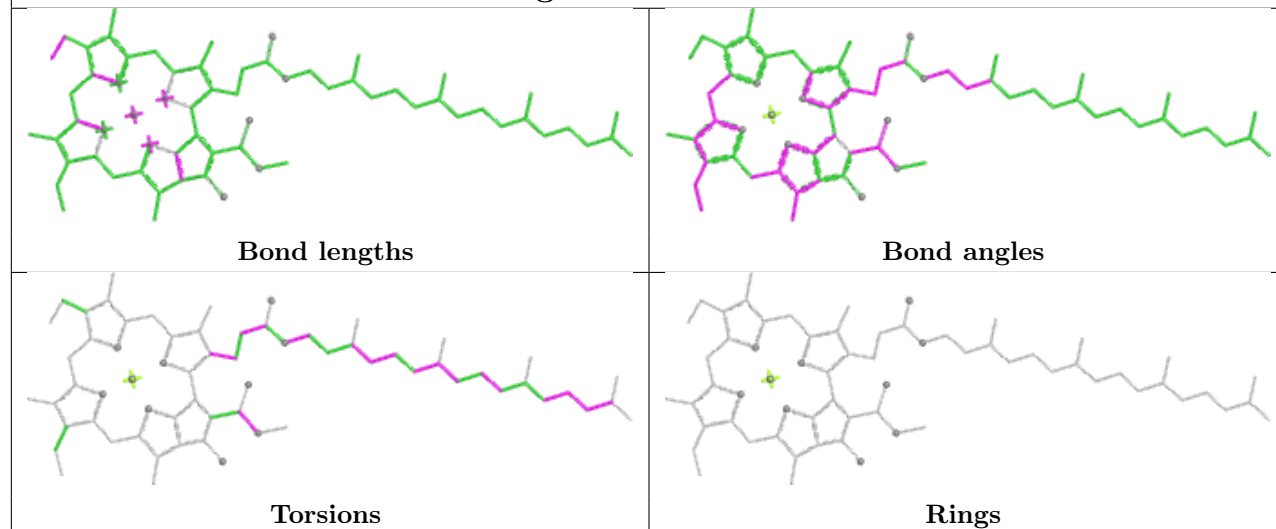
Rings

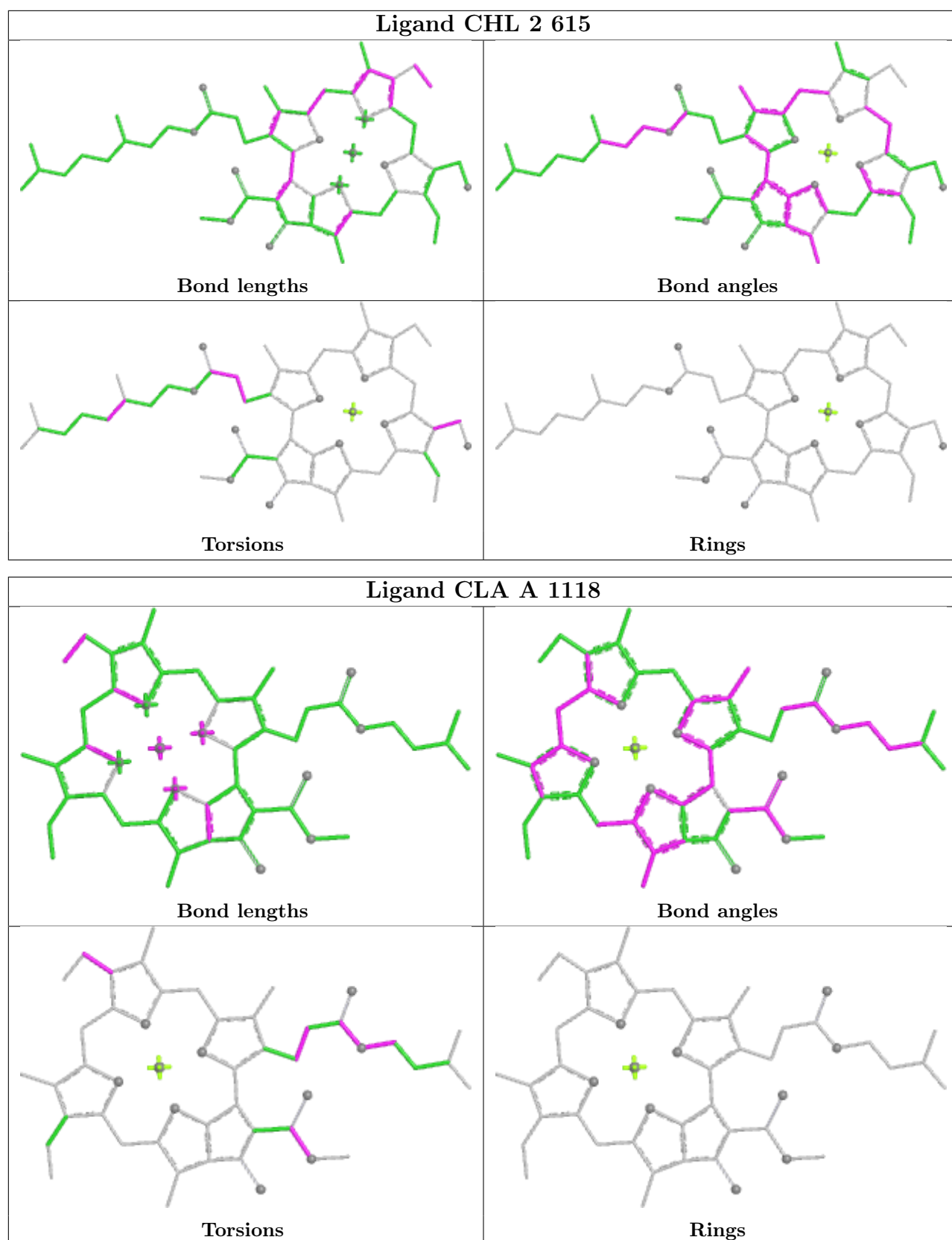


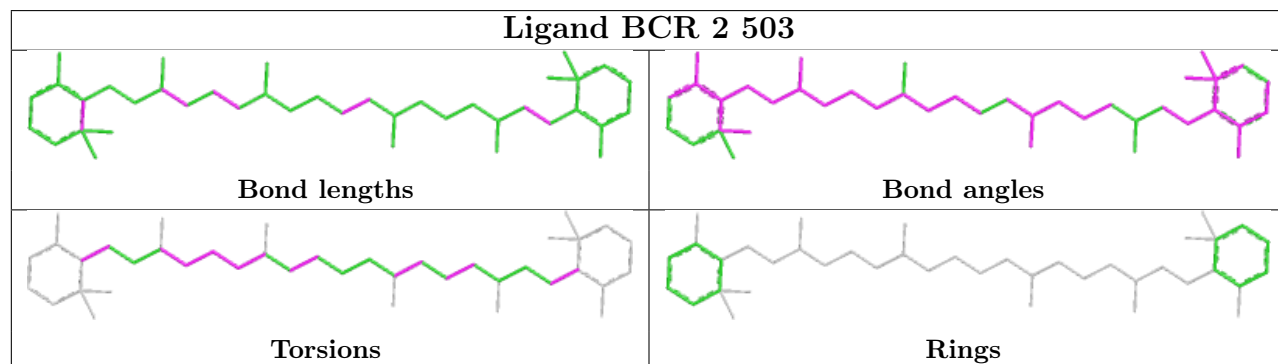
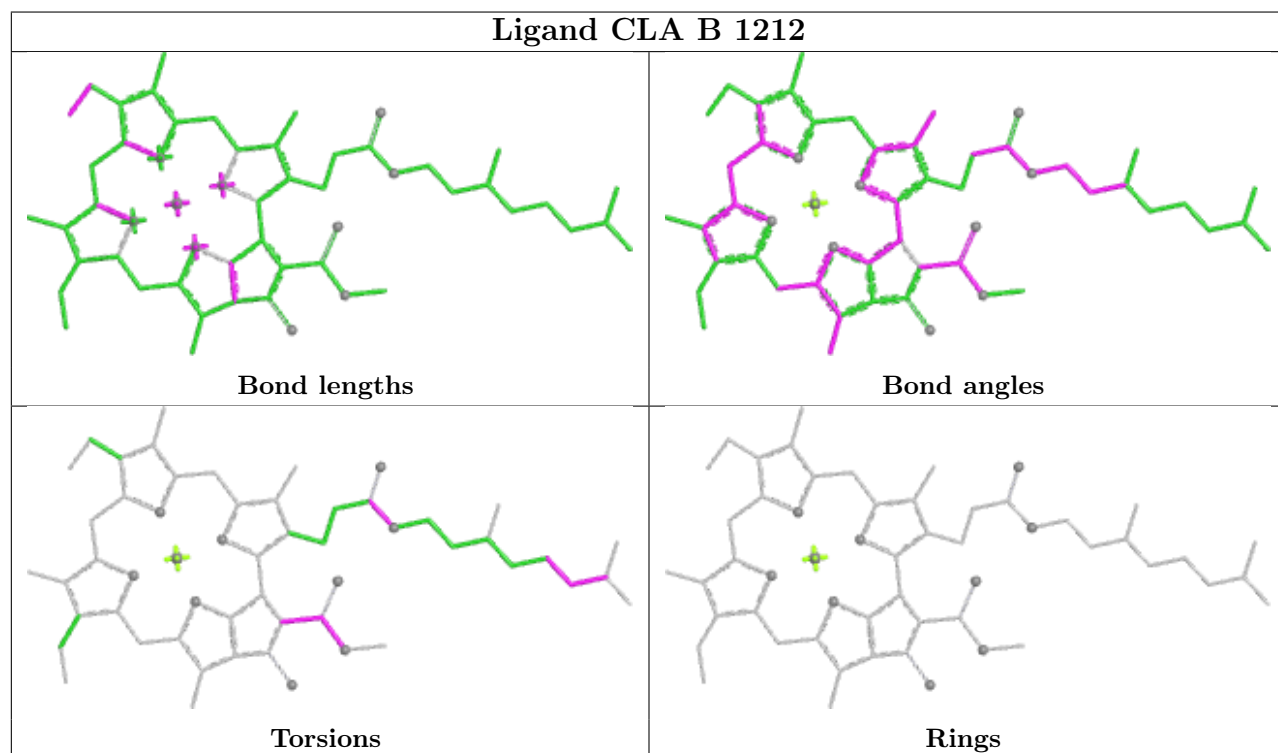
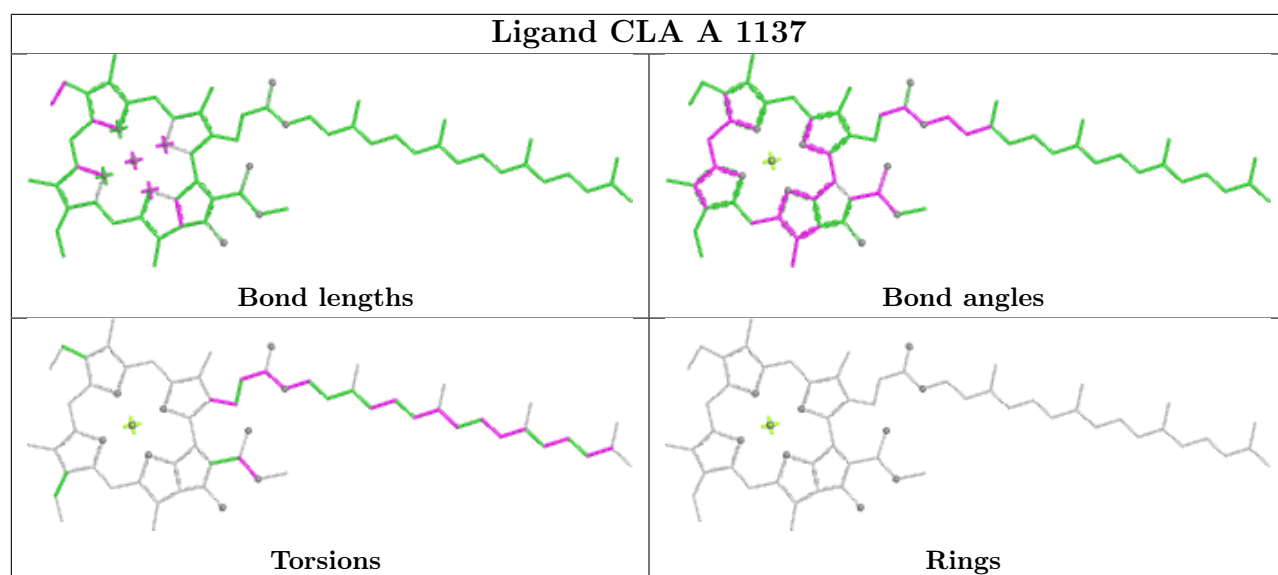
Ligand CHL 2 613

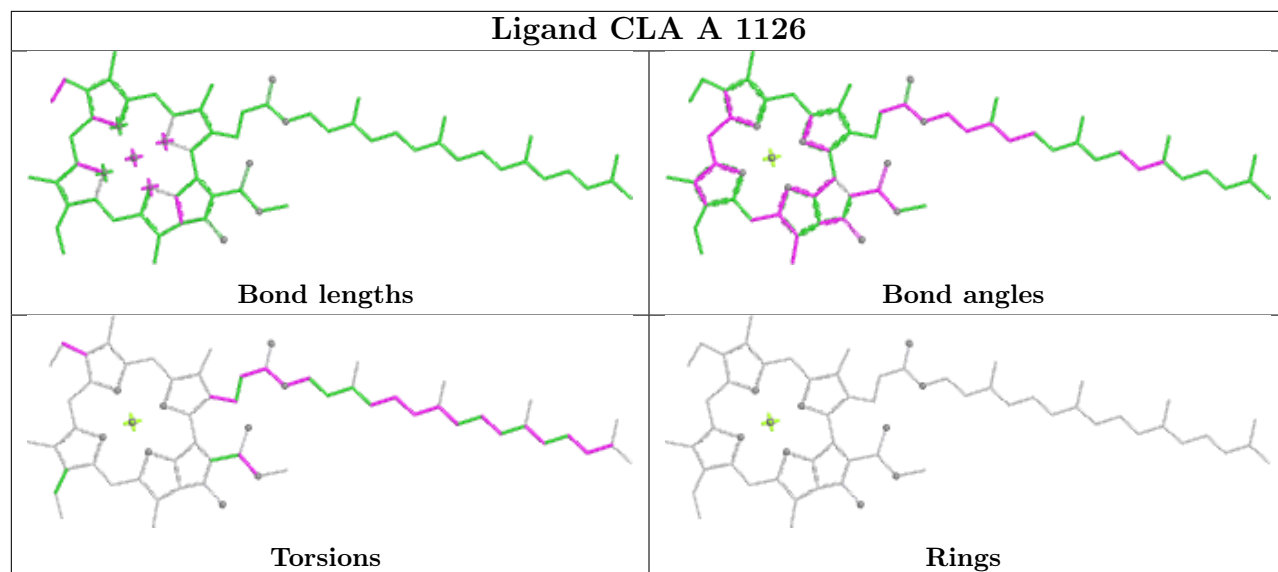
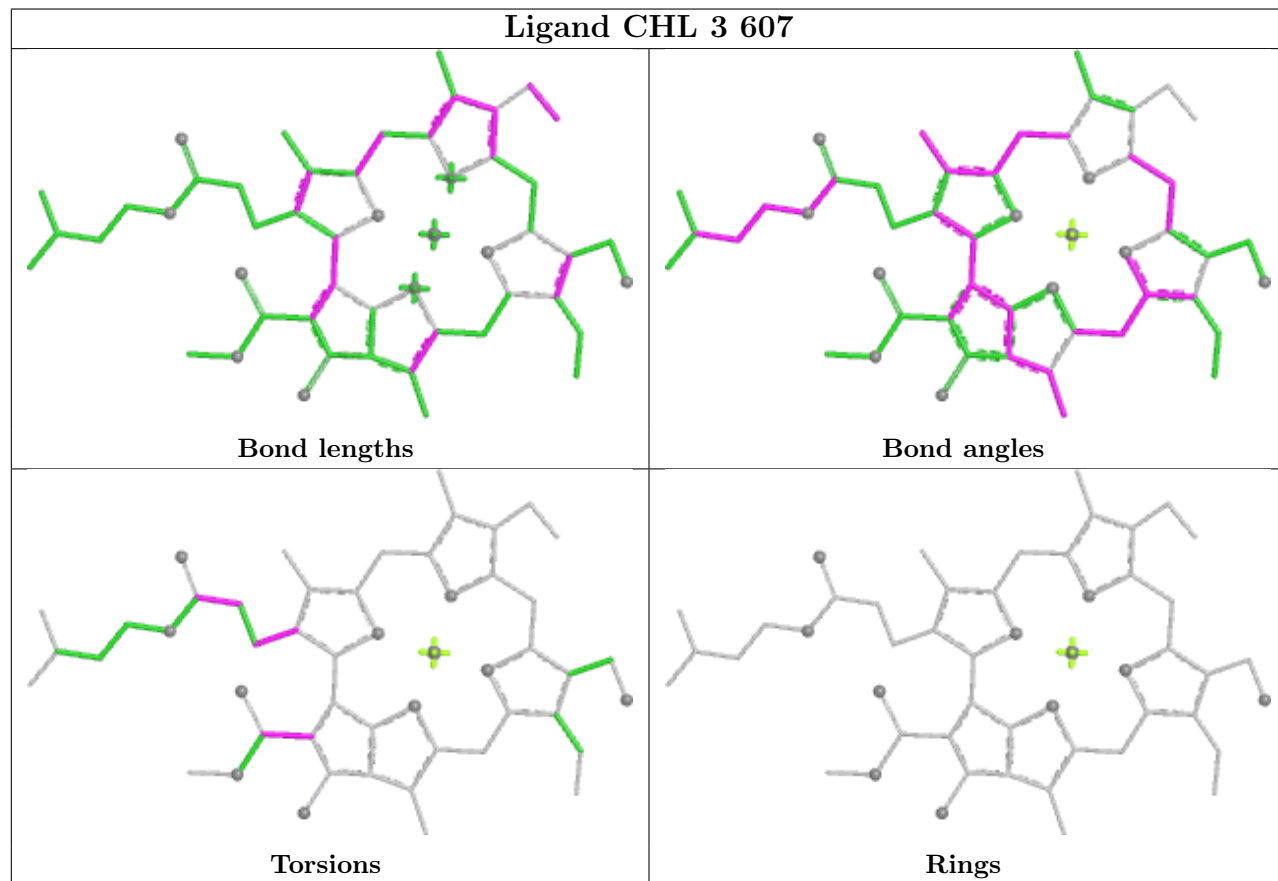


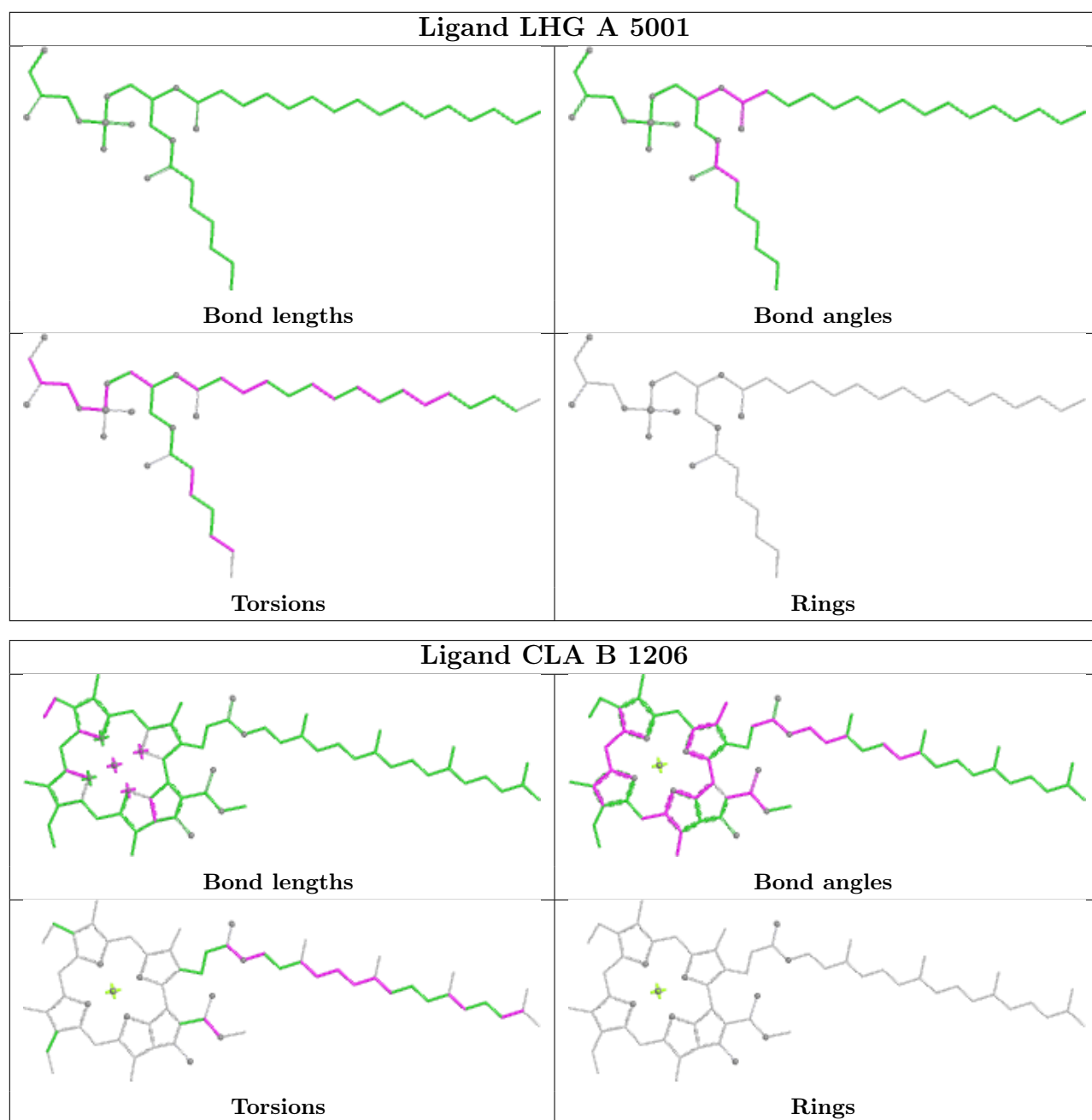
Ligand CLA 1 605

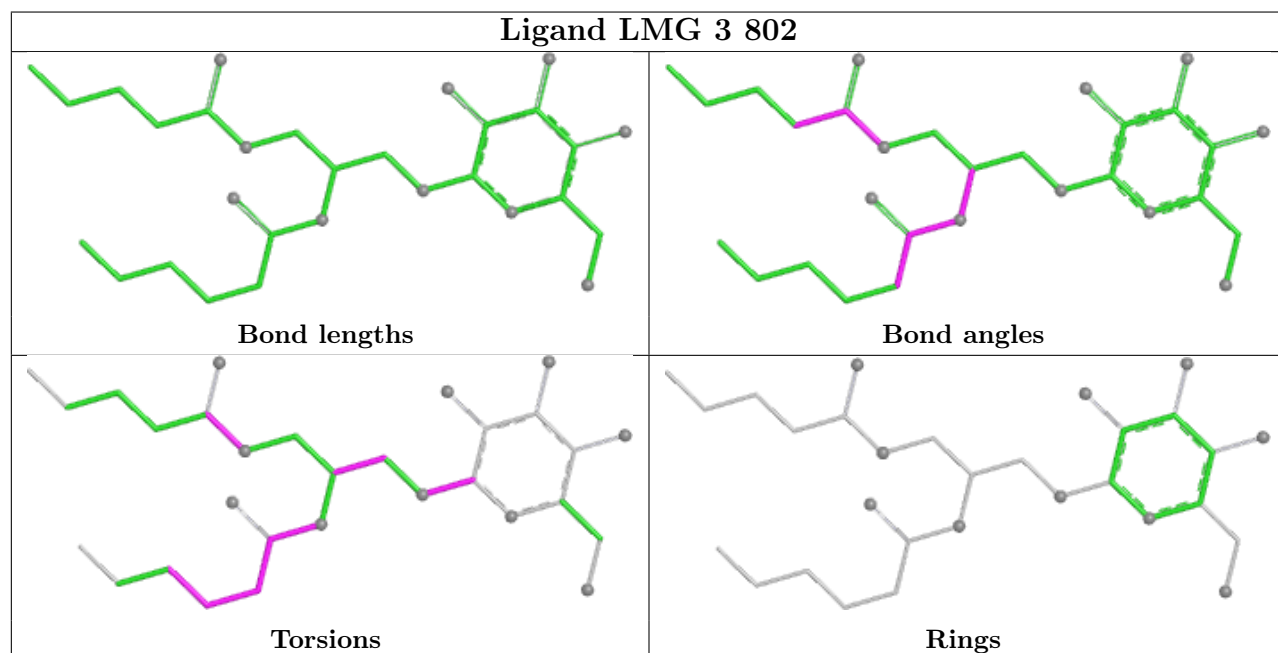
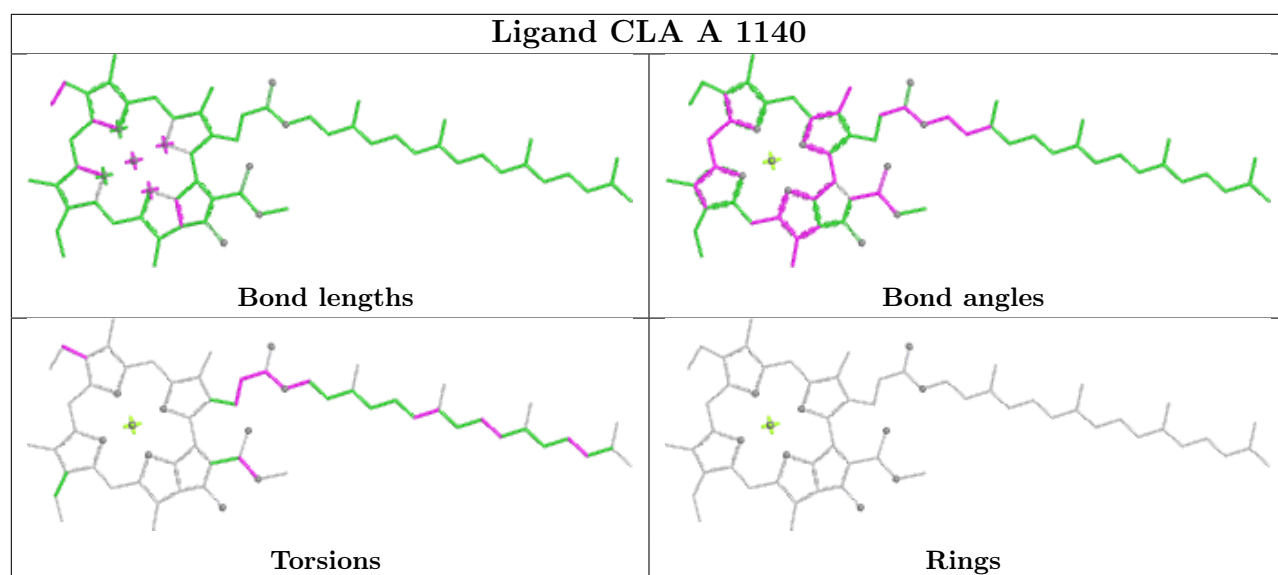


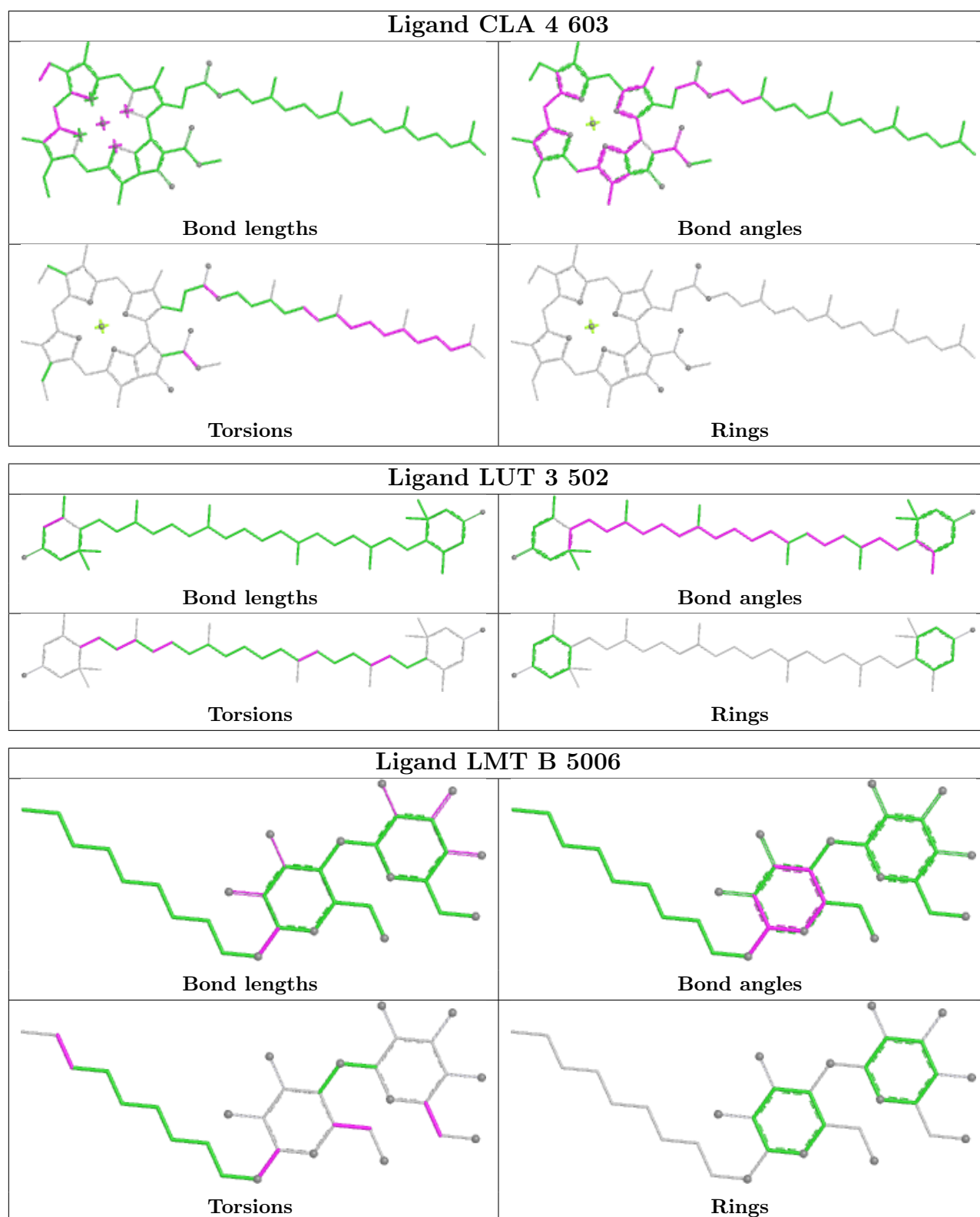


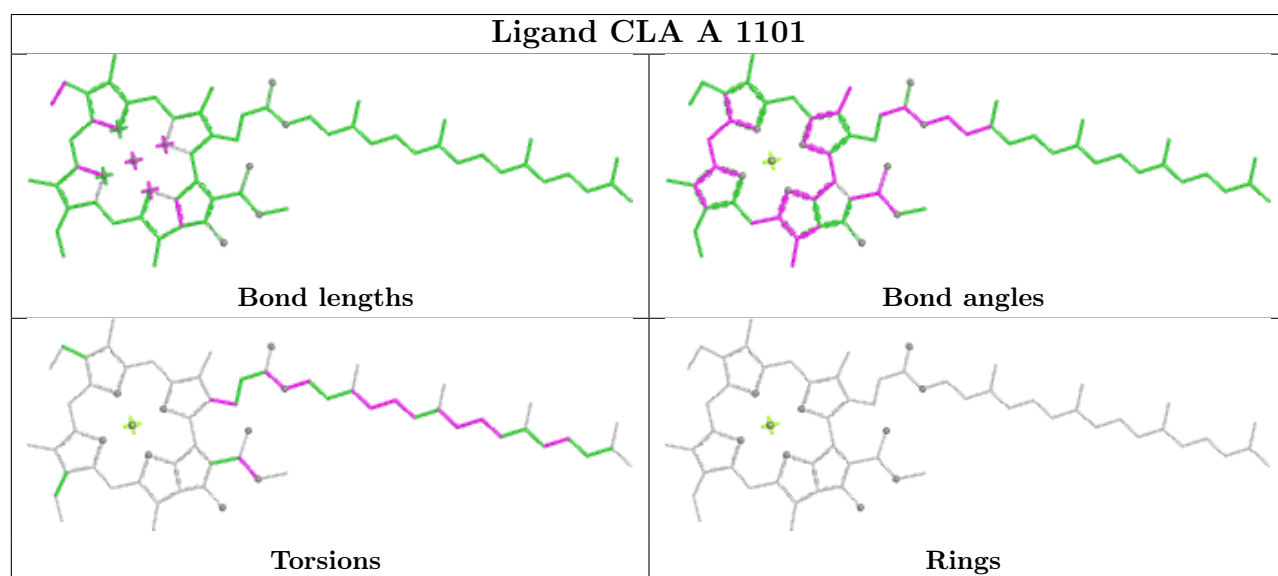
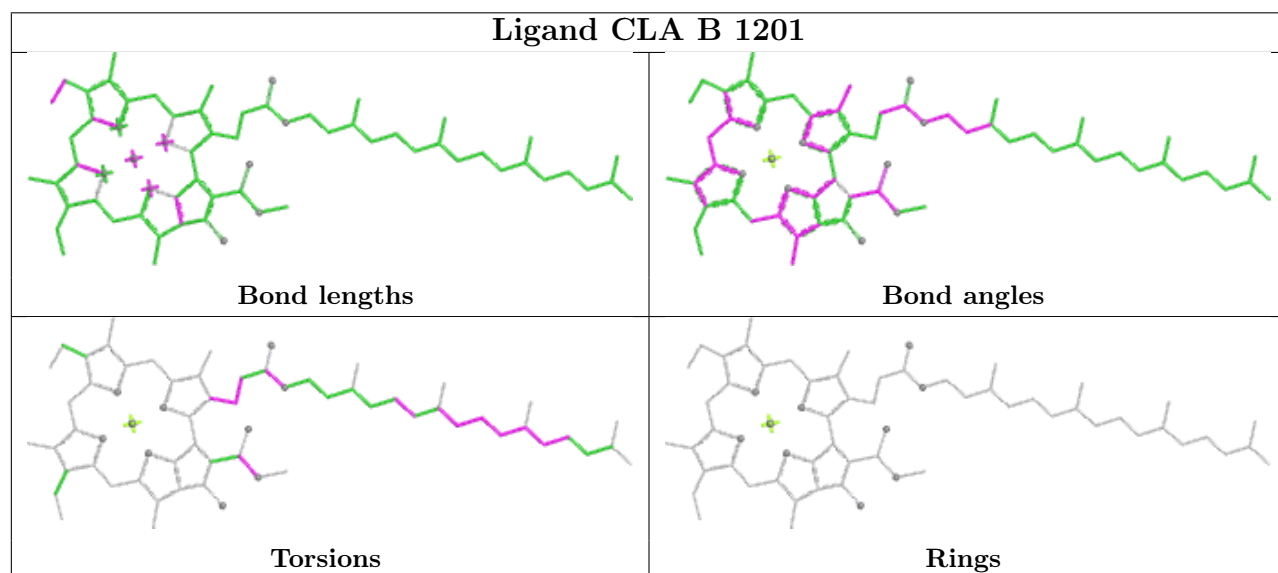
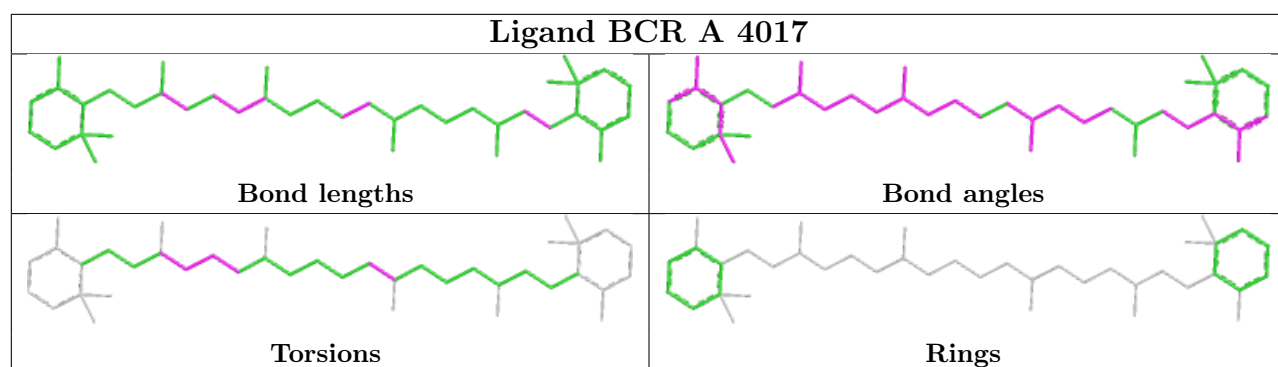


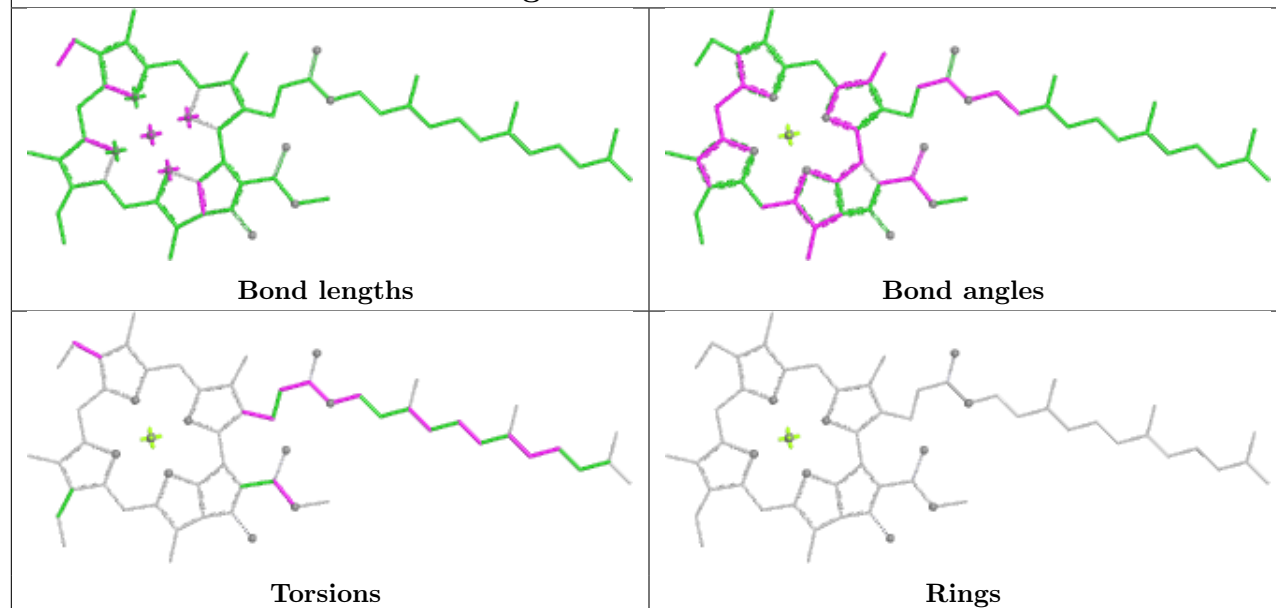
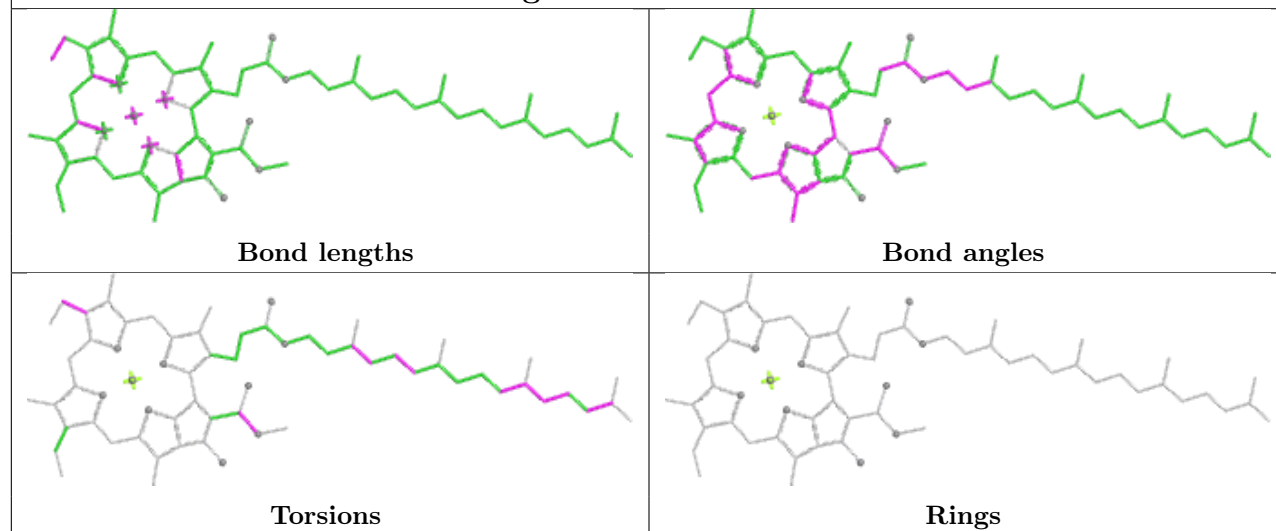
Ligand CLA A 1126**Ligand CHL 3 607**

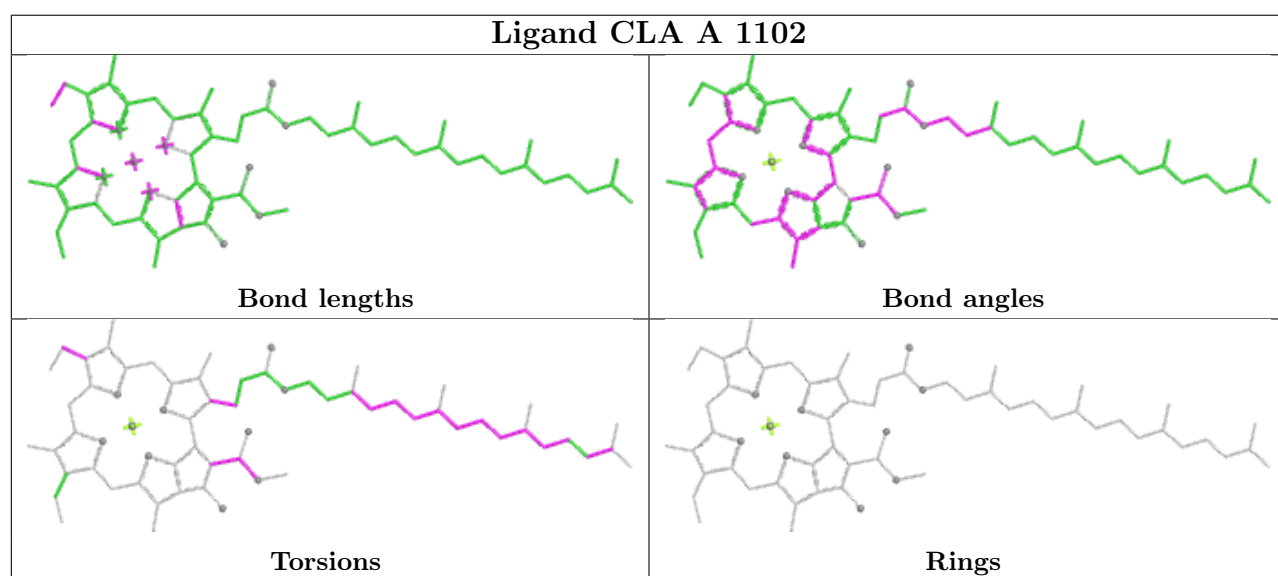
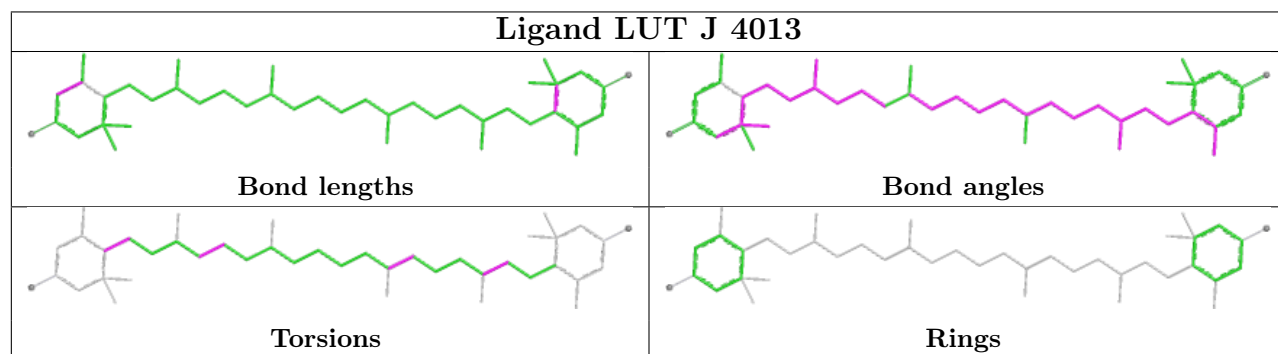
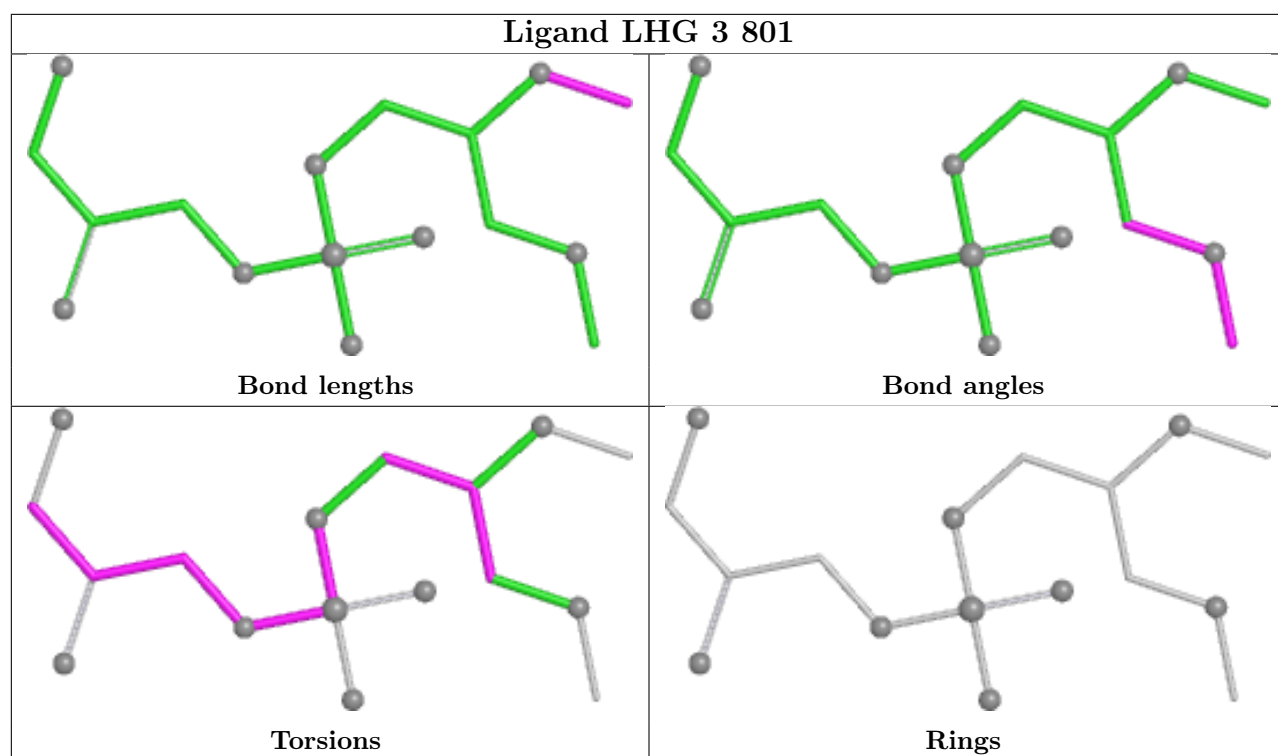




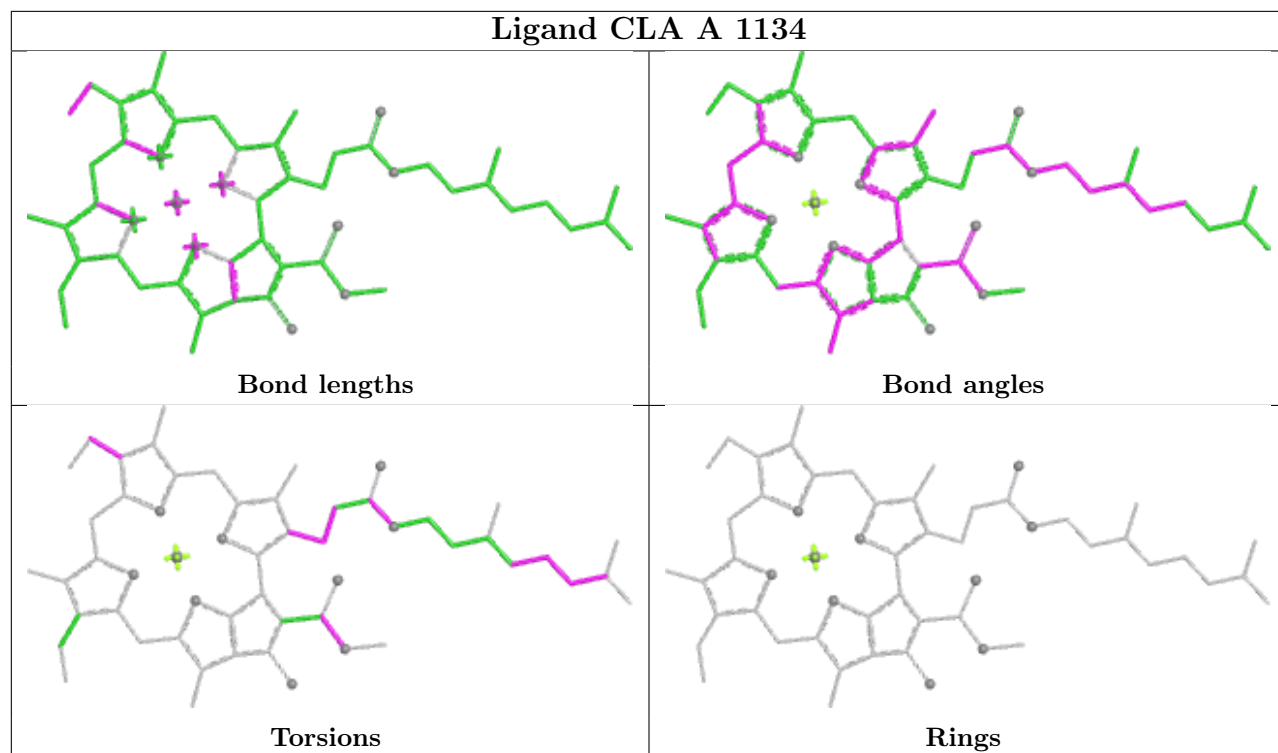




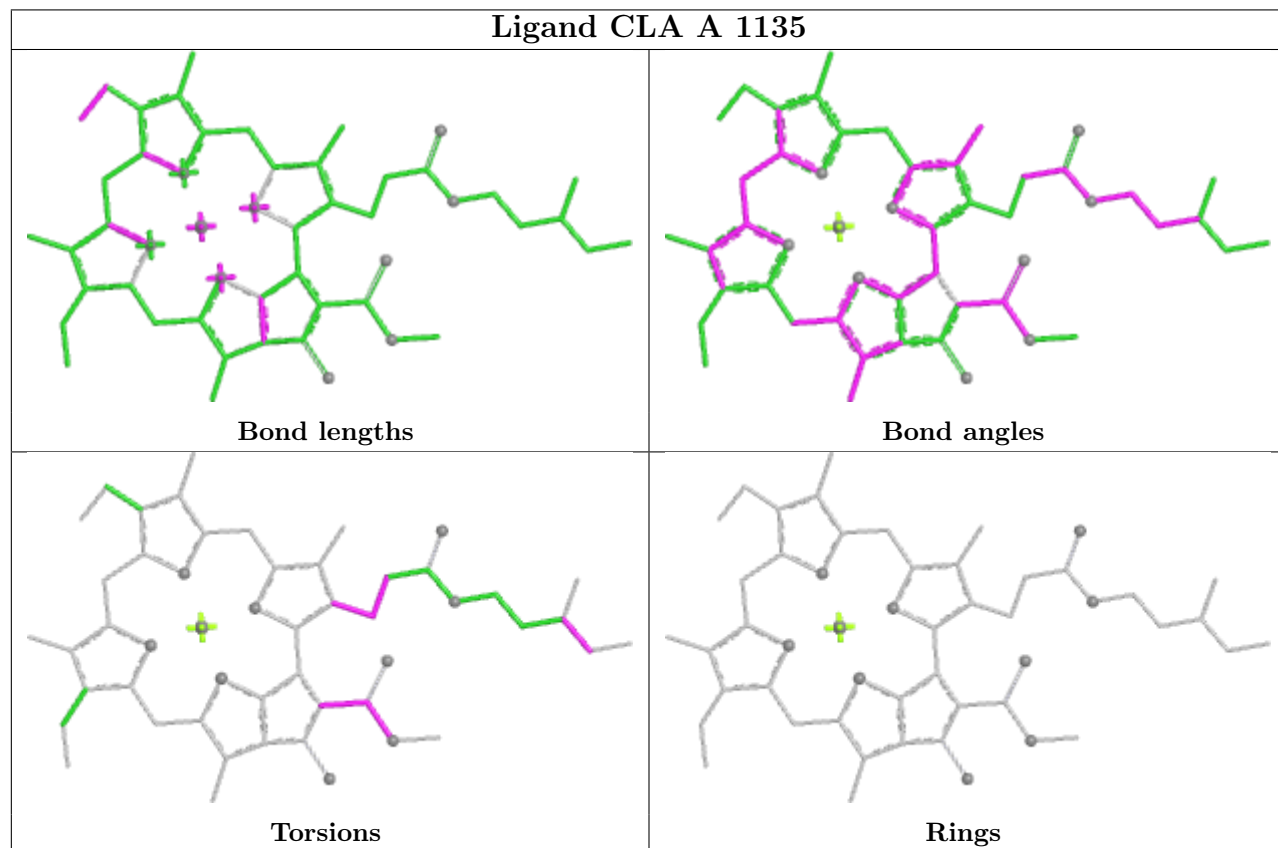
Ligand CLA A 1120**Ligand CLA A 1131**

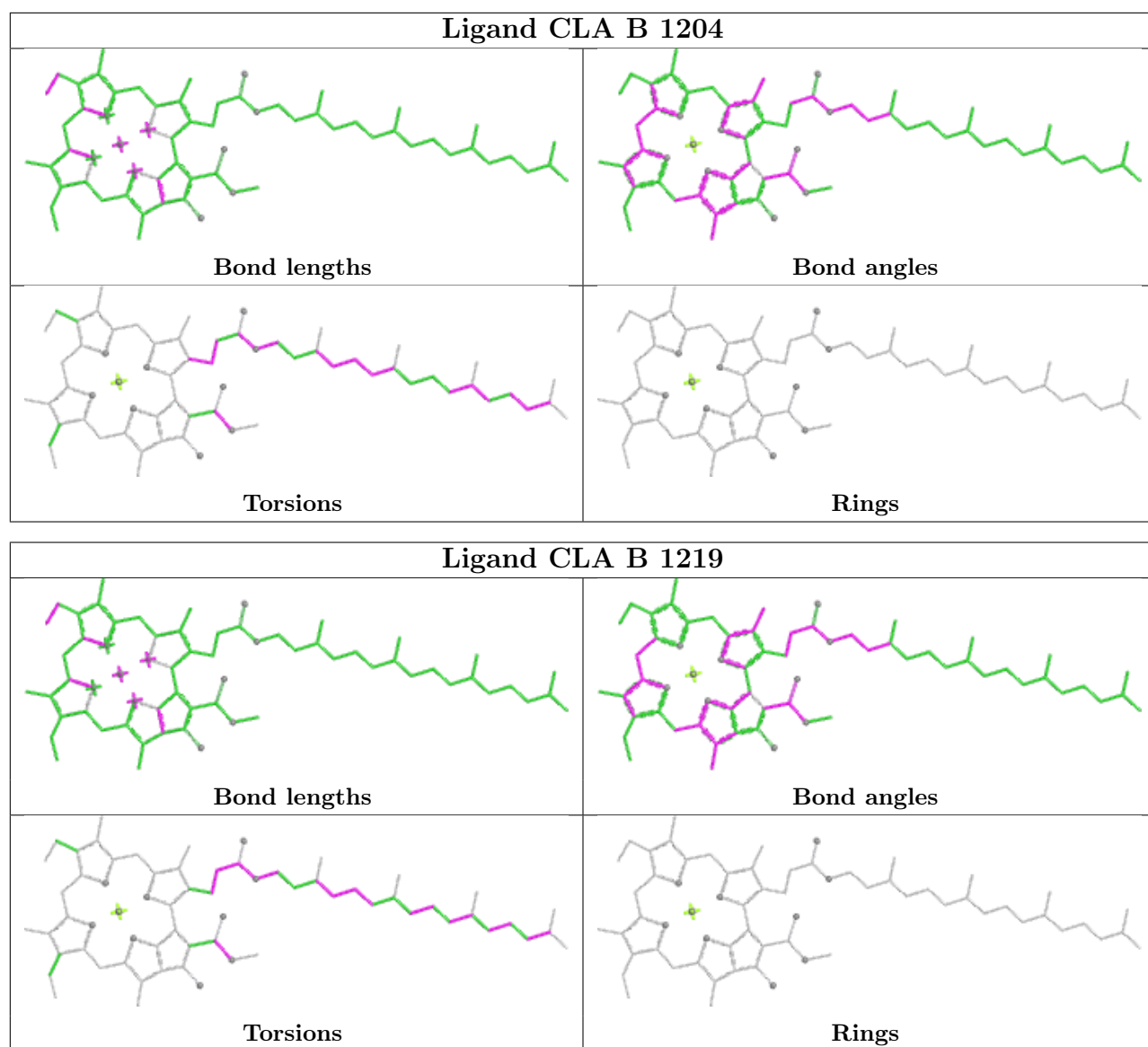


Ligand CLA A 1134

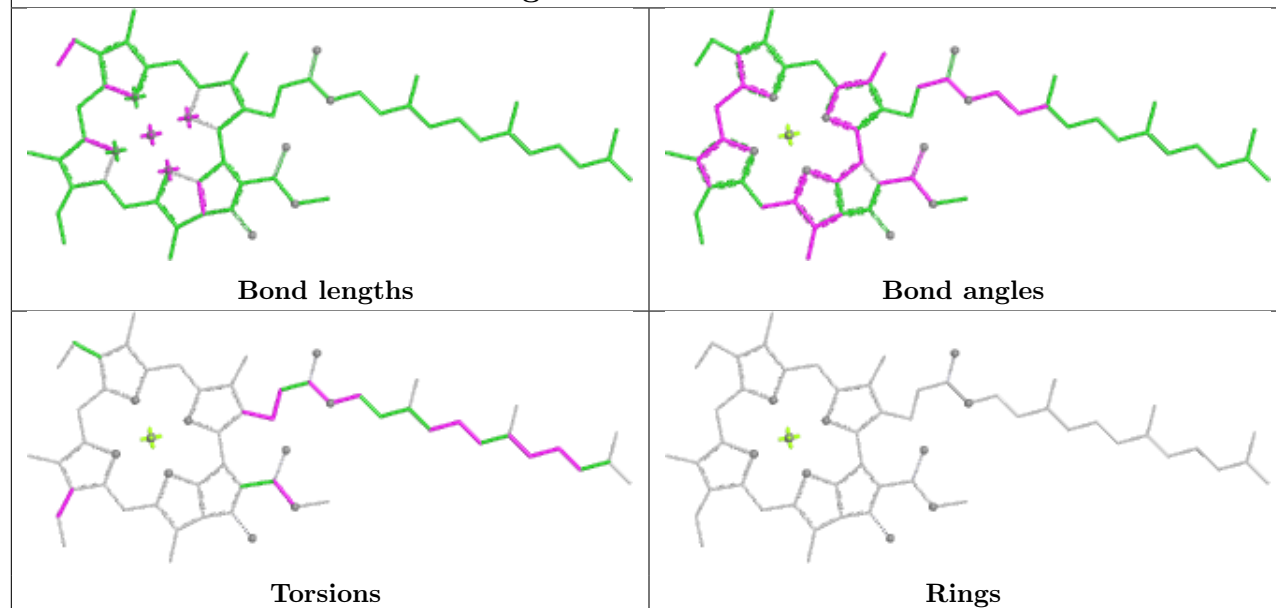


Ligand CLA A 1135

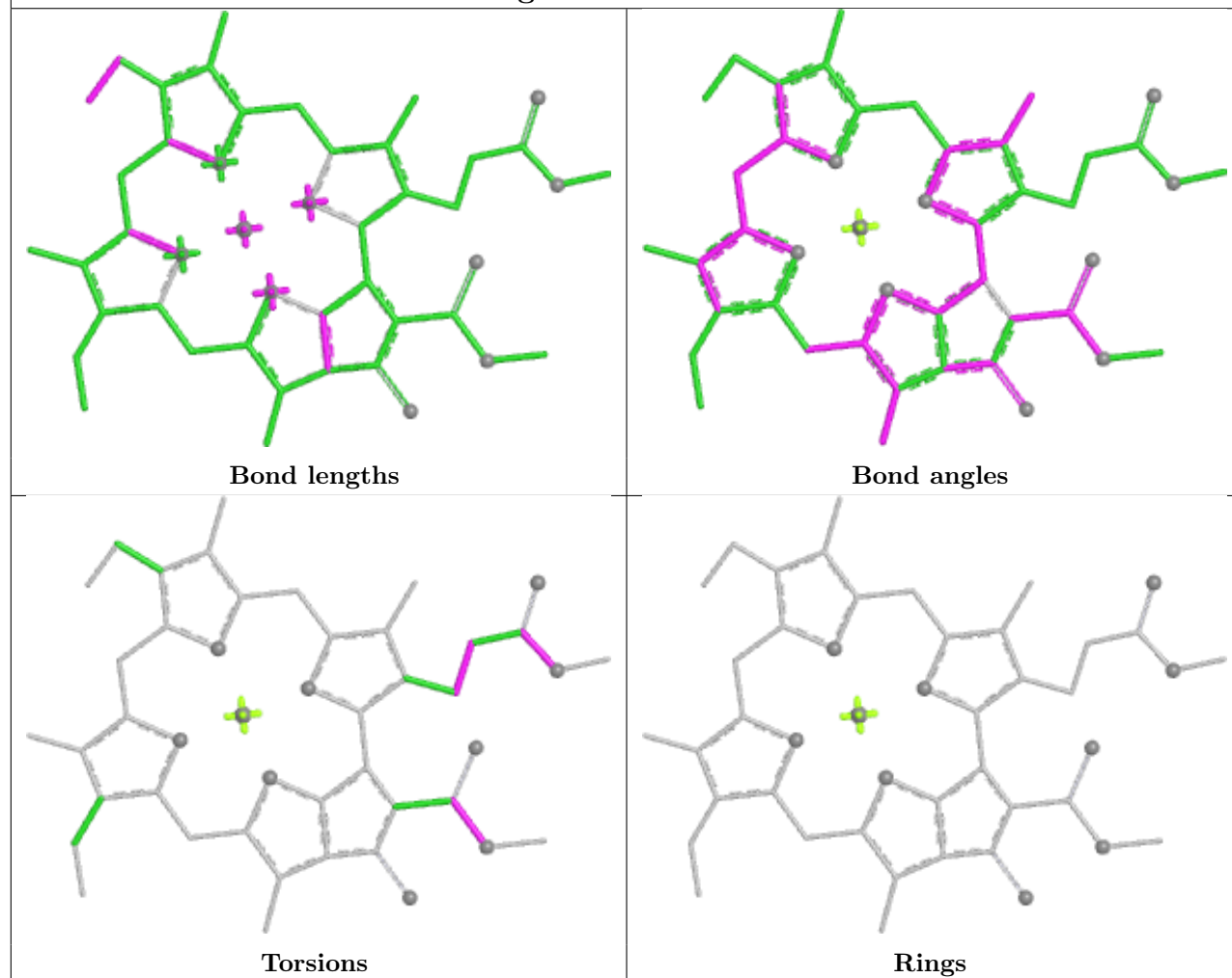




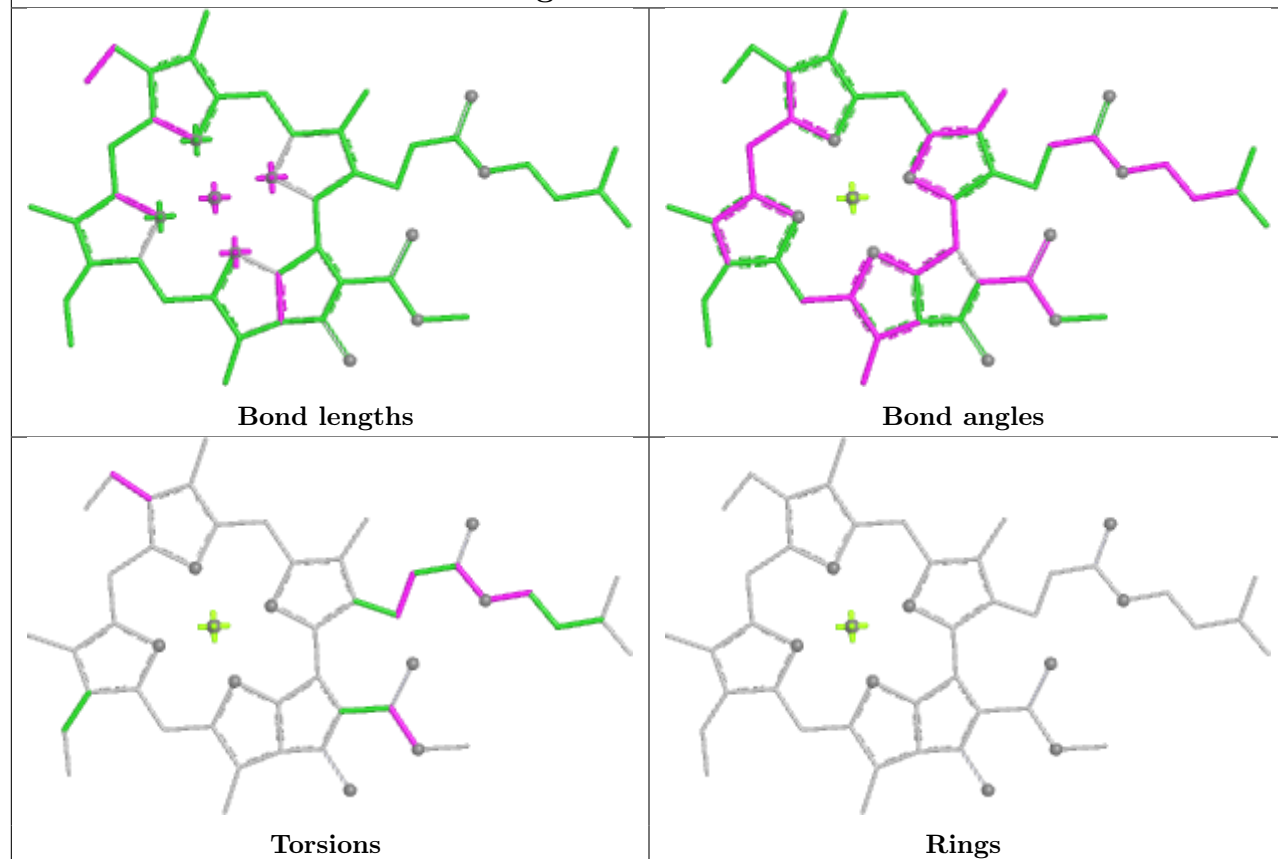
Ligand CLA H 1701



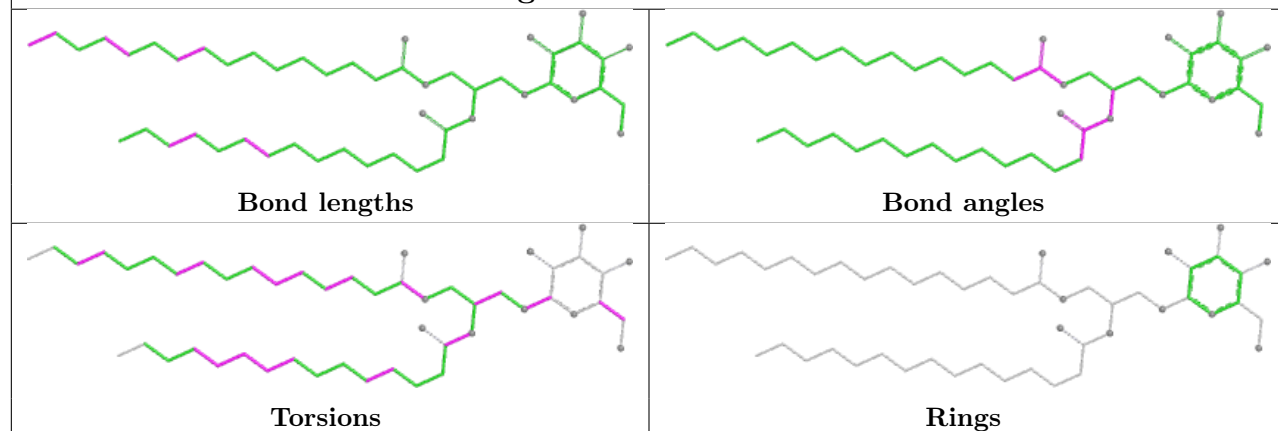
Ligand CLA 1 607

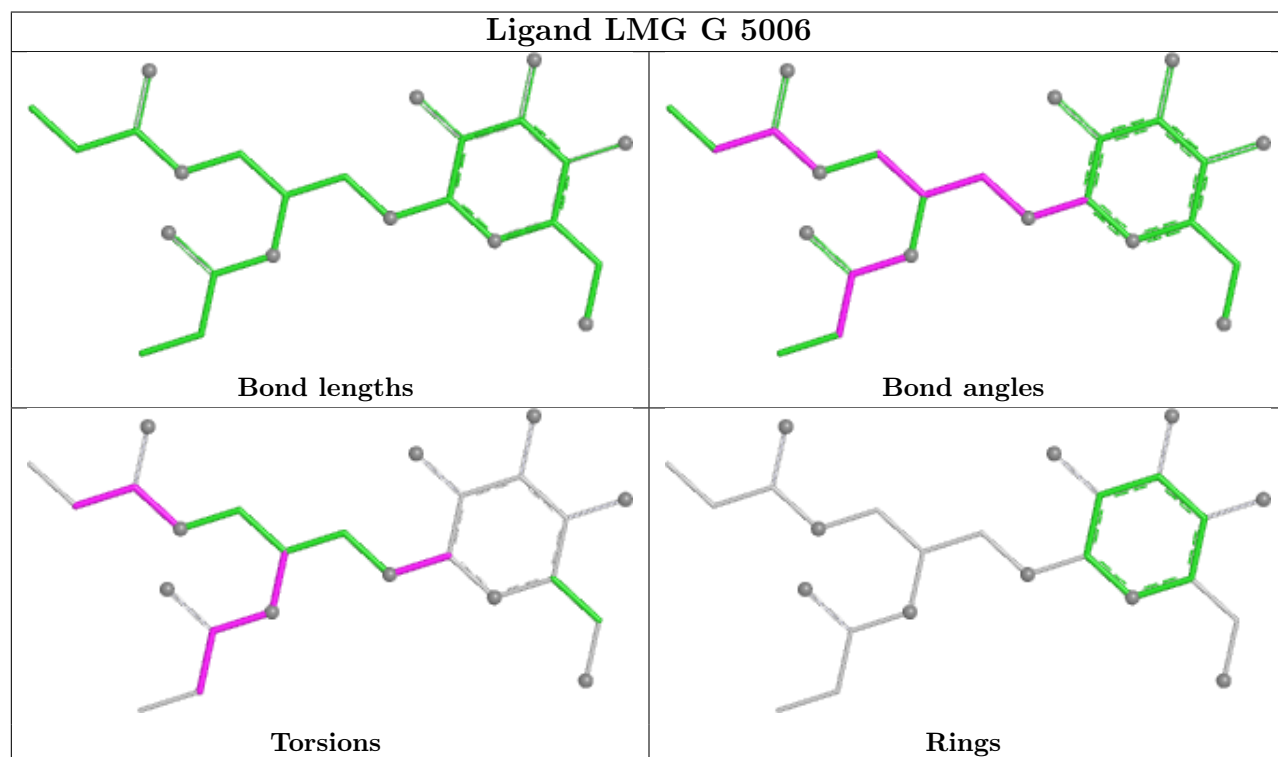
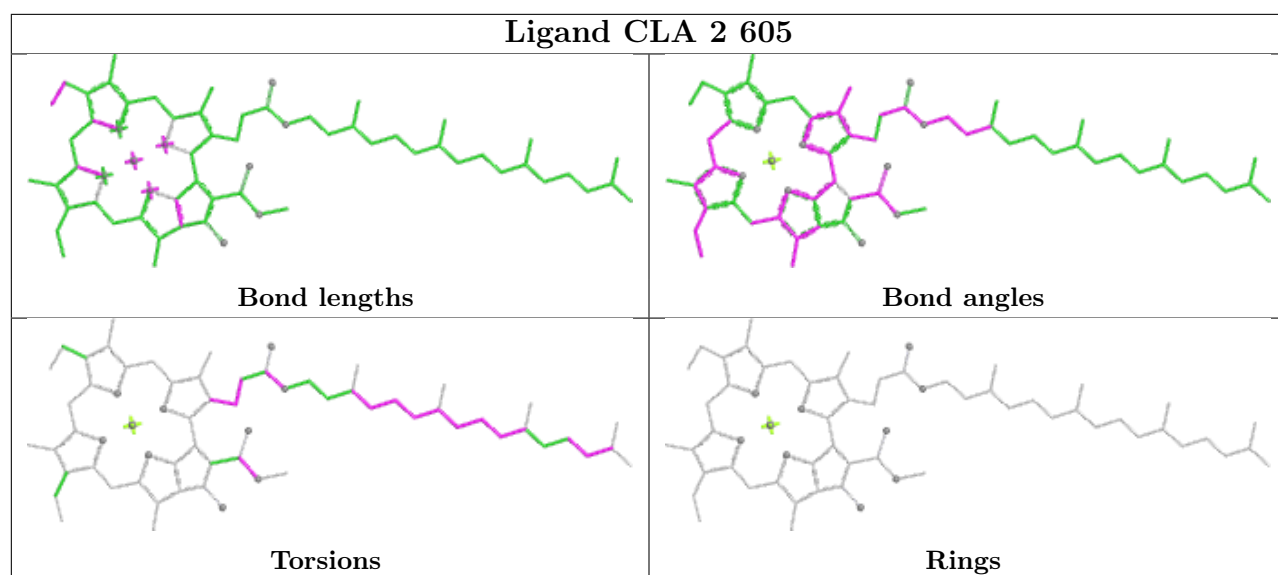


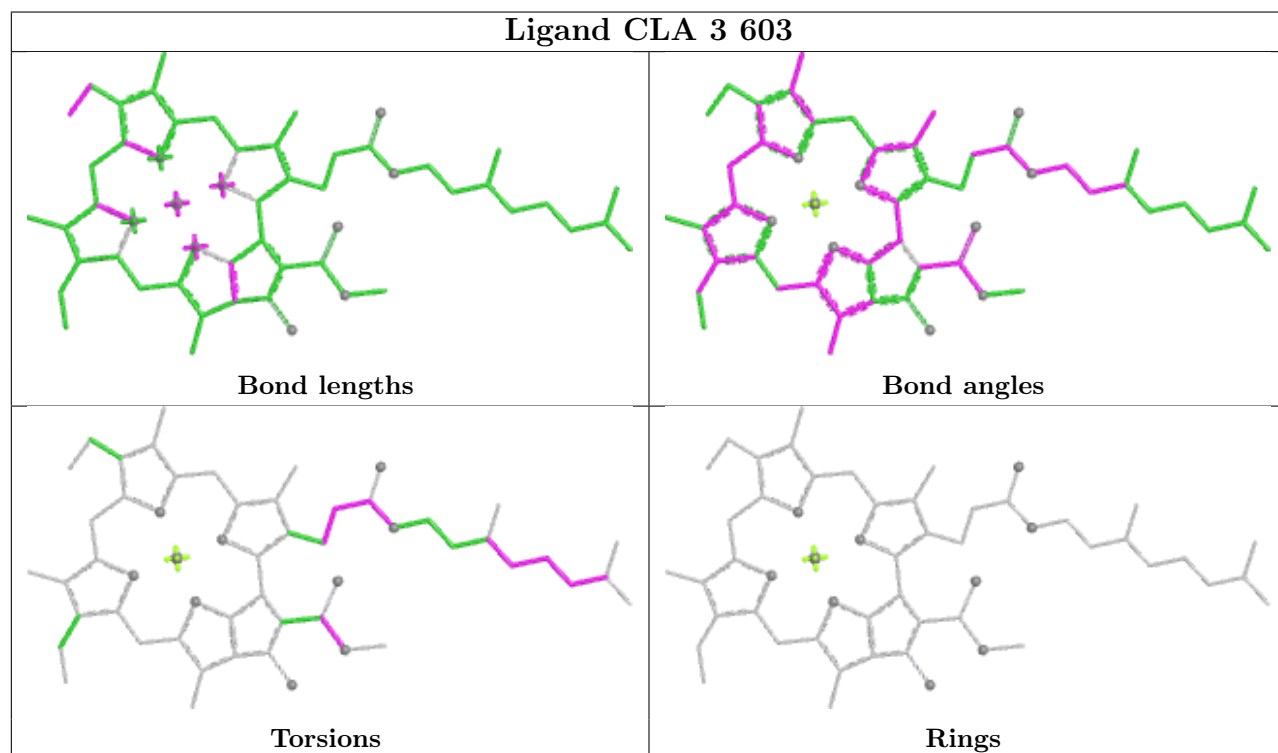
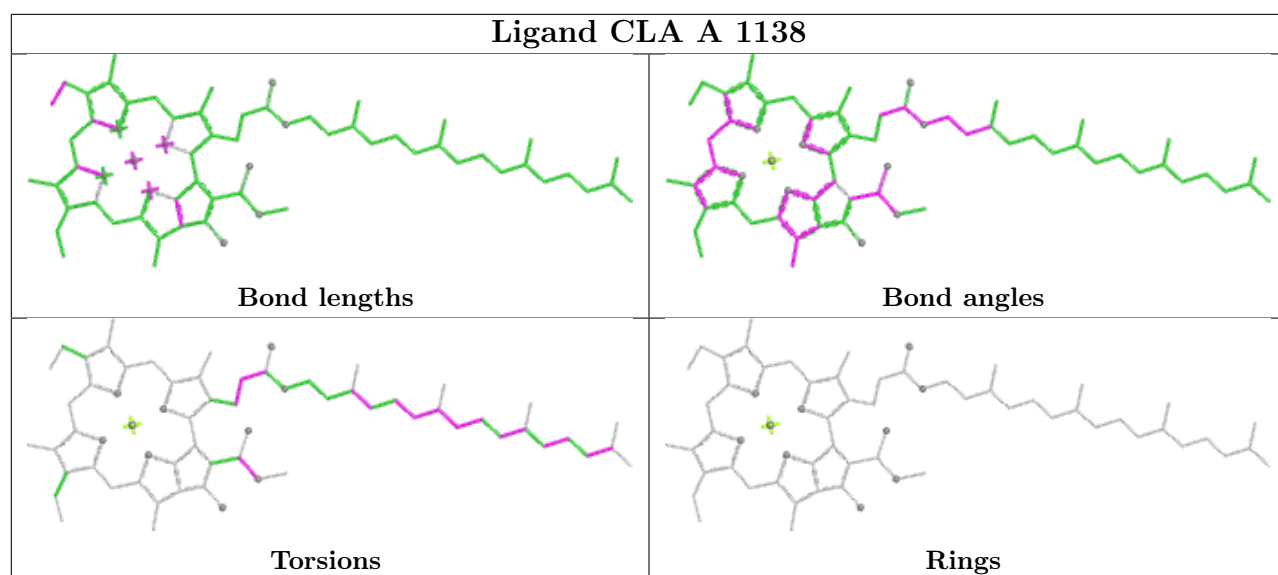
Ligand CLA 2 608

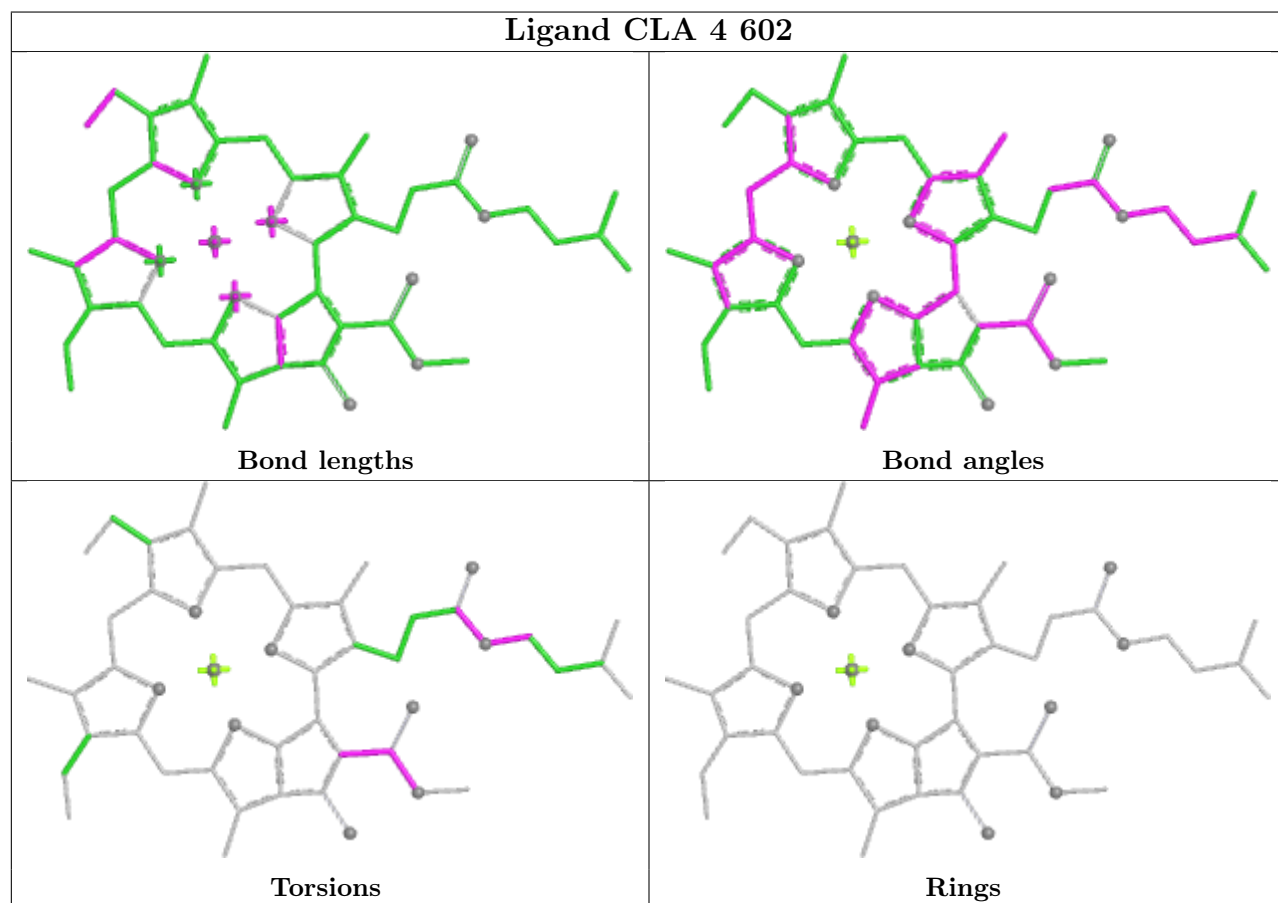
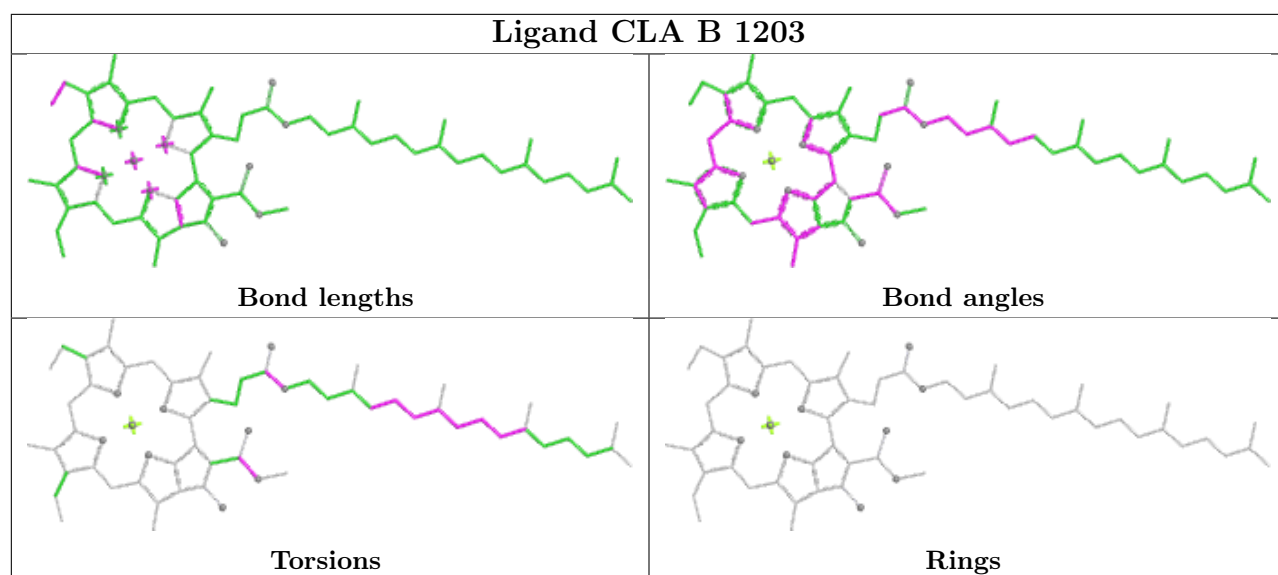


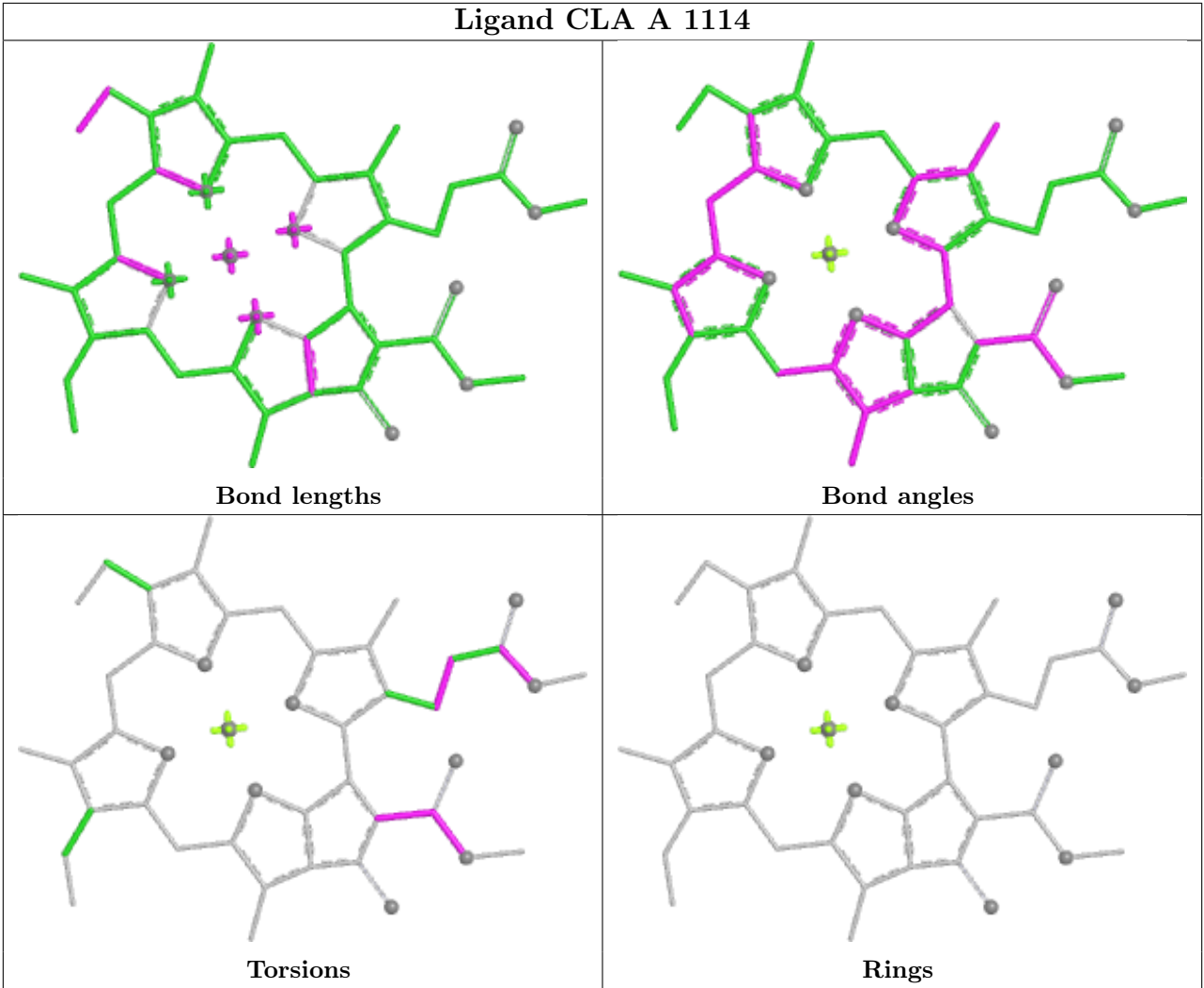
Ligand LMG G 5002











5.7 Other polymers ⓘ

There are no such residues in this entry.

5.8 Polymer linkage issues ⓘ

The following chains have linkage breaks:

Mol	Chain	Number of breaks
17	N	1

All chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	N	35:LEU	C	36:PRO	N	1.19

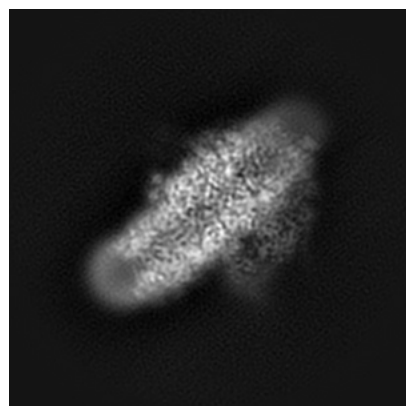
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-10798. These allow visual inspection of the internal detail of the map and identification of artifacts.

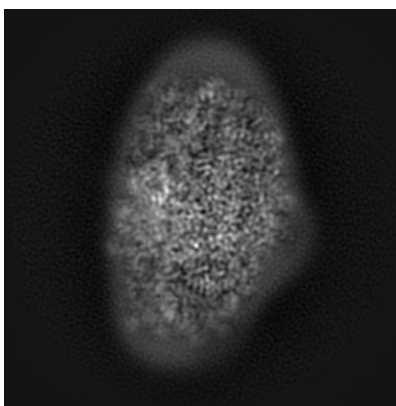
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

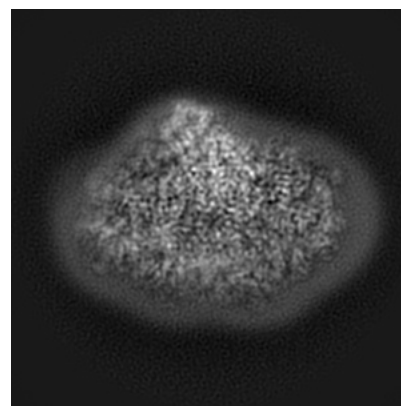
6.1.1 Primary map



X

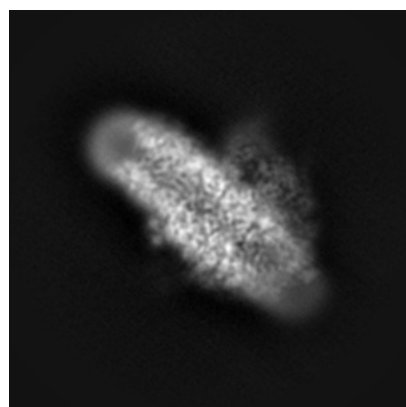


Y

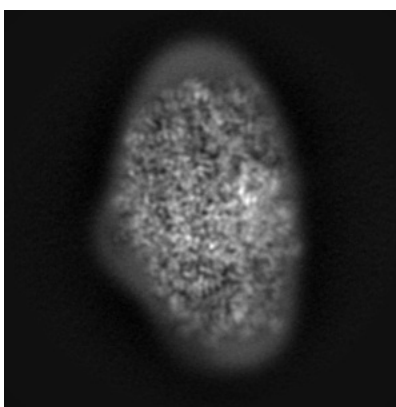


Z

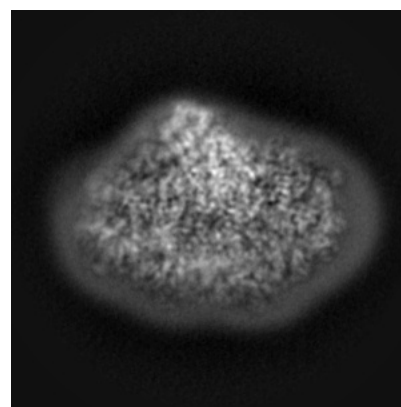
6.1.2 Raw map



X



Y

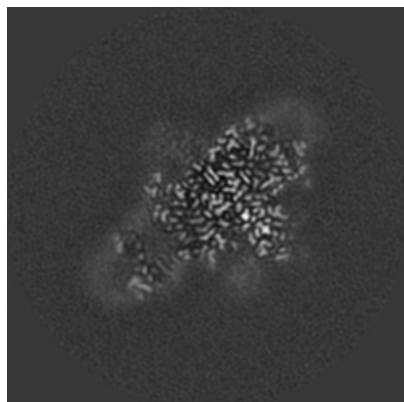


Z

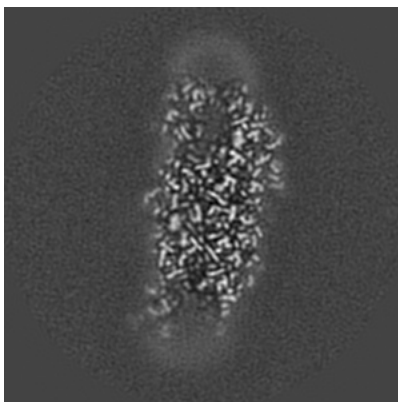
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

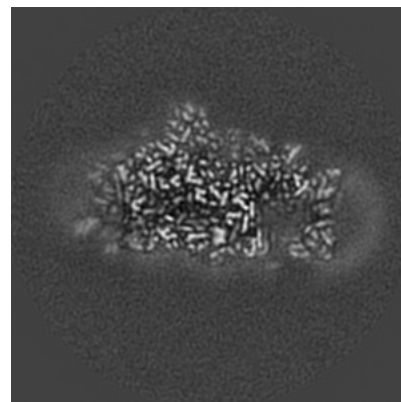
6.2.1 Primary map



X Index: 150

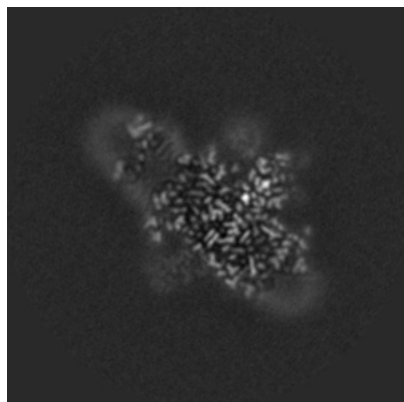


Y Index: 150

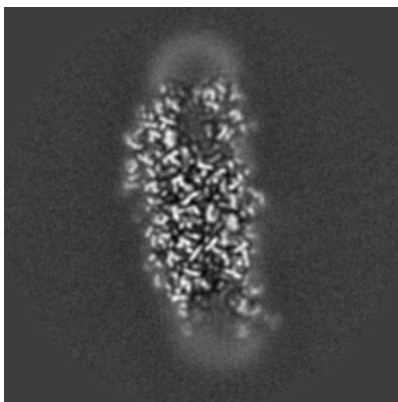


Z Index: 150

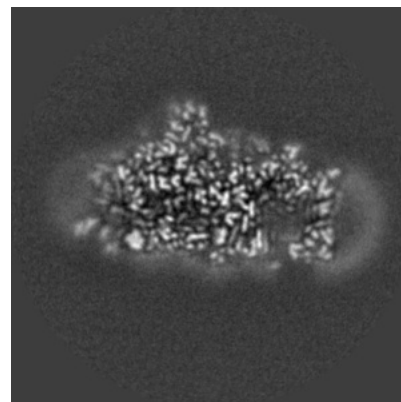
6.2.2 Raw map



X Index: 150



Y Index: 150

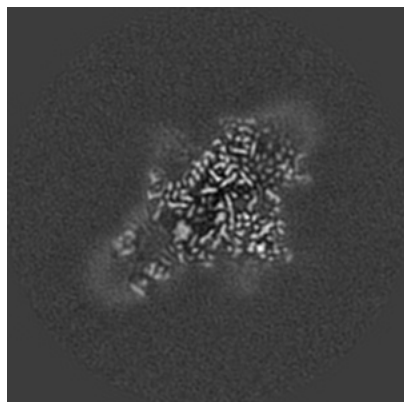


Z Index: 150

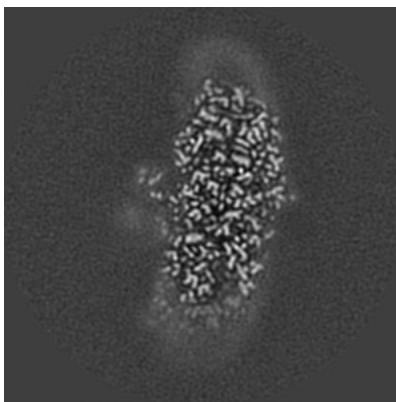
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

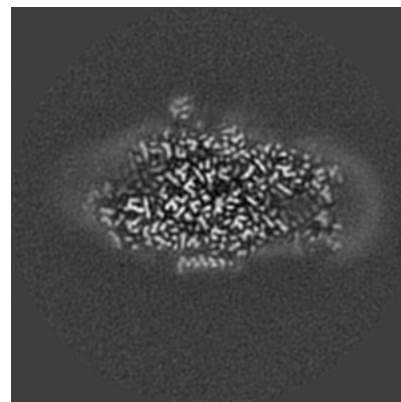
6.3.1 Primary map



X Index: 154

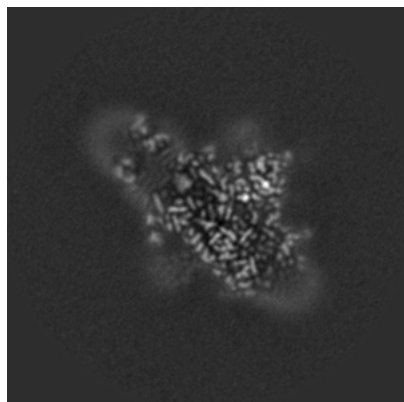


Y Index: 164

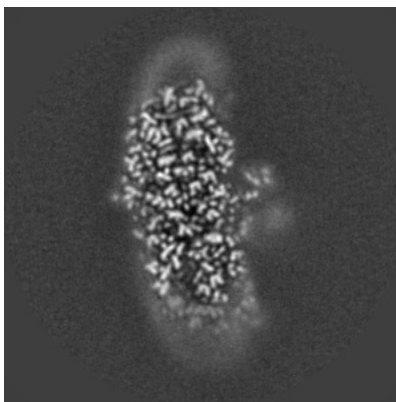


Z Index: 161

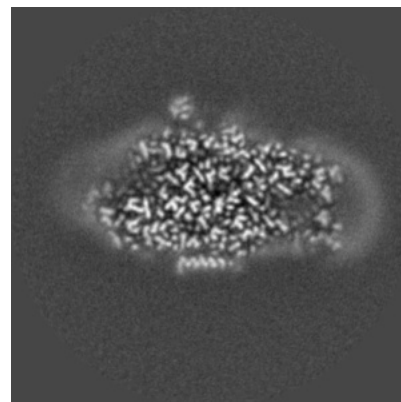
6.3.2 Raw map



X Index: 153



Y Index: 164

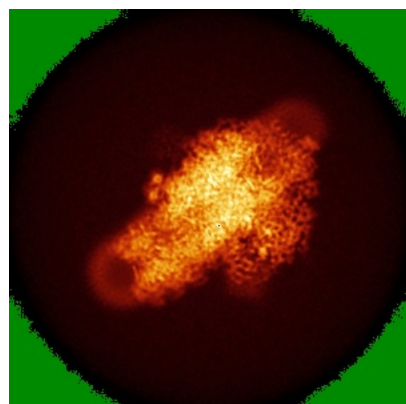


Z Index: 138

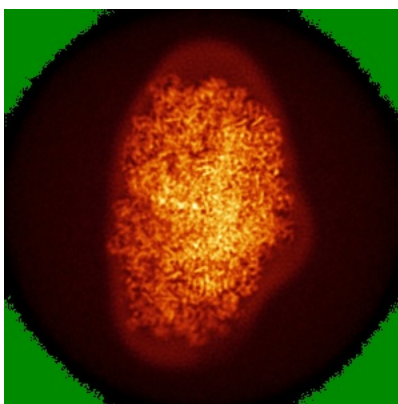
The images above show the largest variance slices of the map in three orthogonal directions.

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

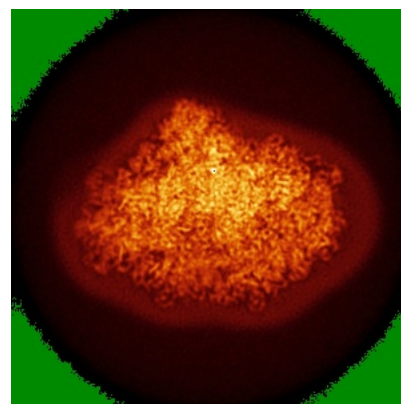
6.4.1 Primary map



X

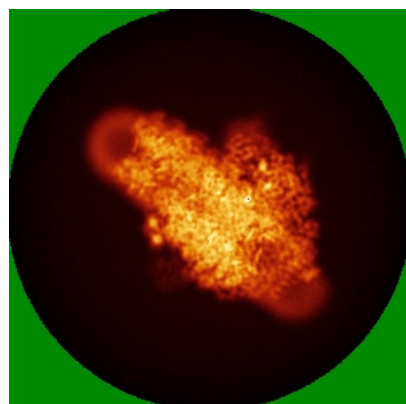


Y

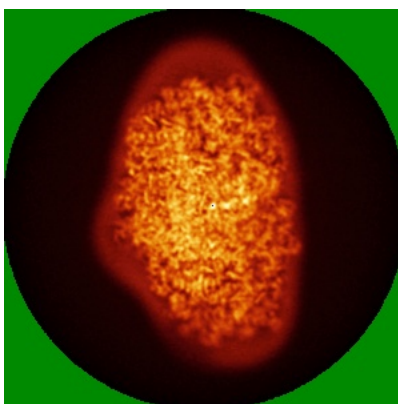


Z

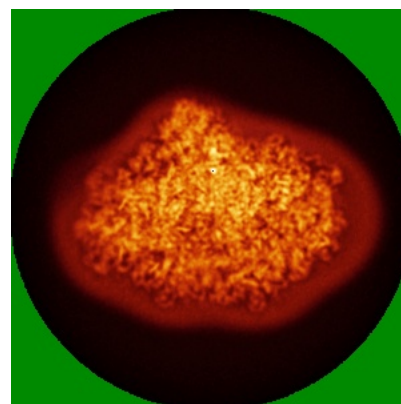
6.4.2 Raw map



X



Y

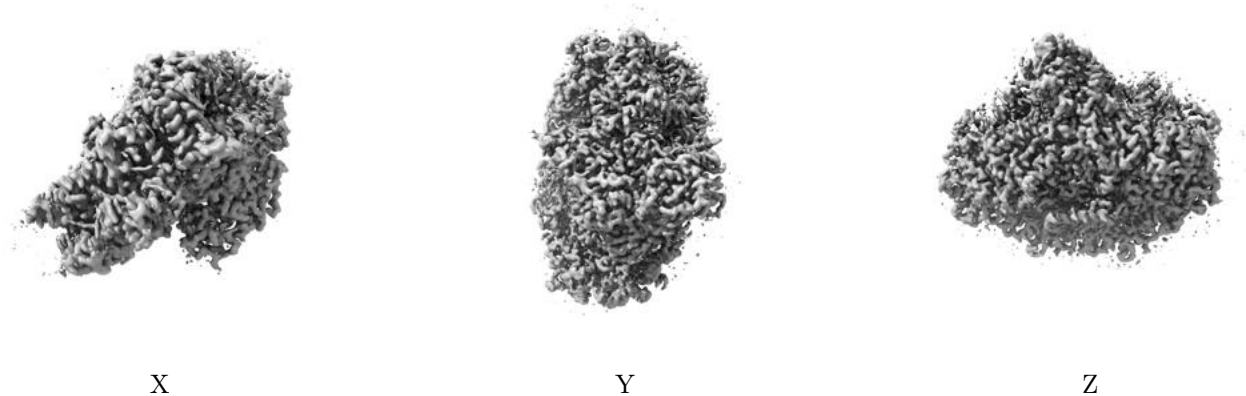


Z

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.

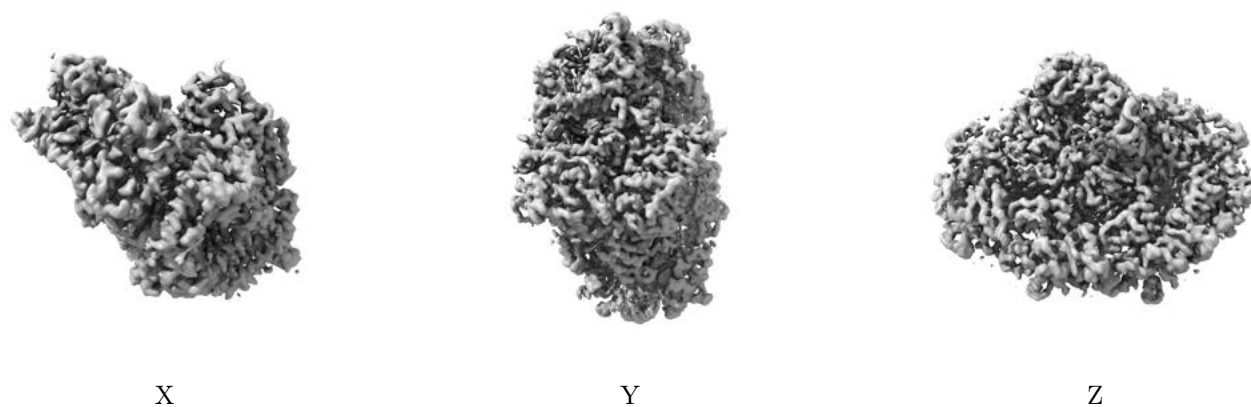
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.022. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

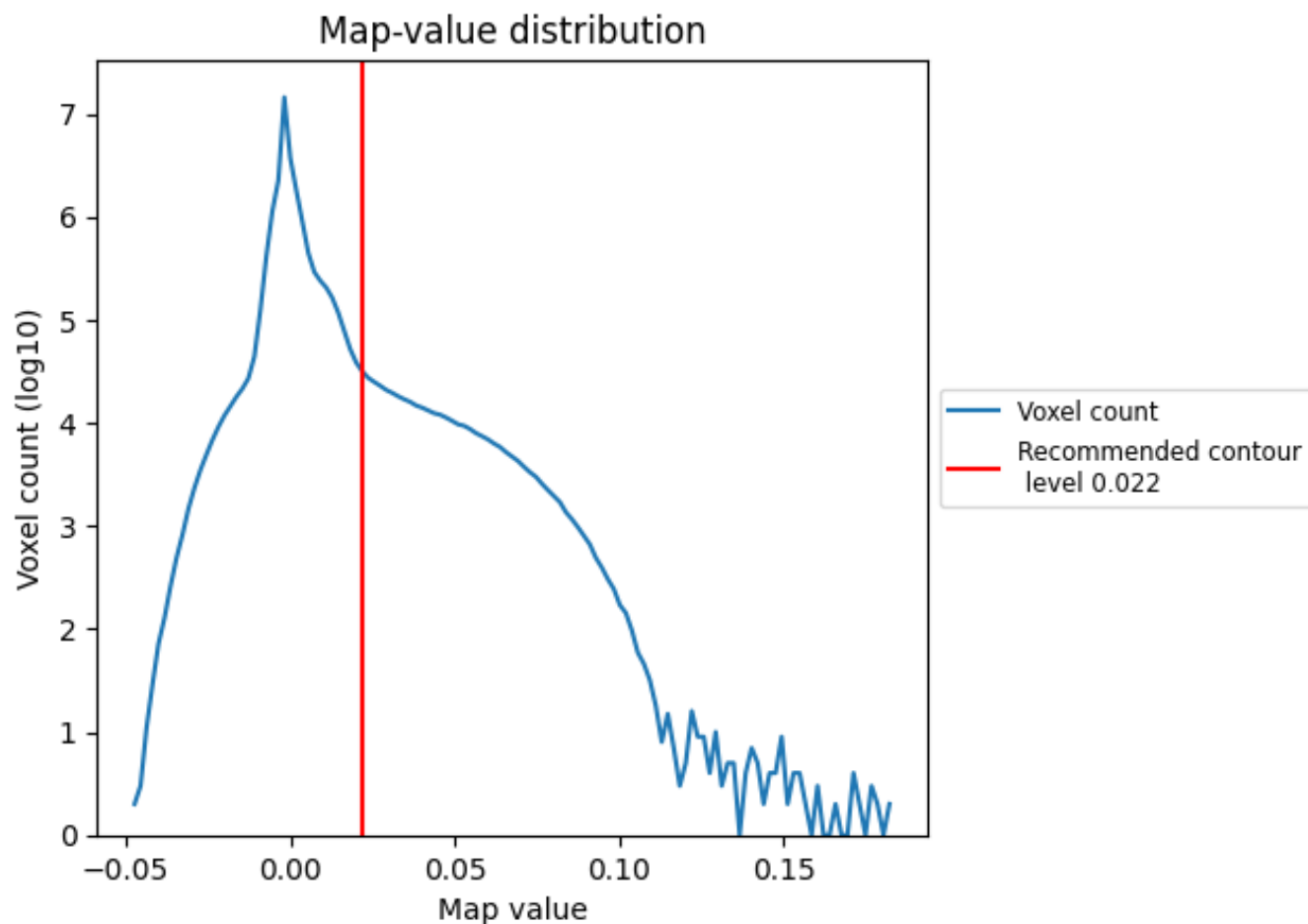
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

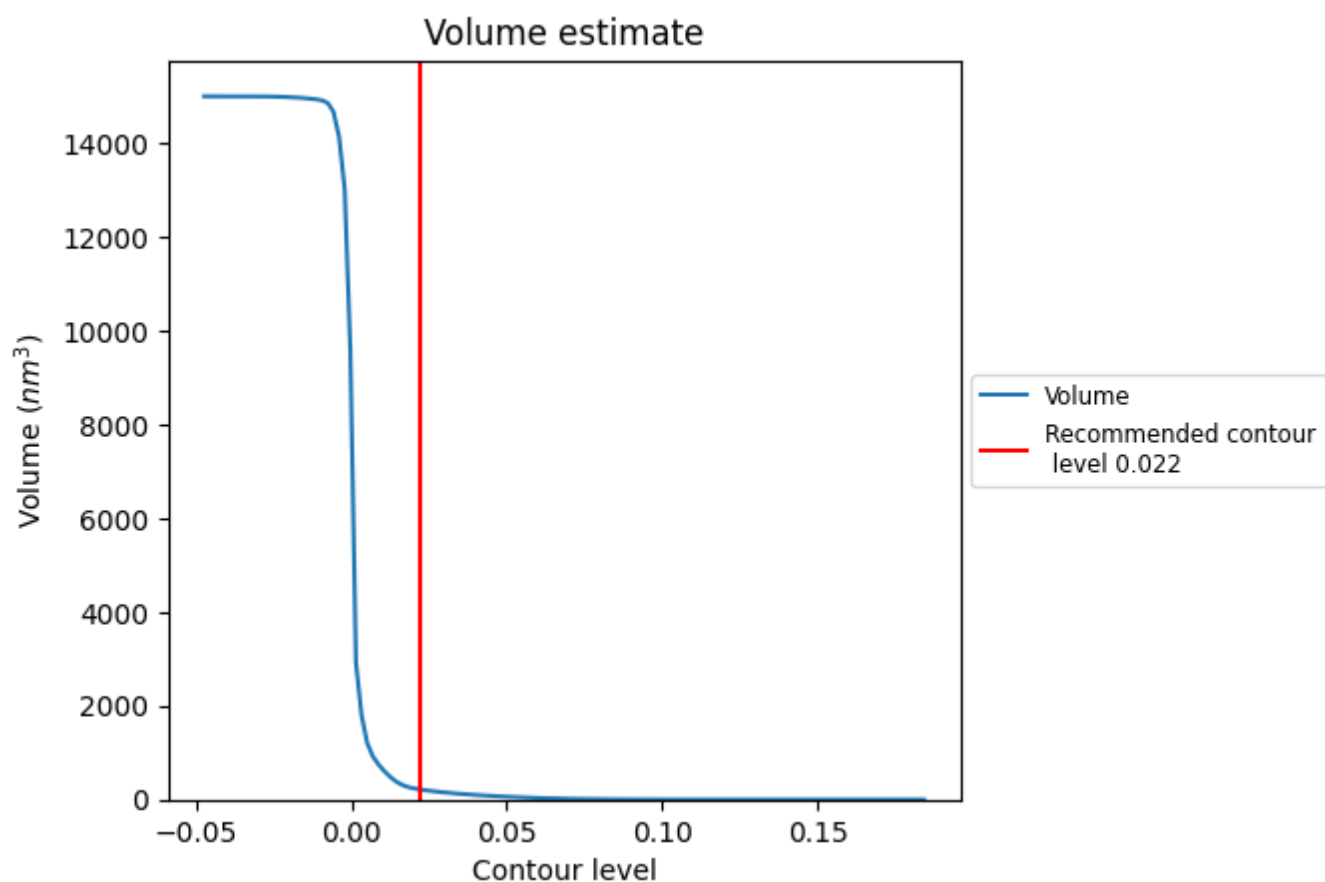
This section contains the results of statistical analysis of the map.

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.

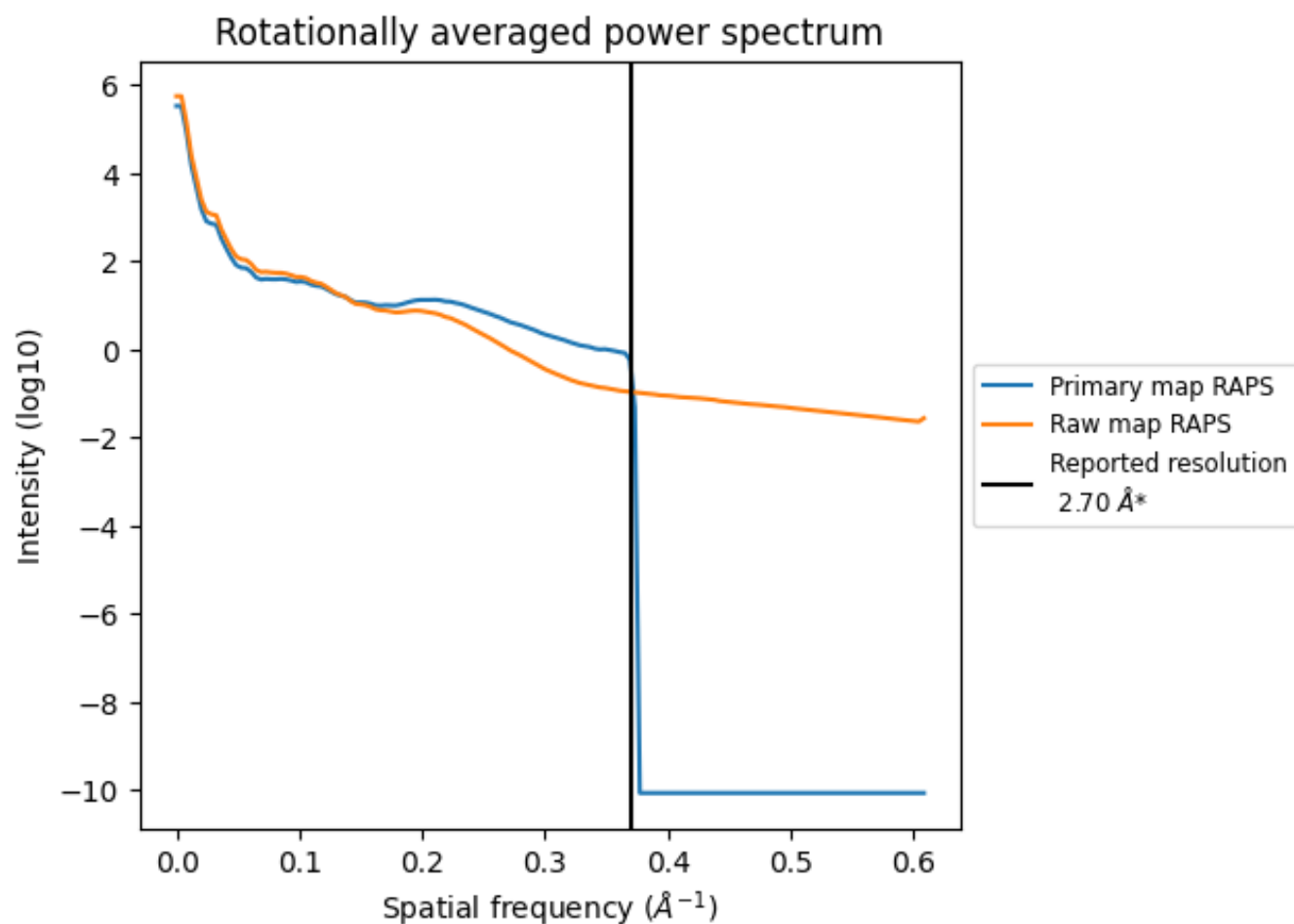
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 215 nm^3 ; this corresponds to an approximate mass of 195 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 ⓘ

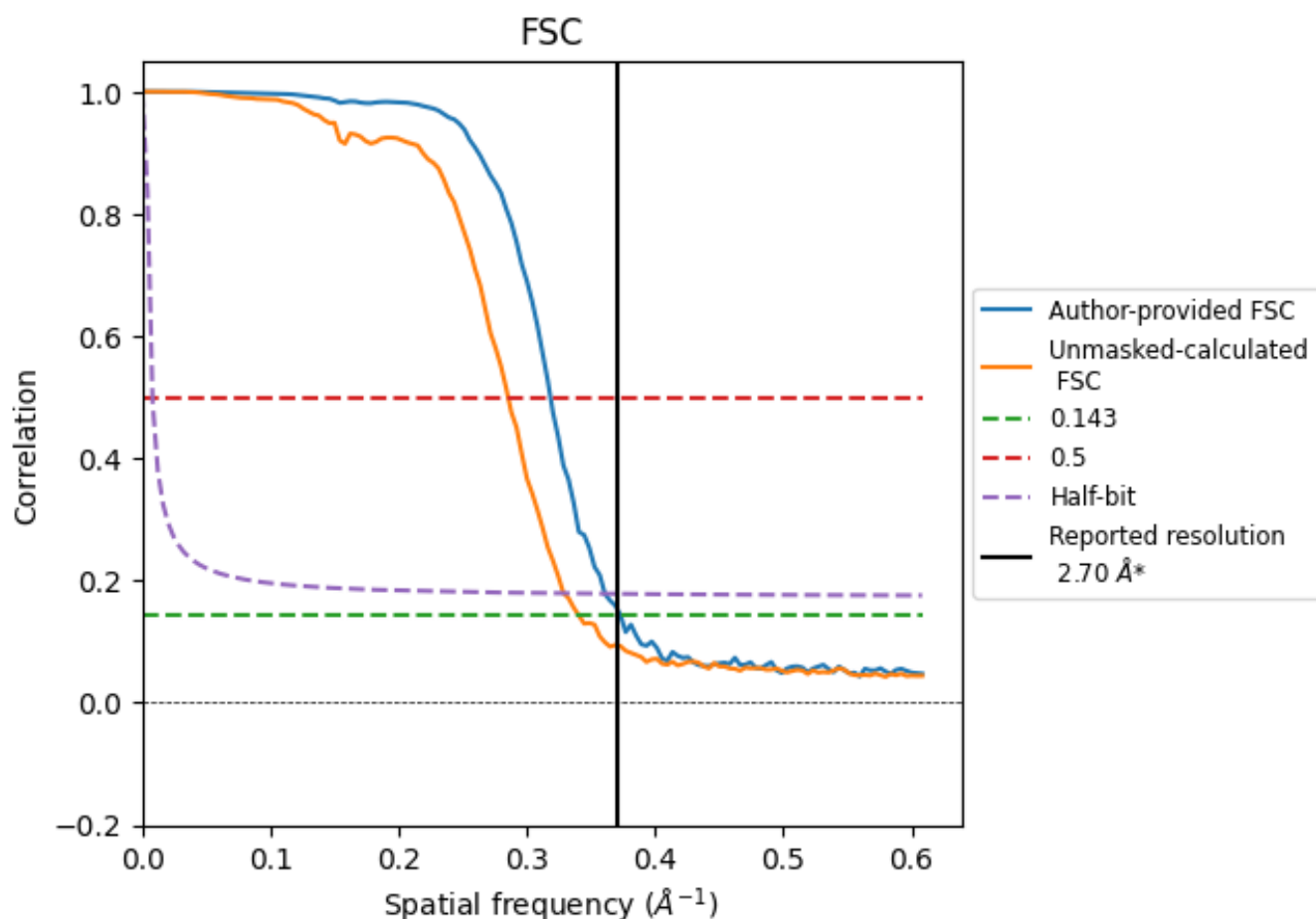


*Reported resolution corresponds to spatial frequency of 0.370 Å⁻¹

8 Fourier-Shell correlation [i](#)

Fourier-Shell Correlation (FSC) is the most commonly used method to estimate the resolution of single-particle and subtomogram-averaged maps. The shape of the curve depends on the imposed symmetry, mask and whether or not the two 3D reconstructions used were processed from a common reference. The reported resolution is shown as a black line. A curve is displayed for the half-bit criterion in addition to lines showing the 0.143 gold standard cut-off and 0.5 cut-off.

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.370 $\text{\AA}^{-1}$

8.2 Resolution estimates [i](#)

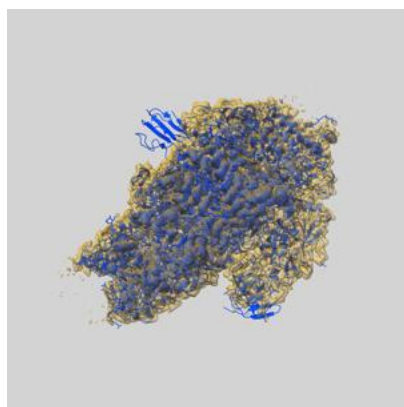
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	2.70	-	-
Author-provided FSC curve	2.68	3.14	2.77
Unmasked-calculated*	2.94	3.50	3.04

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

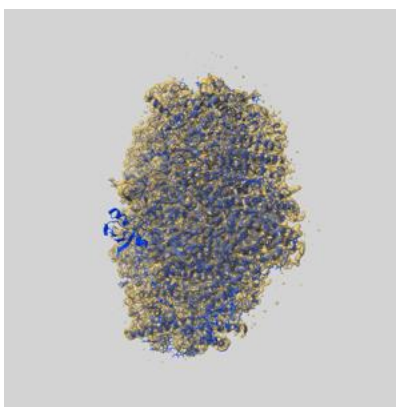
9 Map-model fit [i](#)

This section contains information regarding the fit between EMDB map EMD-10798 and PDB model 6YEZ. Per-residue inclusion information can be found in section 3 on page 31.

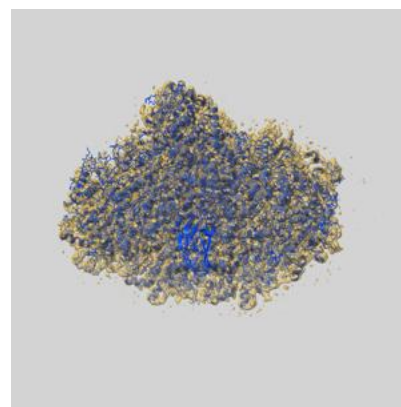
9.1 Map-model overlay [i](#)



X



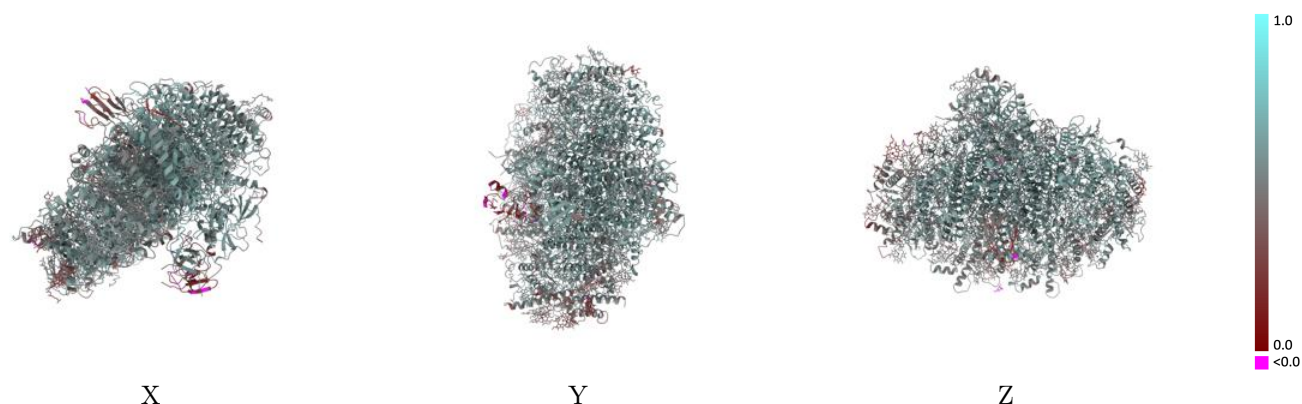
Y



Z

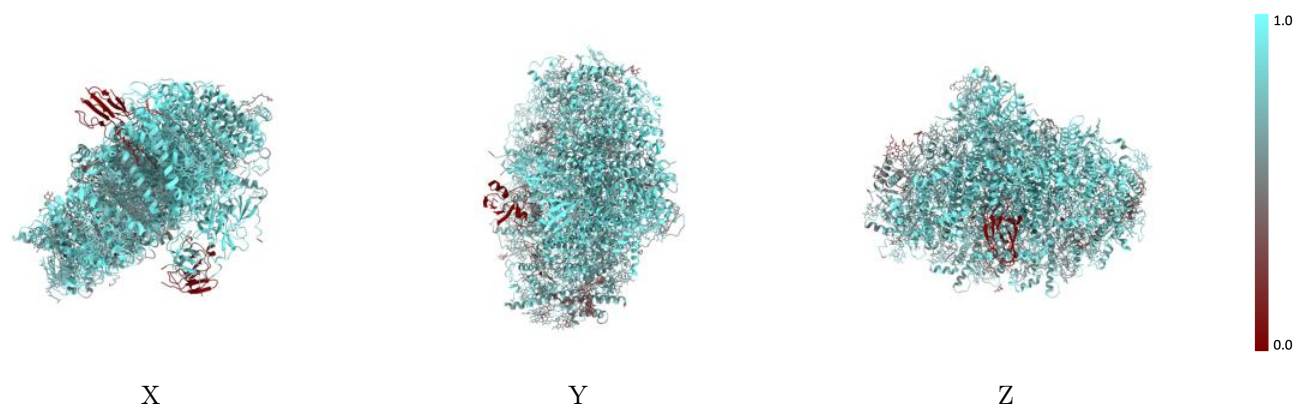
The images above show the 3D surface view of the map at the recommended contour level 0.022 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



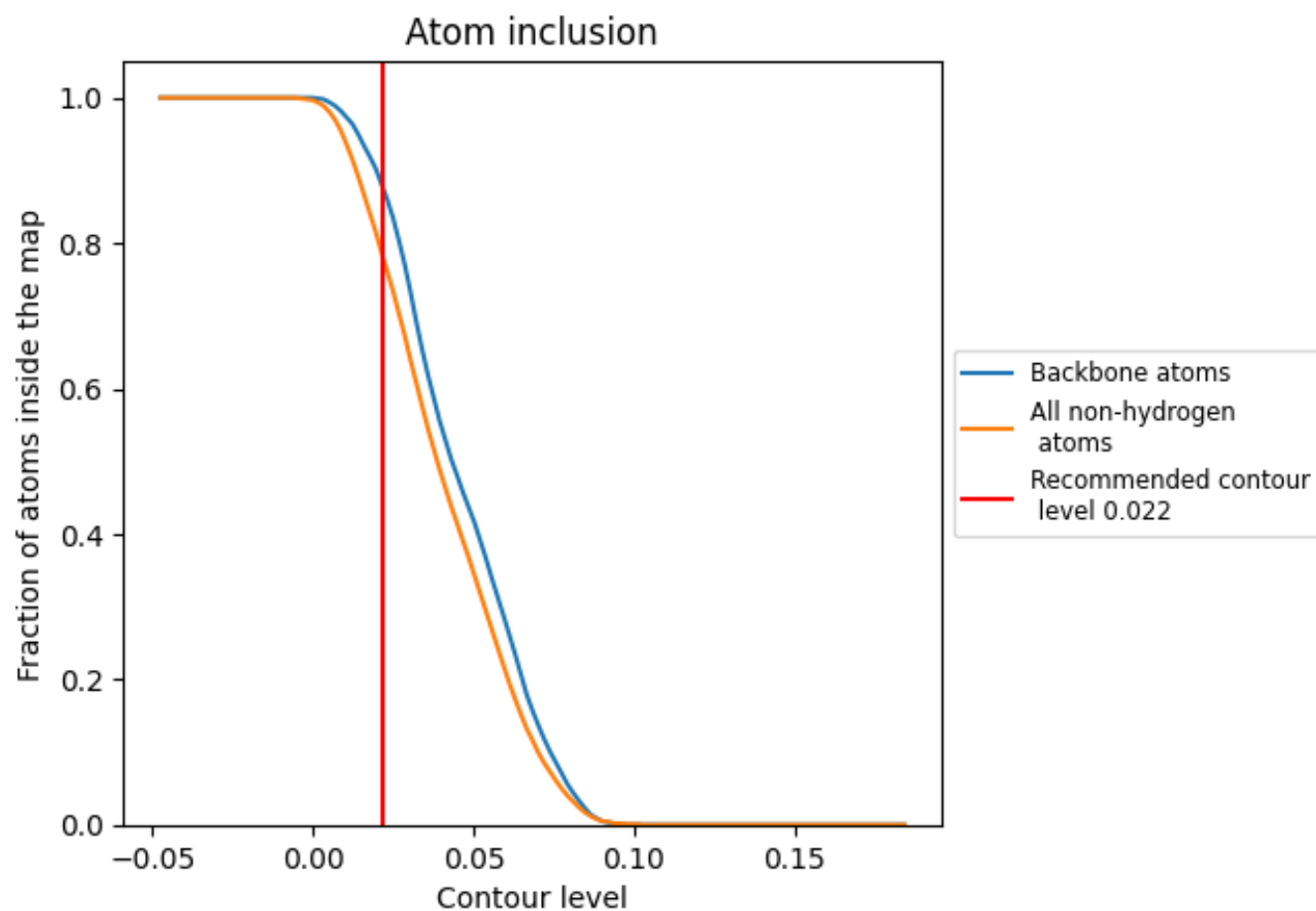
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.022).

9.4 Atom inclusion [i](#)



At the recommended contour level, 87% of all backbone atoms, 78% 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 (0.022) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.7780	 0.5360
1	 0.7920	 0.5340
2	 0.7140	 0.4750
3	 0.7210	 0.4840
4	 0.7870	 0.5180
A	 0.8470	 0.5710
B	 0.8780	 0.5920
C	 0.9270	 0.5930
D	 0.8900	 0.5660
E	 0.8310	 0.5480
F	 0.7670	 0.5400
G	 0.7330	 0.5090
H	 0.7480	 0.4990
I	 0.7890	 0.5460
J	 0.7540	 0.5320
K	 0.4650	 0.4130
L	 0.8330	 0.5480
N	 0.0570	 0.2140
P	 0.0040	 0.3360

