



wwPDB EM Validation Summary Report ⓘ

May 27, 2024 – 02:24 PM JST

PDB ID : 7EWK
EMDB ID : EMD-31350
Title : Barley photosystem I-LHCI-Lhca6 supercomplex
Authors : Wang, W.D.; Shen, L.; Tang, K.; Han, G.Y.; Zhang, X.; Shen, J.R.
Deposited on : 2021-05-25
Resolution : 3.88 Å(reported)

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

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

A user guide is available at

<https://www.wwpdb.org/validation/2017/EMValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

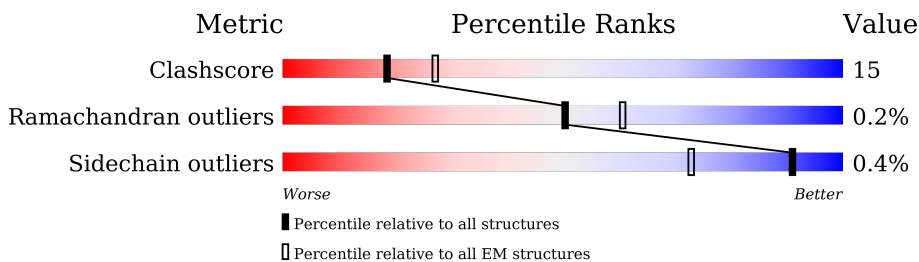
EMDB validation analysis : 0.0.1.dev92
Mogul : 1.8.5 (274361), CSD as541be (2020)
MolProbity : 4.02b-467
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
MapQ : 1.9.13
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

1 Overall quality at a glance

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

The reported resolution of this entry is 3.88 Å.

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



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

The table below summarises the geometric issues observed across the polymeric chains and their fit to the map. The red, orange, yellow and green segments of the bar indicate the fraction of residues that contain outliers for ≥ 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	742	
2	B	733	
3	C	81	
4	D	142	
5	E	68	
6	F	158	
7	H	61	
8	I	30	

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Mol	Chain	Length	Quality of chain
9	J	42	
10	K	84	
11	L	146	
12	1	193	
13	3	222	
14	4	197	
15	6	211	

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
16	CL0	A	801	X	-	-	-
17	CLA	1	504	X	-	-	-
17	CLA	1	506	X	-	-	-
17	CLA	1	507	X	-	-	-
17	CLA	1	508	X	-	-	-
17	CLA	1	509	X	-	-	-
17	CLA	1	510	X	-	-	-
17	CLA	1	511	X	-	-	-
17	CLA	1	513	X	-	-	-
17	CLA	1	515	X	-	-	-
17	CLA	3	304	X	-	-	-
17	CLA	3	305	X	-	-	-
17	CLA	3	306	X	-	-	-
17	CLA	3	308	X	-	-	-
17	CLA	3	309	X	-	-	-
17	CLA	3	310	X	-	-	-
17	CLA	3	311	X	-	-	-
17	CLA	3	312	X	-	-	-
17	CLA	3	314	X	-	-	-
17	CLA	3	316	X	-	-	-
17	CLA	4	304	X	-	-	-
17	CLA	4	305	X	-	-	-
17	CLA	4	306	X	-	-	-
17	CLA	4	307	X	-	-	-
17	CLA	4	308	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
17	CLA	4	309	X	-	-	-
17	CLA	4	310	X	-	-	-
17	CLA	4	311	X	-	-	-
17	CLA	4	312	X	-	-	-
17	CLA	4	315	X	-	-	-
17	CLA	4	317	X	-	-	-
17	CLA	6	504	X	-	-	-
17	CLA	6	506	X	-	-	-
17	CLA	6	507	X	-	-	-
17	CLA	6	508	X	-	-	-
17	CLA	6	509	X	-	-	-
17	CLA	6	510	X	-	-	-
17	CLA	6	511	X	-	-	-
17	CLA	6	514	X	-	-	-
17	CLA	A	802	X	-	-	-
17	CLA	A	803	X	-	-	-
17	CLA	A	804	X	-	-	-
17	CLA	A	805	X	-	-	-
17	CLA	A	806	X	-	-	-
17	CLA	A	807	X	-	-	-
17	CLA	A	808	X	-	-	-
17	CLA	A	809	X	-	-	-
17	CLA	A	810	X	-	-	-
17	CLA	A	811	X	-	-	-
17	CLA	A	812	X	-	-	-
17	CLA	A	813	X	-	-	-
17	CLA	A	814	X	-	-	-
17	CLA	A	815	X	-	-	-
17	CLA	A	816	X	-	-	-
17	CLA	A	817	X	-	-	-
17	CLA	A	818	X	-	-	-
17	CLA	A	819	X	-	-	-
17	CLA	A	820	X	-	-	-
17	CLA	A	821	X	-	-	-
17	CLA	A	822	X	-	-	-
17	CLA	A	823	X	-	-	-
17	CLA	A	824	X	-	-	-
17	CLA	A	825	X	-	-	-
17	CLA	A	826	X	-	-	-
17	CLA	A	827	X	-	-	-
17	CLA	A	828	X	-	-	-
17	CLA	A	829	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
17	CLA	A	830	X	-	-	-
17	CLA	A	831	X	-	-	-
17	CLA	A	832	X	-	-	-
17	CLA	A	833	X	-	-	-
17	CLA	A	834	X	-	-	-
17	CLA	A	835	X	-	-	-
17	CLA	A	836	X	-	-	-
17	CLA	A	837	X	-	-	-
17	CLA	A	838	X	-	-	-
17	CLA	A	839	X	-	-	-
17	CLA	A	840	X	-	-	-
17	CLA	A	841	X	-	-	-
17	CLA	A	843	X	-	-	-
17	CLA	A	852	X	-	-	-
17	CLA	A	853	X	-	-	-
17	CLA	B	801	X	-	-	-
17	CLA	B	803	X	-	-	-
17	CLA	B	804	X	-	-	-
17	CLA	B	805	X	-	-	-
17	CLA	B	806	X	-	-	-
17	CLA	B	807	X	-	-	-
17	CLA	B	808	X	-	-	-
17	CLA	B	809	X	-	-	-
17	CLA	B	810	X	-	-	-
17	CLA	B	811	X	-	-	-
17	CLA	B	812	X	-	-	-
17	CLA	B	813	X	-	-	-
17	CLA	B	814	X	-	-	-
17	CLA	B	815	X	-	-	-
17	CLA	B	816	X	-	-	-
17	CLA	B	817	X	-	-	-
17	CLA	B	818	X	-	-	-
17	CLA	B	819	X	-	-	-
17	CLA	B	820	X	-	-	-
17	CLA	B	821	X	-	-	-
17	CLA	B	822	X	-	-	-
17	CLA	B	823	X	-	-	-
17	CLA	B	824	X	-	-	-
17	CLA	B	825	X	-	-	-
17	CLA	B	826	X	-	-	-
17	CLA	B	827	X	-	-	-
17	CLA	B	828	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
17	CLA	B	829	X	-	-	-
17	CLA	B	830	X	-	-	-
17	CLA	B	831	X	-	-	-
17	CLA	B	832	X	-	-	-
17	CLA	B	833	X	-	-	-
17	CLA	B	834	X	-	-	-
17	CLA	B	835	X	-	-	-
17	CLA	B	836	X	-	-	-
17	CLA	B	837	X	-	-	-
17	CLA	B	838	X	-	-	-
17	CLA	B	839	X	-	-	-
17	CLA	B	840	X	-	-	-
17	CLA	B	841	X	-	-	-
17	CLA	B	842	X	-	-	-
17	CLA	B	843	X	-	-	-
17	CLA	F	802	X	-	-	-
17	CLA	J	101	X	-	-	-
17	CLA	J	102	X	-	-	-
17	CLA	K	201	X	-	-	-
17	CLA	K	202	X	-	-	-
17	CLA	K	203	X	-	-	-
17	CLA	K	205	X	-	-	-
17	CLA	L	302	X	-	-	-
17	CLA	L	303	X	-	-	-
17	CLA	L	304	X	-	-	-
24	LUT	1	502	X	-	-	-
25	CHL	1	512	X	-	-	-
25	CHL	1	514	X	-	-	-
25	CHL	1	517	X	-	-	-
25	CHL	3	313	X	-	-	-
25	CHL	4	313	X	-	-	-
25	CHL	4	314	X	-	-	-
25	CHL	4	316	X	-	-	-
25	CHL	6	512	X	-	-	-
25	CHL	6	513	X	-	-	-
25	CHL	6	515	X	-	-	-
25	CHL	6	517	X	-	-	-

2 Entry composition i

There are 26 unique types of molecules in this entry. The entry contains 32765 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
			Total	C	N	O	S		
1	A	742	5798	3797	987	996	18	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
2	B	733	5840	3832	991	1003	14	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
3	C	81	613	377	105	119	12	0	0

- Molecule 4 is a protein called Photosystem I reaction center subunit D.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
4	D	142	1112	714	195	200	3	0	0

- Molecule 5 is a protein called Photosystem I reaction center subunit E.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
5	E	68	540	341	99	100	0	0

- Molecule 6 is a protein called Photosystem I reaction center subunit F.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
6	F	158	1188	768	203	216	1	0	0

- Molecule 7 is a protein called Photosystem I reaction center subunit H.

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
7	H	61	447	298	73	76	0	0

- Molecule 8 is a protein called Photosystem I reaction center subunit VIII.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
8	I	30	235	163	35	36	1	0	0

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

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
9	J	42	332	225	51	55	1	0	0

- Molecule 10 is a protein called Photosystem I reaction center subunit K.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
10	K	84	576	361	101	110	4	0	0

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

Mol	Chain	Residues	Atoms				AltConf	Trace
			Total	C	N	O		
11	L	146	1090	720	174	196	0	0

- Molecule 12 is a protein called Chlorophyll a-b binding protein Lhca1.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
12	1	193	1484	965	248	269	2	0	0

- Molecule 13 is a protein called Chlorophyll a-b binding protein Lhca3.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
13	3	222	1672	1097	272	298	5	0	0

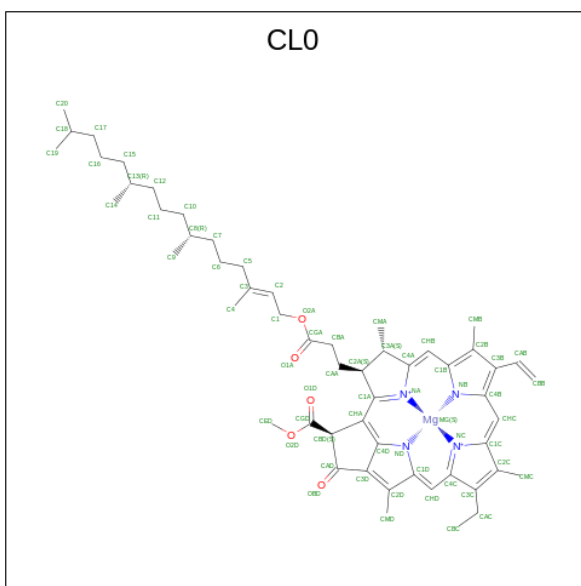
- Molecule 14 is a protein called Chlorophyll a-b binding protein, chloroplastic.

Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
14	4	197	1529	991	253	282	3	0	0

- Molecule 15 is a protein called Chlorophyll a-b binding protein Lhca6.

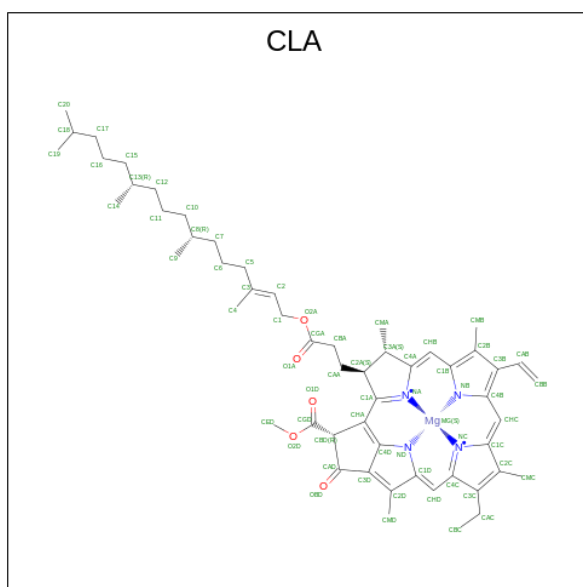
Mol	Chain	Residues	Atoms					AltConf	Trace
			Total	C	N	O	S		
15	6	211	1655	1088	274	285	8	0	0

- Molecule 16 is CHLOROPHYLL A ISOMER (three-letter code: CL0) (formula: $C_{55}H_{72}MgN_4O_5$).



Mol	Chain	Residues	Atoms					AltConf
			Total	C	Mg	N	O	
16	A	1	61	52	1	4	4	0

- Molecule 17 is CHLOROPHYLL A (three-letter code: CLA) (formula: $C_{55}H_{72}MgN_4O_5$).



Mol	Chain	Residues	Atoms				AltConf	
17	A	1	Total	C	Mg	N	O	0
			40	32	1	4	3	
17	A	1	Total	C	Mg	N	O	0
			52	42	1	4	5	
17	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
17	A	1	Total	C	Mg	N	O	0
			46	36	1	4	5	
17	A	1	Total	C	Mg	N	O	0
			41	33	1	4	3	
17	A	1	Total	C	Mg	N	O	0
			42	34	1	4	3	
17	A	1	Total	C	Mg	N	O	0
			41	33	1	4	3	
17	A	1	Total	C	Mg	N	O	0
			50	40	1	4	5	
17	A	1	Total	C	Mg	N	O	0
			55	45	1	4	5	
17	A	1	Total	C	Mg	N	O	0
			41	33	1	4	3	
17	A	1	Total	C	Mg	N	O	0
			39	32	1	4	2	
17	A	1	Total	C	Mg	N	O	0
			42	34	1	4	3	
17	A	1	Total	C	Mg	N	O	0
			42	34	1	4	3	
17	A	1	Total	C	Mg	N	O	0
			45	35	1	4	5	

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Mol	Chain	Residues	Atoms					AltConf
			Total	C	Mg	N	O	
17	A	1	60	50	1	4	5	0
17	A	1	39	32	1	4	2	0
17	A	1	39	31	1	4	3	0
17	A	1	45	35	1	4	5	0
17	A	1	41	33	1	4	3	0
17	A	1	42	34	1	4	3	0
17	A	1	41	33	1	4	3	0
17	A	1	42	34	1	4	3	0
17	A	1	56	46	1	4	5	0
17	A	1	41	33	1	4	3	0
17	A	1	64	54	1	4	5	0
17	A	1	41	33	1	4	3	0
17	A	1	45	36	1	4	4	0
17	A	1	58	48	1	4	5	0
17	A	1	50	40	1	4	5	0
17	A	1	56	46	1	4	5	0
17	A	1	43	35	1	4	3	0
17	A	1	41	33	1	4	3	0
17	A	1	45	35	1	4	5	0
17	A	1	42	34	1	4	3	0
17	A	1	51	41	1	4	5	0

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Mol	Chain	Residues	Atoms					AltConf
			Total	C	Mg	N	O	
17	A	1	55	45	1	4	5	0
17	A	1	52	42	1	4	5	0
17	A	1	65	55	1	4	5	0
17	A	1	41	33	1	4	3	0
17	A	1	41	33	1	4	3	0
17	A	1	49	39	1	4	5	0
17	A	1	57	47	1	4	5	0
17	A	1	45	35	1	4	5	0
17	B	1	64	54	1	4	5	0
17	B	1	41	33	1	4	3	0
17	B	1	42	33	1	4	4	0
17	B	1	38	31	1	4	2	0
17	B	1	41	33	1	4	3	0
17	B	1	41	33	1	4	3	0
17	B	1	54	45	1	4	4	0
17	B	1	52	42	1	4	5	0
17	B	1	65	55	1	4	5	0
17	B	1	51	41	1	4	5	0
17	B	1	40	32	1	4	3	0
17	B	1	54	44	1	4	5	0
17	B	1	43	35	1	4	3	0

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Mol	Chain	Residues	Atoms					AltConf
			Total	C	Mg	N	O	
17	B	1	43	35	1	4	3	0
17	B	1	38	32	1	4	1	0
17	B	1	43	35	1	4	3	0
17	B	1	55	45	1	4	5	0
17	B	1	41	33	1	4	3	0
17	B	1	50	40	1	4	5	0
17	B	1	55	45	1	4	5	0
17	B	1	41	33	1	4	3	0
17	B	1	41	33	1	4	3	0
17	B	1	42	34	1	4	3	0
17	B	1	42	34	1	4	3	0
17	B	1	65	55	1	4	5	0
17	B	1	40	32	1	4	3	0
17	B	1	41	33	1	4	3	0
17	B	1	50	40	1	4	5	0
17	B	1	39	31	1	4	3	0
17	B	1	56	46	1	4	5	0
17	B	1	40	32	1	4	3	0
17	B	1	43	35	1	4	3	0
17	B	1	43	35	1	4	3	0
17	B	1	45	35	1	4	5	0

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Mol	Chain	Residues	Atoms					AltConf
			Total	C	Mg	N	O	
17	B	1	58	49	1	4	4	0
17	B	1	42	34	1	4	3	0
17	B	1	40	32	1	4	3	0
17	B	1	41	33	1	4	3	0
17	B	1	41	33	1	4	3	0
17	B	1	41	33	1	4	3	0
17	B	1	41	33	1	4	3	0
17	B	1	38	32	1	4	1	0
17	F	1	41	33	1	4	3	0
17	J	1	41	33	1	4	3	0
17	J	1	42	34	1	4	3	0
17	K	1	37	31	1	4	1	0
17	K	1	45	35	1	4	5	0
17	K	1	41	33	1	4	3	0
17	K	1	39	31	1	4	3	0
17	L	1	40	32	1	4	3	0
17	L	1	60	50	1	4	5	0
17	L	1	45	35	1	4	5	0
17	1	1	41	33	1	4	3	0
17	1	1	46	36	1	4	5	0
17	1	1	55	45	1	4	5	0

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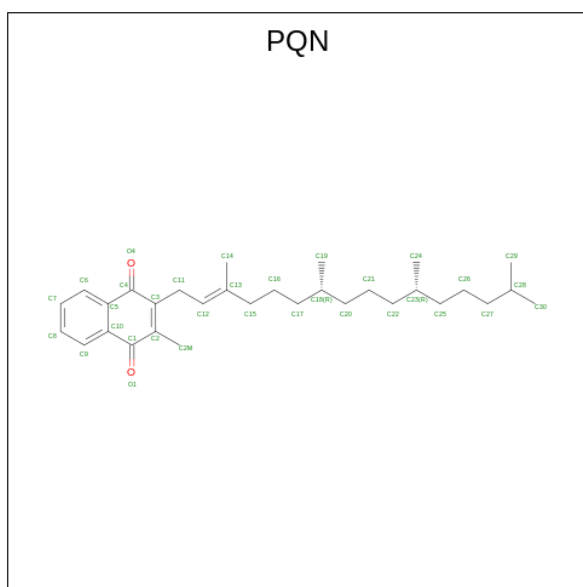
Mol	Chain	Residues	Atoms					AltConf
			Total	C	Mg	N	O	
17	1	1	40	32	1	4	3	0
17	1	1	65	55	1	4	5	0
17	1	1	50	40	1	4	5	0
17	1	1	46	36	1	4	5	0
17	1	1	39	31	1	4	3	0
17	1	1	52	42	1	4	5	0
17	1	1	40	32	1	4	3	0
17	3	1	38	30	1	4	3	0
17	3	1	52	42	1	4	5	0
17	3	1	41	33	1	4	3	0
17	3	1	40	32	1	4	3	0
17	3	1	42	34	1	4	3	0
17	3	1	50	40	1	4	5	0
17	3	1	41	33	1	4	3	0
17	3	1	48	38	1	4	5	0
17	3	1	60	50	1	4	5	0
17	3	1	40	32	1	4	3	0
17	3	1	45	35	1	4	5	0
17	3	1	46	36	1	4	5	0
17	4	1	39	31	1	4	3	0
17	4	1	50	40	1	4	5	0

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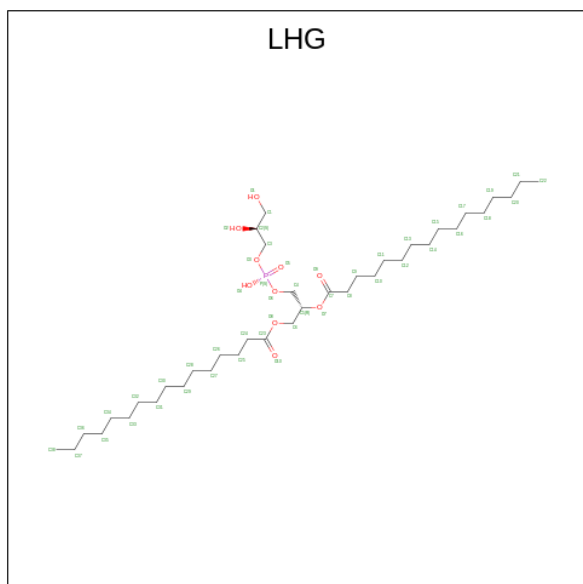
Mol	Chain	Residues	Atoms					AltConf
			Total	C	Mg	N	O	
17	4	1	65	55	1	4	5	0
17	4	1	60	50	1	4	5	0
17	4	1	41	33	1	4	3	0
17	4	1	50	40	1	4	5	0
17	4	1	42	34	1	4	3	0
17	4	1	46	36	1	4	5	0
17	4	1	41	33	1	4	3	0
17	4	1	41	33	1	4	3	0
17	4	1	50	40	1	4	5	0
17	6	1	39	31	1	4	3	0
17	6	1	52	42	1	4	5	0
17	6	1	65	55	1	4	5	0
17	6	1	51	41	1	4	5	0
17	6	1	40	32	1	4	3	0
17	6	1	50	40	1	4	5	0
17	6	1	60	50	1	4	5	0
17	6	1	50	40	1	4	5	0
17	6	1	40	32	1	4	3	0

- Molecule 18 is PHYLLOQUINONE (three-letter code: PQN) (formula: C₃₁H₄₆O₂).



Mol	Chain	Residues	Atoms			AltConf
18	A	1	Total	C	O	0
			14	12	2	
18	B	1	Total	C	O	0
			16	14	2	

- Molecule 19 is 1,2-DIPALMITOYL-PHOSPHATIDYL-GLYCEROLE (three-letter code: LHG) (formula: C₃₈H₇₅O₁₀P).



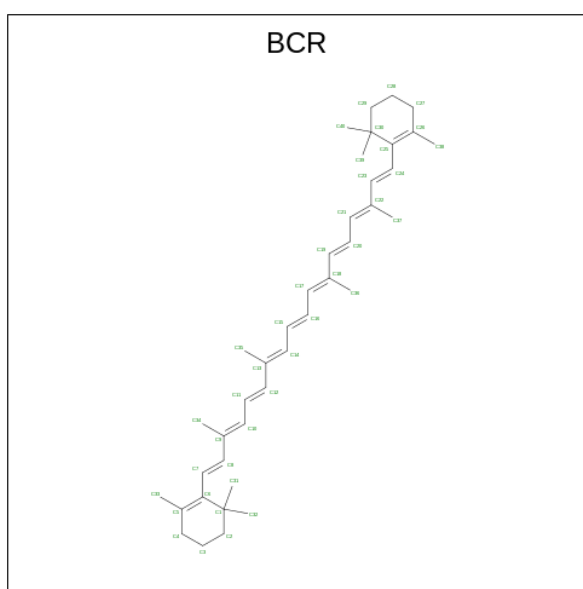
Mol	Chain	Residues	Atoms				AltConf
19	A	1	Total	C	O	P	0
			40	29	10	1	

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Mol	Chain	Residues	Atoms				AltConf
			Total	C	O	P	
19	A	1	Total 25	C 15	O 9	P 1	0
19	B	1	Total 38	C 27	O 10	P 1	0
19	1	1	Total 49	C 38	O 10	P 1	0
19	6	1	Total 35	C 24	O 10	P 1	0

- Molecule 20 is BETA-CAROTENE (three-letter code: BCR) (formula: C₄₀H₅₆).



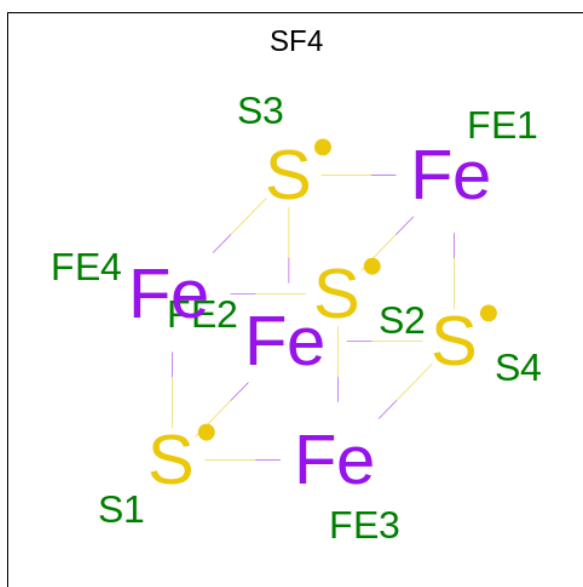
Mol	Chain	Residues	Atoms		AltConf
20	A	1	Total 40	C 40	0
20	A	1	Total 40	C 40	0
20	A	1	Total 40	C 40	0
20	A	1	Total 40	C 40	0
20	A	1	Total 40	C 40	0
20	A	1	Total 40	C 40	0
20	A	1	Total 40	C 40	0

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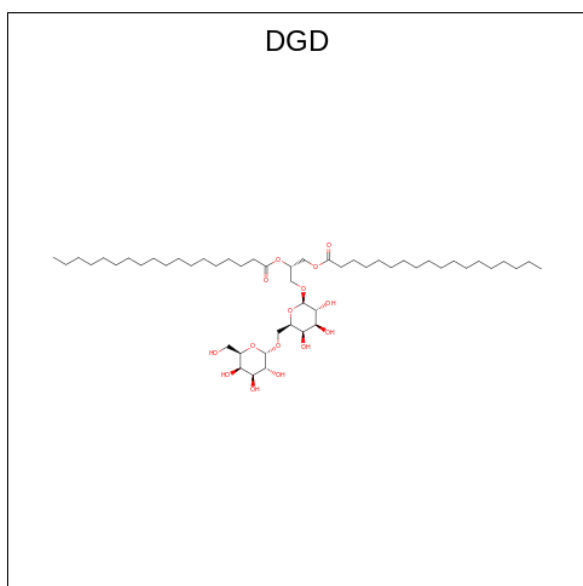
Mol	Chain	Residues	Atoms	AltConf
20	B	1	Total C 40 40	0
20	B	1	Total C 40 40	0
20	B	1	Total C 40 40	0
20	B	1	Total C 40 40	0
20	B	1	Total C 40 40	0
20	B	1	Total C 39 39	0
20	F	1	Total C 40 40	0
20	F	1	Total C 40 40	0
20	I	1	Total C 40 40	0
20	I	1	Total C 40 40	0
20	J	1	Total C 40 40	0
20	K	1	Total C 40 40	0
20	L	1	Total C 40 40	0
20	L	1	Total C 40 40	0
20	1	1	Total C 11 11	0
20	3	1	Total C 40 40	0
20	4	1	Total C 40 40	0
20	6	1	Total C 40 40	0

- Molecule 21 is IRON/SULFUR CLUSTER (three-letter code: SF4) (formula: Fe₄S₄).



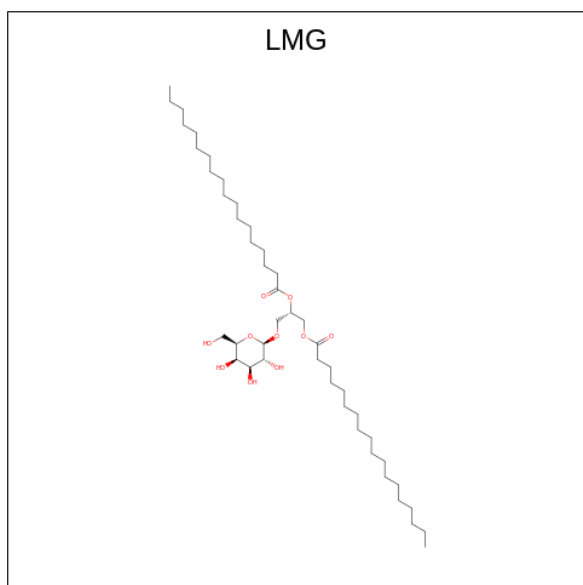
Mol	Chain	Residues	Atoms			AltConf
			Total	Fe	S	
21	B	1	8	4	4	0
21	C	1	8	4	4	0
21	C	1	8	4	4	0

- Molecule 22 is DIGALACTOSYL DIACYL GLYCEROL (DGDG) (three-letter code: DGD) (formula: $C_{51}H_{96}O_{15}$).



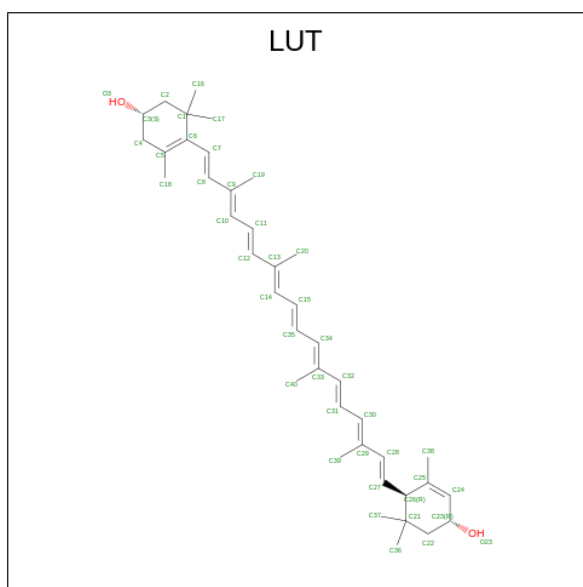
Mol	Chain	Residues	Atoms			AltConf
22	B	1	Total	C	O	0
			52	37	15	
22	J	1	Total	C	O	0
			66	51	15	

- Molecule 23 is 1,2-DISTEAROYL-MONOGALACTOSYL-DIGLYCERIDE (three-letter code: LMG) (formula: C₄₅H₈₆O₁₀).



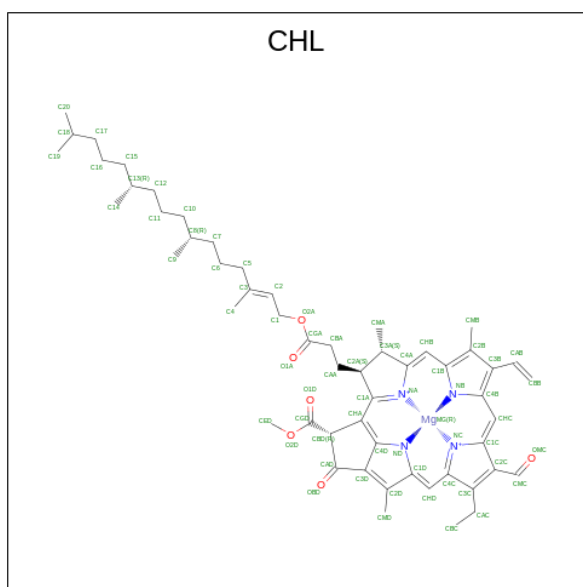
Mol	Chain	Residues	Atoms			AltConf
23	J	1	Total	C	O	0
			30	20	10	
23	4	1	Total	C	O	0
			18	10	8	

- Molecule 24 is (3R,3'R,6S)-4,5-DIDEHYDRO-5,6-DIHYDRO-BETA,BETA-CAROTENE-3,3'-DIOL (three-letter code: LUT) (formula: C₄₀H₅₆O₂).



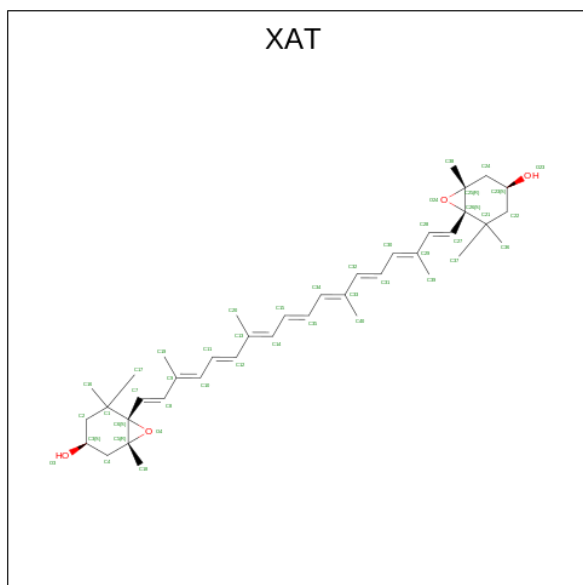
Mol	Chain	Residues	Atoms			AltConf
24	1	1	Total	C	O	0
			42	40	2	
24	1	1	Total	C	O	0
			42	40	2	
24	3	1	Total	C	O	0
			42	40	2	
24	3	1	Total	C	O	0
			42	40	2	
24	4	1	Total	C	O	0
			42	40	2	
24	6	1	Total	C	O	0
			42	40	2	

- Molecule 25 is CHLOROPHYLL B (three-letter code: CHL) (formula: $C_{55}H_{70}MgN_4O_6$).



Mol	Chain	Residues	Atoms				AltConf	
			Total	C	Mg	N		O
25	1	1	47	36	1	4	6	0
25	1	1	47	36	1	4	6	0
25	1	1	43	34	1	4	4	0
25	3	1	46	36	1	4	5	0
25	4	1	47	36	1	4	6	0
25	4	1	51	40	1	4	6	0
25	4	1	43	34	1	4	4	0
25	6	1	47	36	1	4	6	0
25	6	1	48	37	1	4	6	0
25	6	1	41	32	1	4	4	0
25	6	1	47	36	1	4	6	0

- Molecule 26 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 (three-letter code: XAT) (formula: C₄₀H₅₆O₄).

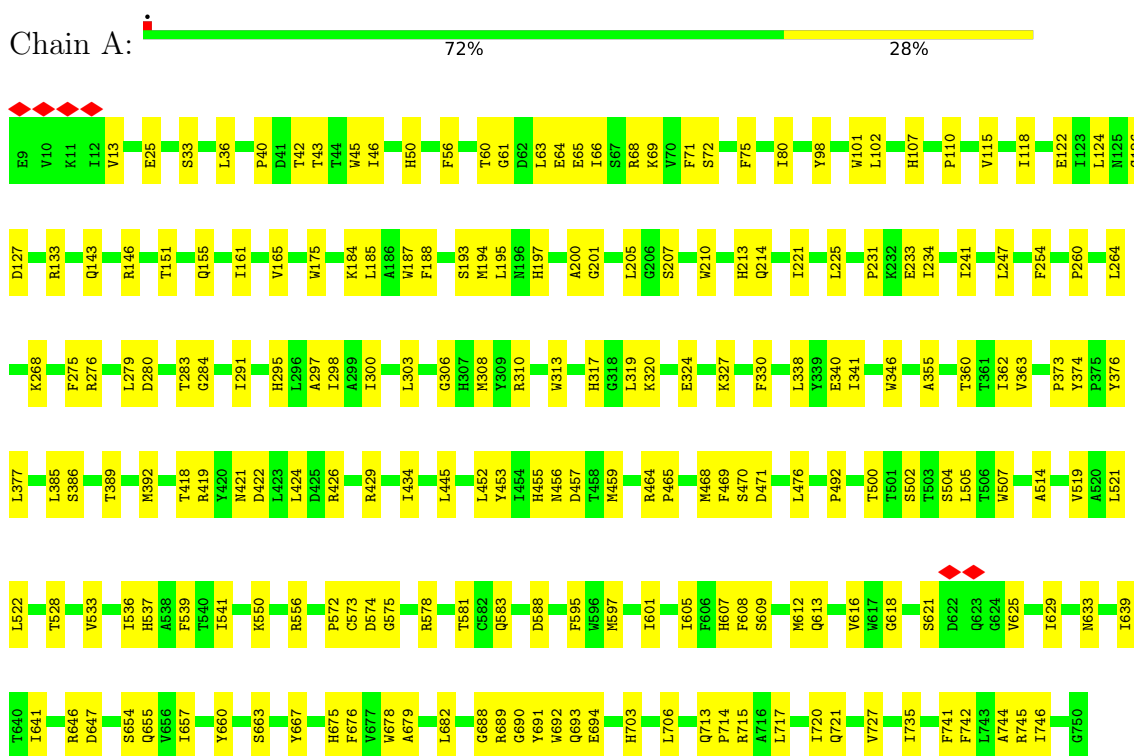


Mol	Chain	Residues	Atoms			AltConf
26	4	1	Total	C	O	0
			44	40	4	
26	6	1	Total	C	O	0
			44	40	4	

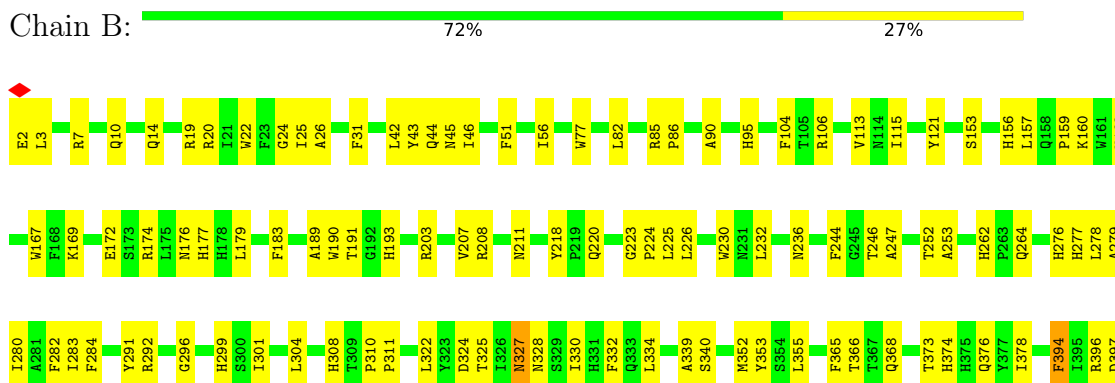
3 Residue-property plots

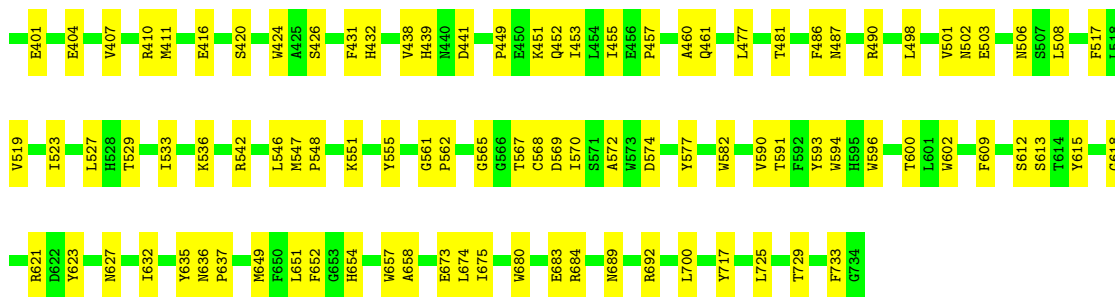
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and atom inclusion 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

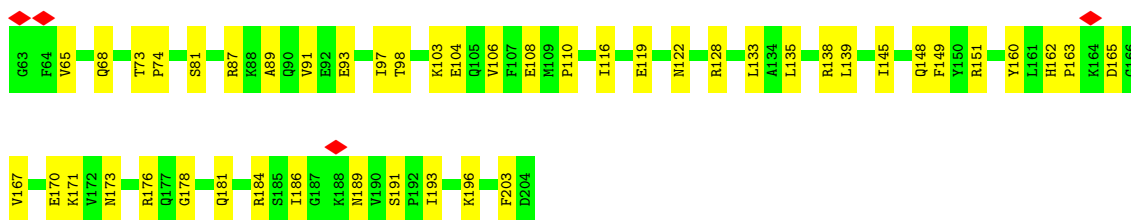




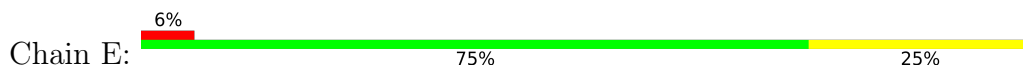
• Molecule 3: Photosystem I iron-sulfur center



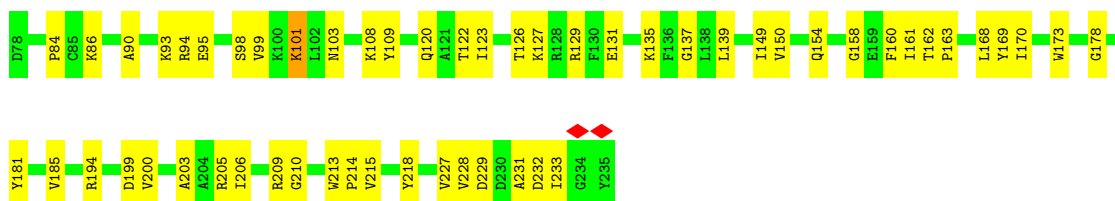
• Molecule 4: Photosystem I reaction center subunit D



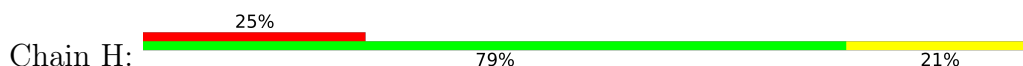
• Molecule 5: Photosystem I reaction center subunit E

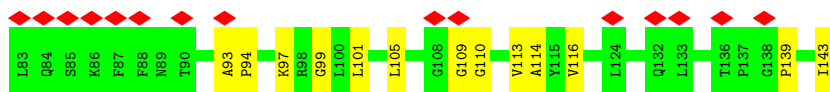


• Molecule 6: Photosystem I reaction center subunit F

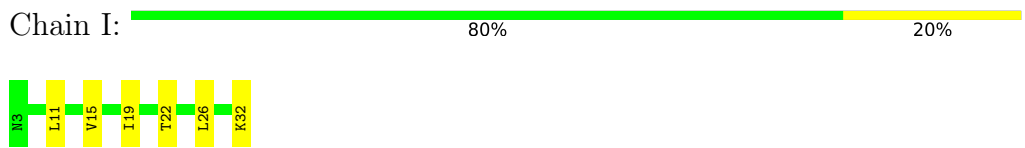


• Molecule 7: Photosystem I reaction center subunit H

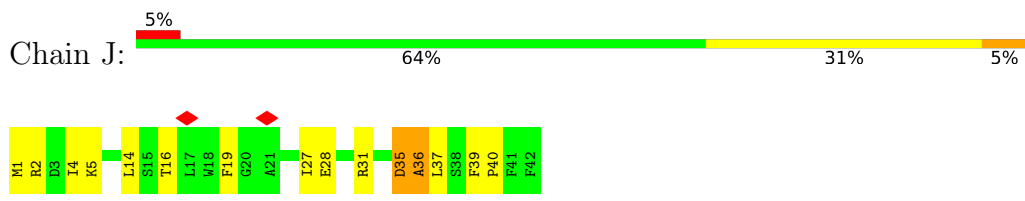




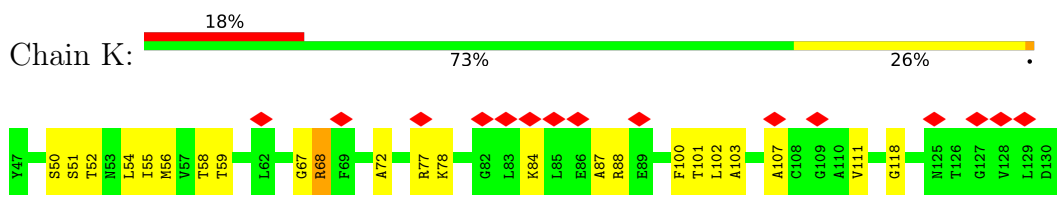
• Molecule 8: Photosystem I reaction center subunit VIII



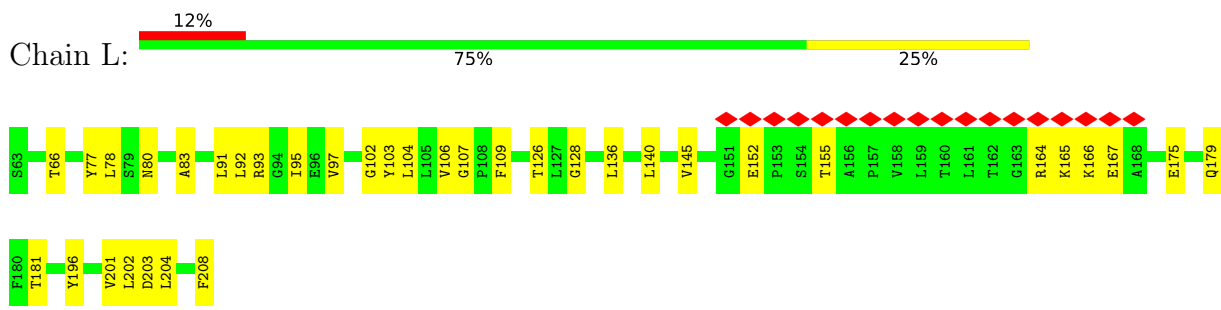
• Molecule 9: Photosystem I reaction center subunit IX



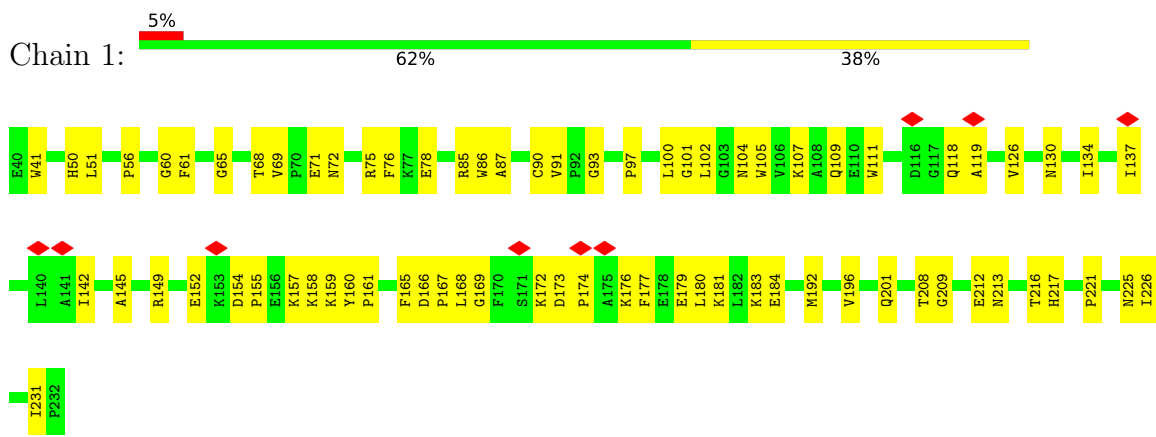
• Molecule 10: Photosystem I reaction center subunit K



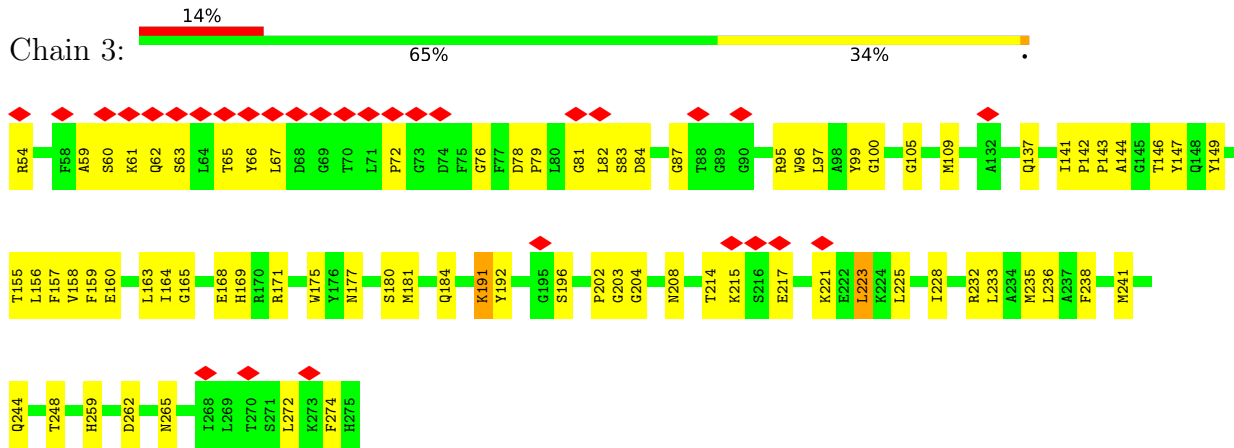
• Molecule 11: Photosystem I reaction center subunit L



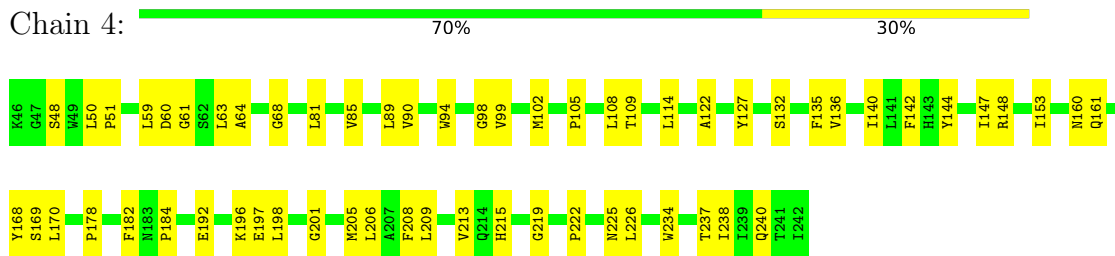
• Molecule 12: Chlorophyll a-b binding protein Lhca1



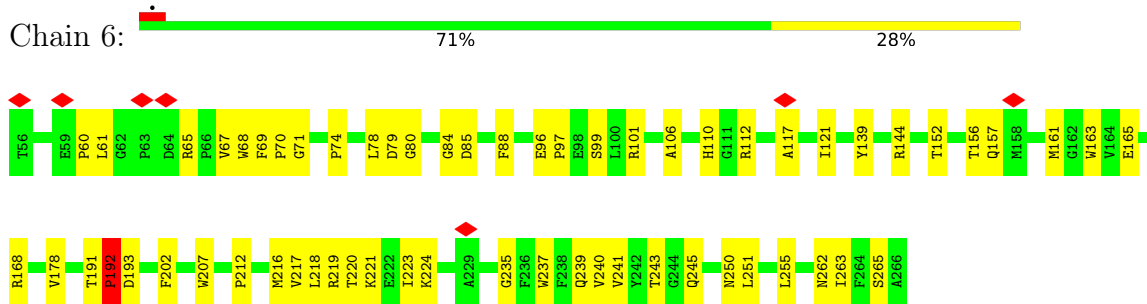
• Molecule 13: Chlorophyll a-b binding protein Lhca3



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



• Molecule 15: Chlorophyll a-b binding protein Lhca6



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, Not provided	
Number of particles used	103844	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	FEI TITAN KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	2000	Depositor
Maximum defocus (nm)	2500	Depositor
Magnification	Not provided	
Image detector	GATAN K2 SUMMIT (4k x 4k)	Depositor
Maximum map value	4.561	Depositor
Minimum map value	-1.733	Depositor
Average map value	-0.001	Depositor
Map value standard deviation	0.074	Depositor
Recommended contour level	0.43	Depositor
Map size (Å)	575.08, 575.08, 575.08	wwPDB
Map dimensions	440, 440, 440	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.307, 1.307, 1.307	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: DGD, LUT, CL0, LHG, SF4, PQN, BCR, CHL, CLA, XAT, LMG

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	0.31	0/5993	0.43	0/8178
2	B	0.36	3/6052 (0.0%)	0.43	1/8267 (0.0%)
3	C	0.30	0/624	0.46	0/844
4	D	0.27	0/1141	0.47	0/1540
5	E	0.30	0/553	0.47	0/751
6	F	0.26	0/1214	0.45	0/1643
7	H	0.26	0/460	0.45	0/623
8	I	0.27	0/241	0.41	0/327
9	J	0.26	0/342	0.45	0/465
10	K	0.26	0/581	0.51	0/786
11	L	0.27	0/1122	0.46	0/1537
12	1	0.28	0/1535	0.45	0/2097
13	3	0.30	0/1727	0.48	0/2353
14	4	0.27	0/1577	0.44	0/2153
15	6	0.27	0/1723	0.48	2/2358 (0.1%)
All	All	0.31	3/24885 (0.0%)	0.45	3/33922 (0.0%)

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
9	J	0	1
12	1	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	657	TRP	CB-CG	-10.12	1.32	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	657	TRP	C-O	-6.31	1.11	1.23
2	B	657	TRP	CG-CD1	-6.28	1.27	1.36

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	657	TRP	CB-CA-C	-5.93	98.55	110.40
15	6	192	PRO	N-CA-C	5.69	126.89	112.10
15	6	192	PRO	C-N-CA	-5.19	108.73	121.70

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
12	1	231	ILE	Peptide
9	J	35	ASP	Peptide

5.2 Too-close contacts [\(i\)](#)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5798	0	5613	173	0
2	B	5840	0	5580	173	0
3	C	613	0	607	27	0
4	D	1112	0	1113	40	0
5	E	540	0	539	12	0
6	F	1188	0	1178	39	0
7	H	447	0	472	9	0
8	I	235	0	261	6	0
9	J	332	0	335	17	0
10	K	576	0	584	17	0
11	L	1090	0	1075	30	0
12	1	1484	0	1410	62	0
13	3	1672	0	1575	69	0
14	4	1529	0	1448	47	0
15	6	1655	0	1530	54	0
16	A	61	0	62	6	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
17	1	474	0	373	27	0
17	3	543	0	404	28	0
17	4	525	0	424	23	0
17	6	447	0	371	25	0
17	A	2028	0	1626	85	0
17	B	1921	0	1521	85	0
17	F	41	0	29	0	0
17	J	83	0	59	1	0
17	K	162	0	112	3	0
17	L	145	0	118	7	0
18	A	14	0	7	2	0
18	B	16	0	10	7	0
19	1	49	0	74	7	0
19	6	35	0	40	1	0
19	A	65	0	77	2	0
19	B	38	0	46	1	0
20	1	11	0	16	1	0
20	3	40	0	56	3	0
20	4	40	0	56	5	0
20	6	40	0	56	4	0
20	A	280	0	392	23	0
20	B	239	0	333	13	0
20	F	80	0	112	2	0
20	I	80	0	112	5	0
20	J	40	0	56	3	0
20	K	40	0	56	1	0
20	L	80	0	112	3	0
21	B	8	0	0	0	0
21	C	16	0	0	1	0
22	B	52	0	62	2	0
22	J	66	0	96	2	0
23	4	18	0	12	0	0
23	J	30	0	30	1	0
24	1	84	0	111	12	0
24	3	84	0	112	11	0
24	4	42	0	56	9	0
24	6	42	0	56	6	0
25	1	137	0	91	6	0
25	3	46	0	31	1	0
25	4	141	0	97	3	0
25	6	183	0	121	8	0
26	4	44	0	56	3	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
26	6	44	0	56	9	0
All	All	32765	0	31017	927	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 15.

The worst 5 of 927 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:700:LEU:HB2	18:B:844:PQN:H6	1.62	0.81
3:C:54:CYS:SG	3:C:55:GLU:N	2.61	0.72
6:F:215:VAL:O	6:F:218:TYR:HB3	1.89	0.72
1:A:175:TRP:HB2	17:A:810:CLA:HMC3	1.72	0.71
12:1:90:CYS:CB	24:1:502:LUT:H392	2.21	0.70

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	740/742 (100%)	684 (92%)	56 (8%)	0	100	100
2	B	731/733 (100%)	680 (93%)	51 (7%)	0	100	100
3	C	79/81 (98%)	68 (86%)	11 (14%)	0	100	100
4	D	140/142 (99%)	125 (89%)	15 (11%)	0	100	100
5	E	66/68 (97%)	61 (92%)	5 (8%)	0	100	100
6	F	156/158 (99%)	148 (95%)	8 (5%)	0	100	100
7	H	59/61 (97%)	56 (95%)	3 (5%)	0	100	100
8	I	28/30 (93%)	28 (100%)	0	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
9	J	40/42 (95%)	38 (95%)	1 (2%)	1 (2%)	5	36
10	K	82/84 (98%)	75 (92%)	7 (8%)	0	100	100
11	L	144/146 (99%)	134 (93%)	10 (7%)	0	100	100
12	1	191/193 (99%)	171 (90%)	18 (9%)	2 (1%)	15	52
13	3	220/222 (99%)	193 (88%)	27 (12%)	0	100	100
14	4	195/197 (99%)	183 (94%)	12 (6%)	0	100	100
15	6	209/211 (99%)	190 (91%)	17 (8%)	2 (1%)	15	52
All	All	3080/3110 (99%)	2834 (92%)	241 (8%)	5 (0%)	50	79

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
15	6	191	THR
15	6	192	PRO
9	J	36	ALA
12	1	101	GLY
12	1	221	PRO

5.3.2 Protein sidechains [i](#)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	588/600 (98%)	588 (100%)	0	100	100
2	B	589/600 (98%)	587 (100%)	2 (0%)	92	95
3	C	71/71 (100%)	71 (100%)	0	100	100
4	D	118/119 (99%)	118 (100%)	0	100	100
5	E	58/59 (98%)	58 (100%)	0	100	100
6	F	115/127 (91%)	114 (99%)	1 (1%)	78	87
7	H	47/49 (96%)	47 (100%)	0	100	100
8	I	27/27 (100%)	27 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
9	J	34/36 (94%)	34 (100%)	0	100	100
10	K	56/61 (92%)	54 (96%)	2 (4%)	35	61
11	L	110/115 (96%)	110 (100%)	0	100	100
12	1	147/155 (95%)	145 (99%)	2 (1%)	67	81
13	3	159/176 (90%)	157 (99%)	2 (1%)	69	81
14	4	156/164 (95%)	156 (100%)	0	100	100
15	6	162/172 (94%)	162 (100%)	0	100	100
All	All	2437/2531 (96%)	2428 (100%)	9 (0%)	91	94

5 of 9 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
13	3	191	LYS
13	3	223	LEU
10	K	68	ARG
10	K	88	ARG
12	1	217	HIS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 11 such sidechains are listed below:

Mol	Chain	Res	Type
5	E	135	ASN
13	3	255	ASN
15	6	250	ASN
13	3	265	ASN
4	D	68	GLN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

5.6 Ligand geometry [i](#)

196 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
17	CLA	A	833	-	41,49,73	1.82	8 (19%)	47,84,113	1.62	9 (19%)
17	CLA	A	834	-	45,53,73	1.74	7 (15%)	52,89,113	1.59	7 (13%)
17	CLA	B	837	-	42,50,73	1.84	6 (14%)	48,85,113	1.58	6 (12%)
17	CLA	1	508	-	65,73,73	1.47	6 (9%)	76,113,113	1.33	6 (7%)
17	CLA	B	806	-	41,49,73	1.81	6 (14%)	47,84,113	1.62	7 (14%)
17	CLA	A	802	-	39,48,73	1.87	7 (17%)	45,82,113	1.76	8 (17%)
17	CLA	1	506	-	55,63,73	1.61	6 (10%)	64,101,113	1.42	9 (14%)
19	LHG	1	516	-	48,48,48	0.93	2 (4%)	51,54,54	1.07	4 (7%)
17	CLA	A	813	-	42,50,73	1.83	7 (16%)	48,85,113	1.57	7 (14%)
17	CLA	J	102	-	42,50,73	1.84	5 (11%)	48,85,113	1.60	6 (12%)
17	CLA	K	202	-	45,53,73	1.78	5 (11%)	52,89,113	1.57	9 (17%)
17	CLA	L	303	-	60,68,73	1.51	6 (10%)	70,107,113	1.43	7 (10%)
20	BCR	B	850	-	40,40,41	0.76	0	54,54,56	2.08	15 (27%)
25	CHL	1	514	-	47,55,74	2.42	16 (34%)	50,91,114	2.77	20 (40%)
17	CLA	A	809	-	50,58,73	1.66	6 (12%)	58,95,113	1.66	9 (15%)
17	CLA	B	828	-	41,49,73	1.81	8 (19%)	47,84,113	1.74	7 (14%)
17	CLA	B	841	-	41,49,73	1.80	7 (17%)	47,84,113	1.69	7 (14%)
23	LMG	J	105	-	30,30,55	1.20	2 (6%)	38,38,63	1.22	3 (7%)
17	CLA	A	825	-	41,49,73	1.78	6 (14%)	47,84,113	1.71	8 (17%)
17	CLA	6	508	-	40,48,73	1.84	6 (15%)	46,83,113	1.73	8 (17%)
20	BCR	L	305	-	41,41,41	0.71	0	56,56,56	2.18	18 (32%)
17	CLA	A	828	-	43,51,73	1.78	7 (16%)	48,85,113	1.61	6 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
17	CLA	3	306	-	41,49,73	1.86	6 (14%)	47,84,113	1.69	8 (17%)
17	CLA	A	840	-	41,49,73	1.80	6 (14%)	47,84,113	1.70	7 (14%)
17	CLA	A	824	-	56,64,73	1.55	7 (12%)	65,102,113	1.45	9 (13%)
17	CLA	A	853	-	45,53,73	1.74	7 (15%)	52,89,113	1.66	8 (15%)
17	CLA	A	827	-	41,49,73	1.80	8 (19%)	47,84,113	1.64	10 (21%)
17	CLA	B	804	-	43,50,73	2.36	9 (20%)	46,83,113	2.07	9 (19%)
17	CLA	B	810	-	65,73,73	1.46	7 (10%)	76,113,113	1.44	8 (10%)
17	CLA	1	505	-	46,54,73	1.65	6 (13%)	53,90,113	1.54	8 (15%)
17	CLA	3	308	-	42,50,73	1.81	5 (11%)	48,85,113	1.69	6 (12%)
19	LHG	A	845	-	24,24,48	0.96	1 (4%)	26,29,54	1.26	3 (11%)
17	CLA	A	832	-	43,51,73	1.80	6 (13%)	49,86,113	1.56	8 (16%)
17	CLA	B	836	-	59,66,73	1.68	7 (11%)	64,103,113	1.42	8 (12%)
17	CLA	4	308	-	41,49,73	1.83	6 (14%)	47,84,113	1.67	7 (14%)
24	LUT	1	502	-	42,43,43	2.39	1 (2%)	51,60,60	1.81	9 (17%)
25	CHL	4	313	-	47,55,74	2.39	16 (34%)	50,91,114	2.74	20 (40%)
17	CLA	B	843	19	37,46,73	1.89	5 (13%)	44,80,113	1.67	7 (15%)
17	CLA	B	830	-	38,47,73	1.93	8 (21%)	48,81,113	1.81	10 (20%)
17	CLA	6	514	-	39,48,73	1.83	6 (15%)	45,82,113	1.74	8 (17%)
20	BCR	F	803	-	41,41,41	0.76	0	56,56,56	2.16	18 (32%)
17	CLA	6	511	-	50,58,73	1.70	6 (12%)	58,95,113	1.53	8 (13%)
17	CLA	A	804	-	65,73,73	1.44	7 (10%)	76,113,113	1.42	6 (7%)
17	CLA	6	507	-	51,59,73	1.66	7 (13%)	59,96,113	1.53	8 (13%)
17	CLA	K	205	-	37,47,73	1.97	6 (16%)	42,81,113	1.63	7 (16%)
20	BCR	B	845	-	41,41,41	0.75	0	56,56,56	2.15	19 (33%)
20	BCR	I	102	-	41,41,41	0.70	0	56,56,56	2.15	16 (28%)
17	CLA	A	810	17	55,63,73	1.61	8 (14%)	64,101,113	1.45	8 (12%)
20	BCR	J	103	-	41,41,41	0.77	1 (2%)	56,56,56	2.11	17 (30%)
25	CHL	3	313	-	47,54,74	2.40	15 (31%)	45,89,114	2.89	19 (42%)
24	LUT	3	302	-	42,43,43	0.77	0	51,60,60	1.85	16 (31%)
17	CLA	3	311	-	48,56,73	1.72	6 (12%)	55,92,113	1.54	7 (12%)
17	CLA	B	819	-	41,49,73	1.83	6 (14%)	47,84,113	1.63	8 (17%)
17	CLA	L	304	-	45,53,73	1.76	6 (13%)	52,89,113	1.60	6 (11%)
17	CLA	4	310	-	42,50,73	1.84	6 (14%)	48,85,113	1.61	7 (14%)
21	SF4	B	802	-	0,12,12	-	-	-	-	-
17	CLA	1	509	-	50,58,73	1.50	6 (12%)	58,95,113	2.06	16 (27%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
17	CLA	B	840	-	41,49,73	1.81	7 (17%)	47,84,113	1.72	9 (19%)
20	BCR	6	503	-	41,41,41	0.71	0	56,56,56	3.61	30 (53%)
17	CLA	B	839	-	41,49,73	1.81	5 (12%)	47,84,113	1.72	8 (17%)
17	CLA	A	823	-	42,50,73	1.83	7 (16%)	52,85,113	1.74	9 (17%)
17	CLA	3	307	13	39,48,73	1.90	5 (12%)	45,82,113	1.71	8 (17%)
17	CLA	A	843	-	49,57,73	1.75	6 (12%)	61,94,113	1.67	8 (13%)
17	CLA	B	833	-	43,51,73	1.77	6 (13%)	49,86,113	1.55	6 (12%)
25	CHL	1	512	-	47,55,74	2.44	16 (34%)	50,91,114	2.79	22 (44%)
17	CLA	3	316	-	46,54,73	1.77	6 (13%)	53,90,113	1.49	7 (13%)
20	BCR	4	301	-	41,41,41	0.72	0	56,56,56	1.85	14 (25%)
17	CLA	B	809	-	52,60,73	1.66	8 (15%)	60,97,113	1.47	8 (13%)
16	CL0	A	801	-	61,69,73	1.64	10 (16%)	70,107,113	2.21	16 (22%)
21	SF4	C	102	-	0,12,12	-	-	-	-	-
17	CLA	A	808	1	41,49,73	1.85	7 (17%)	47,84,113	1.67	8 (17%)
20	BCR	I	101	-	41,41,41	0.80	1 (2%)	56,56,56	2.29	21 (37%)
17	CLA	F	802	-	41,49,73	1.86	6 (14%)	47,84,113	1.63	7 (14%)
19	LHG	6	516	-	34,34,48	1.08	2 (5%)	37,40,54	1.14	3 (8%)
17	CLA	B	814	-	43,51,73	1.77	7 (16%)	49,86,113	1.61	6 (12%)
17	CLA	4	315	14	41,49,73	1.85	6 (14%)	47,84,113	1.70	10 (21%)
17	CLA	B	821	-	55,63,73	1.60	7 (12%)	64,101,113	1.47	7 (10%)
17	CLA	4	306	-	65,73,73	1.48	7 (10%)	76,113,113	1.37	8 (10%)
17	CLA	A	811	-	41,49,73	1.83	8 (19%)	47,84,113	1.64	8 (17%)
20	BCR	B	847	-	41,41,41	0.74	0	56,56,56	2.24	18 (32%)
17	CLA	4	317	-	50,58,73	1.68	6 (12%)	58,95,113	1.51	9 (15%)
17	CLA	6	504	-	40,47,73	2.26	8 (20%)	49,81,113	1.74	9 (18%)
20	BCR	F	801	-	41,41,41	0.74	0	56,56,56	2.22	21 (37%)
17	CLA	1	515	-	39,48,73	1.93	7 (17%)	45,82,113	1.74	8 (17%)
17	CLA	B	831	-	56,64,73	1.59	7 (12%)	65,102,113	1.49	7 (10%)
17	CLA	B	834	-	43,51,73	1.77	6 (13%)	49,86,113	1.65	7 (14%)
17	CLA	1	504	-	41,49,73	1.84	7 (17%)	47,84,113	1.73	9 (19%)
17	CLA	A	805	-	46,54,73	1.73	6 (13%)	53,90,113	1.57	6 (11%)
17	CLA	A	807	-	42,50,73	1.81	6 (14%)	48,85,113	1.66	7 (14%)
17	CLA	4	311	-	46,54,73	1.78	7 (15%)	53,90,113	1.54	8 (15%)
17	CLA	4	312	-	41,49,73	1.82	6 (14%)	47,84,113	1.65	7 (14%)
17	CLA	A	852	-	57,65,73	1.57	7 (12%)	66,103,113	1.43	7 (10%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
17	CLA	K	203	-	41,49,73	1.80	6 (14%)	47,84,113	1.71	8 (17%)
20	BCR	A	846	-	41,41,41	0.81	1 (2%)	56,56,56	2.19	18 (32%)
20	BCR	1	503	-	10,11,41	0.69	0	15,16,56	1.94	4 (26%)
17	CLA	B	826	-	65,73,73	1.45	6 (9%)	76,113,113	1.46	8 (10%)
17	CLA	B	832	-	39,48,73	1.86	7 (17%)	45,82,113	1.76	8 (17%)
17	CLA	1	511	-	38,47,73	2.06	8 (21%)	46,80,113	1.72	10 (21%)
26	XAT	6	502	-	39,47,47	0.91	0	54,74,74	2.89	22 (40%)
17	CLA	B	813	-	54,62,73	1.68	6 (11%)	67,100,113	1.51	10 (14%)
17	CLA	3	304	-	40,46,73	2.85	9 (22%)	47,79,113	1.57	8 (17%)
19	LHG	A	844	-	39,39,48	1.00	2 (5%)	42,45,54	1.23	4 (9%)
17	CLA	6	509	-	50,58,73	1.71	5 (10%)	58,95,113	1.54	9 (15%)
20	BCR	K	204	-	41,41,41	0.70	0	56,56,56	2.09	18 (32%)
17	CLA	B	815	-	43,51,73	1.78	6 (13%)	49,86,113	1.51	6 (12%)
20	BCR	A	848	-	41,41,41	0.80	1 (2%)	56,56,56	2.15	20 (35%)
17	CLA	B	818	-	55,63,73	1.58	6 (10%)	64,101,113	1.46	7 (10%)
17	CLA	3	310	-	41,49,73	1.89	6 (14%)	47,84,113	1.65	7 (14%)
25	CHL	6	517	15	47,55,74	2.27	15 (31%)	50,91,114	2.81	19 (38%)
17	CLA	A	826	-	64,72,73	1.52	8 (12%)	79,112,113	1.50	9 (11%)
17	CLA	6	506	-	65,73,73	1.48	6 (9%)	76,113,113	1.37	6 (7%)
17	CLA	A	838	-	52,60,73	1.60	7 (13%)	60,97,113	1.57	8 (13%)
17	CLA	A	829	-	58,66,73	1.60	8 (13%)	67,104,113	1.46	7 (10%)
17	CLA	A	816	-	60,68,73	1.55	6 (10%)	70,107,113	1.41	6 (8%)
20	BCR	B	848	-	41,41,41	0.78	1 (2%)	56,56,56	2.04	16 (28%)
17	CLA	3	305	-	52,60,73	1.70	6 (11%)	60,97,113	1.52	10 (16%)
17	CLA	A	837	-	55,63,73	1.57	6 (10%)	64,101,113	1.54	10 (15%)
17	CLA	B	820	-	50,58,73	1.65	7 (14%)	58,95,113	1.57	8 (13%)
17	CLA	A	821	-	42,50,73	1.76	7 (16%)	48,85,113	1.64	6 (12%)
17	CLA	A	803	17	52,60,73	1.65	6 (11%)	60,97,113	1.55	9 (15%)
17	CLA	4	307	-	60,68,73	1.52	6 (10%)	70,107,113	1.47	8 (11%)
17	CLA	A	841	-	41,49,73	1.83	8 (19%)	47,84,113	1.66	8 (17%)
17	CLA	B	812	-	39,48,73	1.86	7 (17%)	45,82,113	1.77	8 (17%)
17	CLA	6	510	-	60,68,73	1.54	6 (10%)	70,107,113	1.38	6 (8%)
25	CHL	6	512	-	47,55,74	2.43	16 (34%)	50,91,114	2.76	19 (38%)
17	CLA	B	816	-	39,46,73	1.83	7 (17%)	44,79,113	1.62	5 (11%)
17	CLA	A	835	1	41,50,73	1.88	7 (17%)	49,85,113	1.54	7 (14%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
17	CLA	B	807	-	41,49,73	1.80	6 (14%)	47,84,113	1.69	9 (19%)
17	CLA	B	811	2	51,59,73	1.69	7 (13%)	59,96,113	1.52	10 (16%)
17	CLA	K	201	-	38,45,73	1.91	7 (18%)	43,78,113	1.64	7 (16%)
23	LMG	4	318	-	18,18,55	1.60	2 (11%)	22,23,63	1.61	2 (9%)
24	LUT	6	501	-	42,43,43	0.78	0	51,60,60	1.89	13 (25%)
17	CLA	A	814	-	42,50,73	1.81	7 (16%)	48,85,113	1.59	6 (12%)
22	DGD	J	104	-	67,67,67	0.84	2 (2%)	81,81,81	0.88	3 (3%)
17	CLA	A	815	-	45,53,73	1.81	6 (13%)	52,89,113	1.56	6 (11%)
17	CLA	4	305	-	50,58,73	1.71	6 (12%)	58,95,113	1.50	8 (13%)
24	LUT	1	501	-	42,43,43	0.80	0	51,60,60	1.97	13 (25%)
17	CLA	A	817	-	39,47,73	1.86	7 (17%)	42,81,113	1.73	8 (19%)
17	CLA	1	513	-	52,60,73	1.68	6 (11%)	60,97,113	1.51	8 (13%)
26	XAT	4	303	-	39,47,47	0.88	0	54,74,74	2.72	20 (37%)
17	CLA	3	309	-	50,58,73	1.67	5 (10%)	58,95,113	1.60	7 (12%)
24	LUT	3	301	-	42,43,43	0.79	0	51,60,60	2.00	11 (21%)
22	DGD	B	851	-	53,53,67	0.93	2 (3%)	67,67,81	1.06	3 (4%)
17	CLA	B	838	-	39,48,73	1.83	7 (17%)	45,82,113	1.71	8 (17%)
17	CLA	4	304	-	38,47,73	1.90	7 (18%)	48,81,113	1.93	10 (20%)
20	BCR	A	849	-	41,41,41	0.74	0	56,56,56	2.06	16 (28%)
20	BCR	A	850	-	41,41,41	0.81	2 (4%)	56,56,56	2.33	20 (35%)
21	SF4	C	101	-	0,12,12	-	-	-	-	-
17	CLA	B	808	-	54,62,73	1.58	6 (11%)	62,99,113	1.51	8 (12%)
17	CLA	3	315	-	45,53,73	1.76	7 (15%)	52,89,113	1.61	7 (13%)
18	PQN	A	842	-	15,15,34	1.44	2 (13%)	22,22,45	0.87	2 (9%)
19	LHG	B	852	17	37,37,48	1.05	2 (5%)	40,43,54	1.16	3 (7%)
17	CLA	A	822	-	41,49,73	1.82	7 (17%)	47,84,113	1.70	7 (14%)
17	CLA	B	824	-	42,50,73	1.81	7 (16%)	48,85,113	1.62	6 (12%)
17	CLA	B	817	-	43,51,73	1.77	7 (16%)	49,86,113	1.57	6 (12%)
17	CLA	A	831	-	56,64,73	1.56	6 (10%)	65,102,113	1.46	7 (10%)
17	CLA	A	812	-	37,47,73	1.85	7 (18%)	41,80,113	1.75	8 (19%)
17	CLA	B	842	-	41,49,73	1.83	8 (19%)	47,84,113	1.69	8 (17%)
17	CLA	B	823	-	41,49,73	1.84	8 (19%)	47,84,113	1.66	7 (14%)
17	CLA	B	829	-	50,58,73	1.64	6 (12%)	58,95,113	1.60	10 (17%)
20	BCR	A	854	-	41,41,41	0.70	0	56,56,56	2.26	19 (33%)
24	LUT	4	302	-	42,43,43	0.76	0	51,60,60	1.90	13 (25%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
25	CHL	4	314	-	51,59,74	2.14	15 (29%)	55,96,114	2.75	20 (36%)
17	CLA	B	827	-	41,48,73	1.96	8 (19%)	45,82,113	1.70	8 (17%)
17	CLA	A	836	-	51,59,73	1.65	6 (11%)	59,96,113	1.52	6 (10%)
17	CLA	B	822	-	41,49,73	1.85	6 (14%)	47,84,113	1.62	8 (17%)
17	CLA	B	835	-	45,53,73	1.74	7 (15%)	52,89,113	1.63	9 (17%)
20	BCR	B	849	-	41,41,41	0.79	1 (2%)	56,56,56	2.02	16 (28%)
20	BCR	A	851	-	41,41,41	0.71	0	56,56,56	2.16	19 (33%)
25	CHL	6	513	-	48,56,74	2.31	16 (33%)	51,92,114	2.77	22 (43%)
17	CLA	A	839	-	65,73,73	1.47	6 (9%)	76,113,113	1.42	9 (11%)
17	CLA	A	818	-	40,47,73	2.03	7 (17%)	44,81,113	2.14	12 (27%)
17	CLA	L	302	-	40,48,73	1.93	6 (15%)	50,83,113	1.88	9 (18%)
20	BCR	L	301	-	41,41,41	0.75	1 (2%)	56,56,56	2.04	17 (30%)
17	CLA	B	805	-	37,46,73	1.83	7 (18%)	40,79,113	1.88	7 (17%)
17	CLA	B	801	-	64,72,73	1.50	8 (12%)	79,112,113	1.48	8 (10%)
20	BCR	3	303	-	41,41,41	0.79	1 (2%)	56,56,56	2.00	18 (32%)
20	BCR	B	846	-	41,41,41	0.73	0	56,56,56	1.96	14 (25%)
17	CLA	A	819	-	45,53,73	1.76	6 (13%)	52,89,113	1.61	7 (13%)
17	CLA	B	825	-	42,50,73	1.79	7 (16%)	48,85,113	1.59	6 (12%)
17	CLA	J	101	-	41,49,73	1.93	7 (17%)	51,84,113	1.55	6 (11%)
17	CLA	4	309	-	50,58,73	1.66	6 (12%)	58,95,113	1.54	7 (12%)
20	BCR	A	847	-	41,41,41	0.76	0	56,56,56	1.90	17 (30%)
25	CHL	6	515	-	40,49,74	2.39	14 (35%)	42,83,114	2.99	19 (45%)
17	CLA	3	312	13	60,68,73	1.54	6 (10%)	70,107,113	1.43	7 (10%)
17	CLA	B	803	-	40,49,73	1.73	6 (15%)	44,83,113	1.64	6 (13%)
17	CLA	3	314	-	39,48,73	1.84	6 (15%)	45,82,113	1.84	8 (17%)
17	CLA	A	806	-	41,49,73	1.81	7 (17%)	47,84,113	1.70	9 (19%)
25	CHL	1	517	-	43,51,74	2.37	15 (34%)	45,86,114	2.90	19 (42%)
25	CHL	4	316	-	43,51,74	2.30	14 (32%)	45,86,114	2.85	19 (42%)
17	CLA	A	830	-	50,58,73	1.69	8 (16%)	58,95,113	1.51	8 (13%)
17	CLA	A	820	-	41,49,73	1.78	6 (14%)	47,84,113	1.71	7 (14%)
17	CLA	6	505	-	52,60,73	1.66	7 (13%)	60,97,113	1.51	6 (10%)
18	PQN	B	844	-	17,17,34	2.13	2 (11%)	21,24,45	1.20	3 (14%)
17	CLA	1	510	-	46,54,73	1.76	7 (15%)	53,90,113	1.49	6 (11%)
17	CLA	1	507	12	40,48,73	1.88	7 (17%)	50,83,113	1.74	10 (20%)

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
17	CLA	A	833	-	1/1/10/20	4/8/86/115	-
17	CLA	A	834	-	1/1/11/20	5/13/91/115	-
17	CLA	B	837	-	1/1/10/20	3/10/88/115	-
17	CLA	1	508	-	1/1/15/20	17/37/115/115	-
17	CLA	B	806	-	1/1/10/20	2/8/86/115	-
17	CLA	A	802	-	1/1/9/20	4/8/82/115	-
17	CLA	1	506	-	1/1/13/20	10/25/103/115	-
19	LHG	1	516	-	-	12/53/53/53	-
17	CLA	A	813	-	1/1/10/20	4/10/88/115	-
17	CLA	J	102	-	1/1/10/20	4/10/88/115	-
17	CLA	K	202	-	1/1/11/20	8/13/91/115	-
17	CLA	L	303	-	1/1/14/20	12/31/109/115	-
25	CHL	1	514	-	3/3/16/26	7/17/115/137	-
20	BCR	B	850	-	-	2/27/61/63	0/2/2/2
17	CLA	A	809	-	1/1/12/20	4/19/97/115	-
17	CLA	B	828	-	1/1/10/20	4/8/86/115	-
17	CLA	B	841	-	1/1/10/20	3/8/86/115	-
23	LMG	J	105	-	-	5/25/45/70	0/1/1/1
17	CLA	A	825	-	1/1/10/20	0/8/86/115	-
17	CLA	6	508	-	1/1/10/20	2/6/84/115	-
20	BCR	L	305	-	-	2/29/63/63	0/2/2/2
17	CLA	A	828	-	1/1/10/20	5/10/88/115	-
17	CLA	3	306	-	1/1/10/20	3/8/86/115	-
17	CLA	A	840	-	1/1/10/20	1/8/86/115	-
17	CLA	A	824	-	1/1/13/20	9/27/105/115	-
17	CLA	A	853	-	1/1/11/20	8/13/91/115	-
17	CLA	A	827	-	1/1/10/20	1/8/86/115	-
17	CLA	B	804	-	1/1/9/20	4/10/84/115	-
17	CLA	B	810	-	1/1/15/20	10/37/115/115	-
17	CLA	3	308	-	1/1/10/20	5/10/88/115	-
17	CLA	1	505	-	-	5/15/93/115	-
19	LHG	A	845	-	-	10/28/28/53	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
17	CLA	A	832	-	1/1/10/20	6/11/89/115	-
17	CLA	B	836	-	1/1/12/20	11/29/103/115	-
17	CLA	4	308	-	1/1/10/20	4/8/86/115	-
24	LUT	1	502	-	1/1/12/27	9/29/67/67	0/2/2/2
25	CHL	4	313	-	3/3/16/26	5/17/115/137	-
17	CLA	B	843	19	1/1/9/20	0/2/80/115	-
17	CLA	B	830	-	1/1/9/20	3/8/80/115	-
17	CLA	6	514	-	1/1/9/20	2/8/82/115	-
20	BCR	F	803	-	-	2/29/63/63	0/2/2/2
17	CLA	6	511	-	1/1/12/20	7/19/97/115	-
17	CLA	A	804	-	1/1/15/20	16/37/115/115	-
17	CLA	6	507	-	1/1/12/20	8/21/99/115	-
17	CLA	K	205	-	1/1/9/20	0/6/80/115	-
25	CHL	3	313	-	3/3/15/26	11/17/111/137	-
20	BCR	B	845	-	-	7/29/63/63	0/2/2/2
17	CLA	A	810	17	1/1/13/20	10/25/103/115	-
20	BCR	I	102	-	-	8/29/63/63	0/2/2/2
20	BCR	J	103	-	-	2/29/63/63	0/2/2/2
24	LUT	3	302	-	-	6/29/67/67	0/2/2/2
17	CLA	3	311	-	1/1/11/20	8/17/95/115	-
17	CLA	B	819	-	1/1/10/20	2/8/86/115	-
17	CLA	L	304	-	1/1/11/20	6/13/91/115	-
17	CLA	4	310	-	1/1/10/20	3/10/88/115	-
21	SF4	B	802	-	-	-	0/6/5/5
17	CLA	1	509	-	1/1/12/20	8/19/97/115	-
17	CLA	B	840	-	1/1/10/20	4/8/86/115	-
20	BCR	6	503	-	-	7/29/63/63	0/2/2/2
17	CLA	B	839	-	1/1/10/20	4/8/86/115	-
17	CLA	A	823	-	1/1/10/20	3/11/87/115	-
17	CLA	3	307	13	-	3/8/82/115	-
17	CLA	A	843	-	1/1/12/20	9/19/95/115	-
17	CLA	B	833	-	1/1/10/20	3/11/89/115	-
25	CHL	1	512	-	3/3/16/26	3/17/115/137	-
17	CLA	3	316	-	1/1/11/20	7/15/93/115	-
20	BCR	4	301	-	-	6/29/63/63	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
17	CLA	B	809	-	1/1/12/20	3/22/100/115	-
16	CL0	A	801	-	3/3/18/25	5/33/125/135	-
21	SF4	C	102	-	-	-	0/6/5/5
17	CLA	A	808	1	1/1/10/20	4/8/86/115	-
20	BCR	I	101	-	-	6/29/63/63	0/2/2/2
17	CLA	F	802	-	1/1/10/20	2/8/86/115	-
19	LHG	6	516	-	-	6/39/39/53	-
17	CLA	B	814	-	1/1/10/20	4/11/89/115	-
17	CLA	4	315	14	1/1/10/20	2/8/86/115	-
17	CLA	B	821	-	1/1/13/20	8/25/103/115	-
17	CLA	4	306	-	1/1/15/20	16/37/115/115	-
17	CLA	A	811	-	1/1/10/20	4/8/86/115	-
20	BCR	B	847	-	-	6/29/63/63	0/2/2/2
17	CLA	4	317	-	1/1/12/20	7/19/97/115	-
17	CLA	6	504	-	1/1/10/20	3/8/84/115	-
20	BCR	F	801	-	-	2/29/63/63	0/2/2/2
17	CLA	1	515	-	1/1/9/20	3/8/82/115	-
17	CLA	B	831	-	1/1/13/20	5/27/105/115	-
17	CLA	B	834	-	1/1/10/20	3/11/89/115	-
17	CLA	1	504	-	1/1/10/20	5/8/86/115	-
17	CLA	A	805	-	1/1/11/20	4/15/93/115	-
17	CLA	A	807	-	1/1/10/20	2/10/88/115	-
17	CLA	4	311	-	1/1/11/20	4/15/93/115	-
17	CLA	4	312	-	1/1/10/20	3/8/86/115	-
17	CLA	A	852	-	1/1/13/20	9/28/106/115	-
17	CLA	K	203	-	1/1/10/20	4/8/86/115	-
20	BCR	A	846	-	-	4/29/63/63	0/2/2/2
20	BCR	1	503	-	-	1/1/19/63	0/1/1/2
17	CLA	B	826	-	1/1/15/20	11/37/115/115	-
17	CLA	B	832	-	1/1/9/20	2/8/82/115	-
17	CLA	1	511	-	1/1/8/20	2/8/79/115	-
26	XAT	6	502	-	-	6/31/93/93	0/4/4/4
17	CLA	B	813	-	1/1/13/20	10/25/101/115	-
17	CLA	3	304	-	1/1/10/20	0/6/84/115	-
19	LHG	A	844	-	-	8/44/44/53	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
17	CLA	6	509	-	1/1/12/20	6/19/97/115	-
20	BCR	K	204	-	-	9/29/63/63	0/2/2/2
17	CLA	B	815	-	1/1/10/20	4/11/89/115	-
20	BCR	A	848	-	-	6/29/63/63	0/2/2/2
17	CLA	B	818	-	1/1/13/20	9/25/103/115	-
17	CLA	3	310	-	1/1/10/20	3/8/86/115	-
25	CHL	6	517	15	3/3/16/26	3/17/115/137	-
17	CLA	A	826	-	1/1/15/20	18/37/113/115	-
17	CLA	6	506	-	1/1/15/20	19/37/115/115	-
17	CLA	A	838	-	1/1/12/20	4/22/100/115	-
17	CLA	A	829	-	1/1/13/20	13/29/107/115	-
17	CLA	A	816	-	1/1/14/20	11/31/109/115	-
20	BCR	B	848	-	-	0/29/63/63	0/2/2/2
17	CLA	3	305	-	1/1/12/20	7/22/100/115	-
17	CLA	A	837	-	1/1/13/20	11/25/103/115	-
17	CLA	B	820	-	1/1/12/20	11/19/97/115	-
17	CLA	A	821	-	1/1/10/20	5/10/88/115	-
17	CLA	A	803	17	1/1/12/20	5/22/100/115	-
17	CLA	4	307	-	1/1/14/20	12/31/109/115	-
17	CLA	A	841	-	1/1/10/20	2/8/86/115	-
17	CLA	B	812	-	1/1/9/20	4/8/82/115	-
17	CLA	6	510	-	1/1/14/20	11/31/109/115	-
25	CHL	6	512	-	3/3/16/26	5/17/115/137	-
17	CLA	B	816	-	1/1/8/20	1/4/78/115	-
17	CLA	A	835	1	1/1/10/20	2/7/85/115	-
17	CLA	B	807	-	1/1/10/20	2/8/86/115	-
17	CLA	B	811	2	1/1/12/20	6/21/99/115	-
17	CLA	K	201	-	1/1/8/20	0/2/76/115	-
23	LMG	4	318	-	-	11/11/28/70	0/1/1/1
24	LUT	6	501	-	-	4/29/67/67	0/2/2/2
17	CLA	A	814	-	1/1/10/20	5/10/88/115	-
22	DGD	J	104	-	-	14/55/95/95	0/2/2/2
17	CLA	A	815	-	1/1/11/20	7/13/91/115	-
17	CLA	4	305	-	1/1/12/20	6/19/97/115	-
24	LUT	1	501	-	-	2/29/67/67	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
17	CLA	A	817	-	1/1/9/20	0/2/82/115	-
17	CLA	1	513	-	1/1/12/20	5/22/100/115	-
26	XAT	4	303	-	-	2/31/93/93	0/4/4/4
17	CLA	3	309	-	1/1/12/20	5/19/97/115	-
24	LUT	3	301	-	-	6/29/67/67	0/2/2/2
22	DGD	B	851	-	-	9/41/81/95	0/2/2/2
17	CLA	B	838	-	1/1/9/20	5/8/82/115	-
17	CLA	4	304	-	1/1/9/20	3/8/80/115	-
20	BCR	A	849	-	-	7/29/63/63	0/2/2/2
20	BCR	A	850	-	-	6/29/63/63	0/2/2/2
21	SF4	C	101	-	-	-	0/6/5/5
17	CLA	B	808	-	1/1/12/20	6/23/101/115	-
17	CLA	3	315	-	-	8/13/91/115	-
19	LHG	B	852	17	-	13/42/42/53	-
18	PQN	A	842	-	-	-	0/2/2/2
17	CLA	A	822	-	1/1/10/20	1/8/86/115	-
17	CLA	B	824	-	1/1/10/20	4/10/88/115	-
17	CLA	B	817	-	1/1/10/20	5/11/89/115	-
17	CLA	A	831	-	1/1/13/20	7/27/105/115	-
17	CLA	A	812	-	1/1/8/20	3/5/79/115	-
17	CLA	B	842	-	1/1/10/20	0/8/86/115	-
17	CLA	B	823	-	1/1/10/20	2/8/86/115	-
17	CLA	B	829	-	1/1/12/20	9/19/97/115	-
25	CHL	4	314	-	3/3/17/26	4/21/119/137	-
20	BCR	A	854	-	-	4/29/63/63	0/2/2/2
24	LUT	4	302	-	-	2/29/67/67	0/2/2/2
17	CLA	B	827	-	1/1/9/20	1/8/82/115	-
17	CLA	A	836	-	1/1/12/20	6/21/99/115	-
17	CLA	B	822	-	1/1/10/20	0/8/86/115	-
17	CLA	B	835	-	1/1/11/20	8/13/91/115	-
20	BCR	B	849	-	-	0/29/63/63	0/2/2/2
20	BCR	A	851	-	-	4/29/63/63	0/2/2/2
25	CHL	6	513	-	3/3/16/26	5/18/116/137	-
17	CLA	A	839	-	1/1/15/20	16/37/115/115	-
17	CLA	A	818	-	1/1/9/20	4/6/80/115	-
17	CLA	L	302	-	1/1/10/20	2/8/84/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
20	BCR	L	301	-	-	2/29/63/63	0/2/2/2
17	CLA	B	805	-	1/1/8/20	0/2/78/115	-
17	CLA	B	801	-	1/1/15/20	13/37/113/115	-
20	BCR	3	303	-	-	7/29/63/63	0/2/2/2
20	BCR	B	846	-	-	7/29/63/63	0/2/2/2
17	CLA	A	819	-	1/1/11/20	5/13/91/115	-
17	CLA	B	825	-	1/1/10/20	3/10/88/115	-
17	CLA	J	101	-	1/1/10/20	6/10/86/115	-
17	CLA	4	309	-	1/1/12/20	7/19/97/115	-
20	BCR	A	847	-	-	0/29/63/63	0/2/2/2
25	CHL	6	515	-	3/3/14/26	4/10/104/137	-
17	CLA	3	312	13	1/1/14/20	12/31/109/115	-
17	CLA	B	803	-	1/1/10/20	3/10/88/115	-
17	CLA	3	314	-	1/1/9/20	2/8/82/115	-
17	CLA	A	806	-	1/1/10/20	1/8/86/115	-
25	CHL	1	517	-	3/3/15/26	5/12/110/137	-
25	CHL	4	316	-	3/3/15/26	3/12/110/137	-
17	CLA	A	830	-	1/1/12/20	6/19/97/115	-
17	CLA	A	820	-	1/1/10/20	3/8/86/115	-
17	CLA	6	505	-	-	8/22/100/115	-
18	PQN	B	844	-	-	0/3/23/43	0/2/2/2
17	CLA	1	510	-	1/1/11/20	7/15/93/115	-
17	CLA	1	507	12	1/1/10/20	4/8/84/115	-

The worst 5 of 1113 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
24	1	502	LUT	C24-C25	14.61	1.51	1.33
17	3	304	CLA	C3B-C4B	9.98	1.49	1.43
17	6	504	CLA	C3B-C4B	8.58	1.49	1.39
17	3	304	CLA	C2B-C1B	8.49	1.49	1.39
17	B	804	CLA	CHB-C4A	8.43	1.41	1.34

The worst 5 of 1884 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
20	6	503	BCR	C40-C30-C25	-12.39	90.20	110.30
16	A	801	CL0	C4A-NA-C1A	10.83	111.58	106.71

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	B	804	CLA	C4A-NA-C1A	9.99	111.20	106.71
20	6	503	BCR	C20-C21-C22	-9.47	113.80	127.31
17	1	509	CLA	C4A-NA-C1A	8.63	110.58	106.71

5 of 170 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
16	A	801	CL0	ND
16	A	801	CL0	NC
16	A	801	CL0	NA
17	A	802	CLA	ND
17	A	803	CLA	ND

5 of 1040 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
16	A	801	CL0	C1A-C2A-CAA-CBA
16	A	801	CL0	CHA-CBD-CGD-O2D
17	A	802	CLA	CAD-CBD-CGD-O2D
17	A	804	CLA	C1A-C2A-CAA-CBA
17	A	805	CLA	CBD-CGD-O2D-CED

There are no ring outliers.

169 monomers are involved in 416 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
17	A	833	CLA	3	0
17	A	834	CLA	1	0
17	B	806	CLA	2	0
17	A	802	CLA	1	0
17	1	506	CLA	6	0
19	1	516	LHG	7	0
17	A	813	CLA	2	0
17	K	202	CLA	1	0
17	L	303	CLA	2	0
20	B	850	BCR	4	0
25	1	514	CHL	3	0
17	A	809	CLA	4	0
17	B	828	CLA	1	0
23	J	105	LMG	1	0
17	A	825	CLA	3	0
17	6	508	CLA	1	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	L	305	BCR	2	0
17	A	828	CLA	2	0
17	3	306	CLA	3	0
17	A	824	CLA	1	0
17	A	853	CLA	3	0
17	A	827	CLA	3	0
17	B	804	CLA	6	0
17	B	810	CLA	3	0
17	1	505	CLA	2	0
17	3	308	CLA	1	0
19	A	845	LHG	1	0
17	A	832	CLA	1	0
17	4	308	CLA	2	0
24	1	502	LUT	6	0
25	4	313	CHL	1	0
17	B	843	CLA	1	0
17	6	514	CLA	2	0
20	F	803	BCR	1	0
17	6	511	CLA	6	0
17	A	804	CLA	4	0
20	B	845	BCR	1	0
20	I	102	BCR	4	0
17	A	810	CLA	7	0
20	J	103	BCR	3	0
25	3	313	CHL	1	0
24	3	302	LUT	6	0
17	3	311	CLA	2	0
17	B	819	CLA	4	0
17	L	304	CLA	4	0
17	4	310	CLA	1	0
17	1	509	CLA	3	0
17	B	840	CLA	3	0
20	6	503	BCR	4	0
17	B	839	CLA	1	0
17	3	307	CLA	3	0
17	A	843	CLA	2	0
17	B	833	CLA	3	0
25	1	512	CHL	1	0
17	3	316	CLA	1	0
20	4	301	BCR	5	0
16	A	801	CL0	6	0
17	A	808	CLA	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
20	I	101	BCR	1	0
19	6	516	LHG	1	0
17	4	315	CLA	2	0
17	B	821	CLA	6	0
17	4	306	CLA	7	0
20	B	847	BCR	1	0
17	4	317	CLA	1	0
17	6	504	CLA	1	0
20	F	801	BCR	1	0
17	1	515	CLA	4	0
17	B	831	CLA	2	0
17	B	834	CLA	3	0
17	1	504	CLA	8	0
17	A	805	CLA	1	0
17	A	807	CLA	3	0
17	4	311	CLA	1	0
17	4	312	CLA	1	0
17	A	852	CLA	6	0
17	K	203	CLA	1	0
20	A	846	BCR	5	0
20	1	503	BCR	1	0
17	B	826	CLA	6	0
17	B	832	CLA	4	0
17	1	511	CLA	1	0
26	6	502	XAT	9	0
17	B	813	CLA	4	0
17	3	304	CLA	3	0
19	A	844	LHG	1	0
17	6	509	CLA	2	0
20	K	204	BCR	1	0
17	B	815	CLA	2	0
20	A	848	BCR	2	0
17	B	818	CLA	4	0
17	3	310	CLA	1	0
25	6	517	CHL	1	0
17	A	826	CLA	3	0
17	6	506	CLA	7	0
17	A	838	CLA	7	0
17	A	829	CLA	1	0
17	A	816	CLA	1	0
20	B	848	BCR	3	0
17	3	305	CLA	1	0

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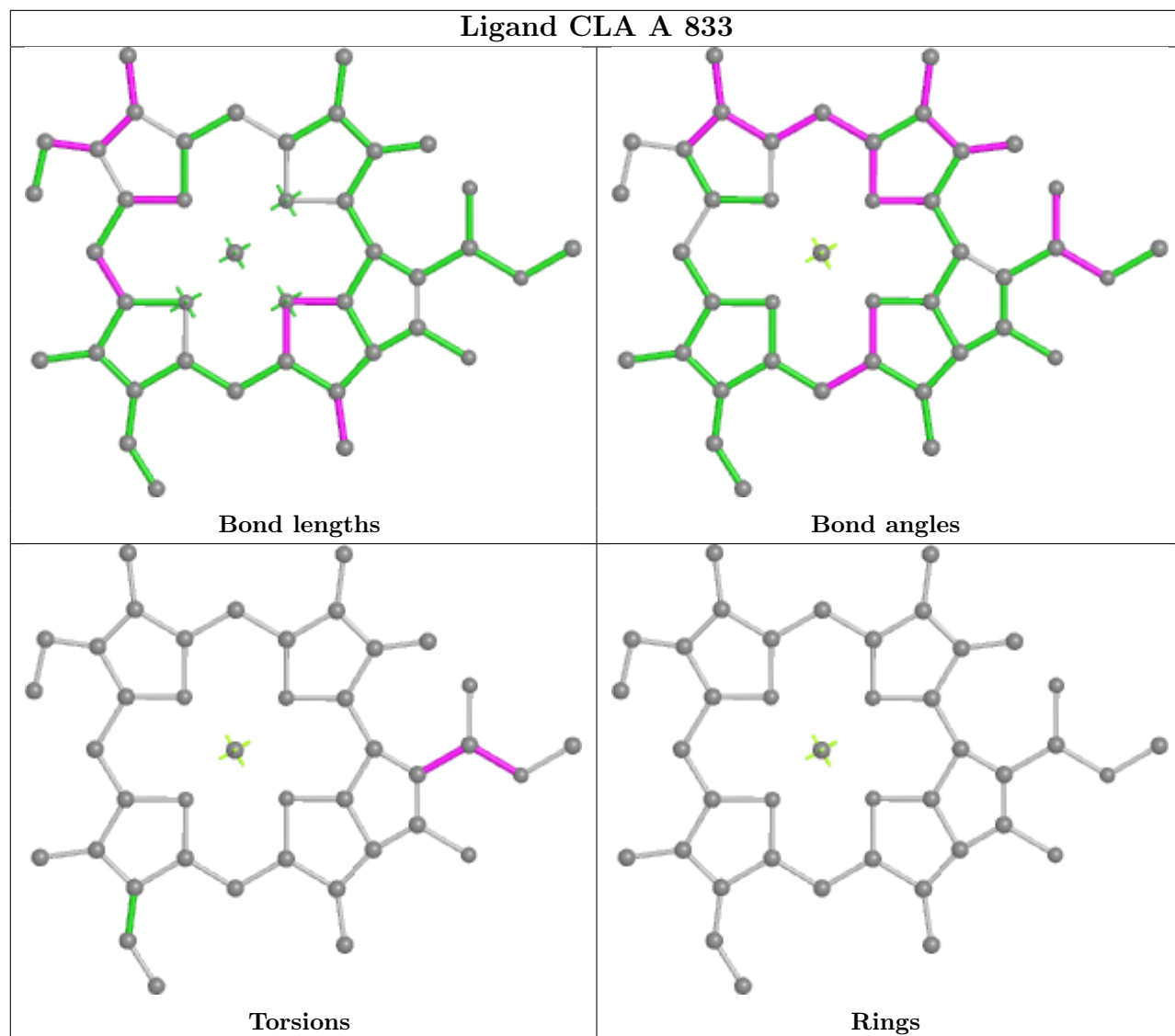
Mol	Chain	Res	Type	Clashes	Symm-Clashes
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17	B	820	CLA	1	0
17	A	821	CLA	2	0
17	A	803	CLA	4	0
17	4	307	CLA	5	0
17	6	510	CLA	6	0
25	6	512	CHL	4	0
17	B	816	CLA	4	0
17	B	807	CLA	5	0
17	B	811	CLA	2	0
17	K	201	CLA	1	0
24	6	501	LUT	6	0
17	A	814	CLA	1	0
22	J	104	DGD	2	0
17	A	815	CLA	1	0
17	4	305	CLA	2	0
24	1	501	LUT	6	0
17	A	817	CLA	2	0
17	1	513	CLA	1	0
26	4	303	XAT	3	0
17	3	309	CLA	1	0
24	3	301	LUT	5	0
22	B	851	DGD	2	0
17	B	838	CLA	3	0
17	4	304	CLA	2	0
20	A	849	BCR	4	0
20	A	850	BCR	2	0
21	C	101	SF4	1	0
17	B	808	CLA	7	0
17	3	315	CLA	6	0
18	A	842	PQN	2	0
19	B	852	LHG	1	0
17	A	812	CLA	5	0
17	B	842	CLA	1	0
17	B	823	CLA	1	0
17	B	829	CLA	1	0
20	A	854	BCR	4	0
24	4	302	LUT	9	0
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17	A	836	CLA	2	0
17	B	835	CLA	6	0
20	B	849	BCR	4	0

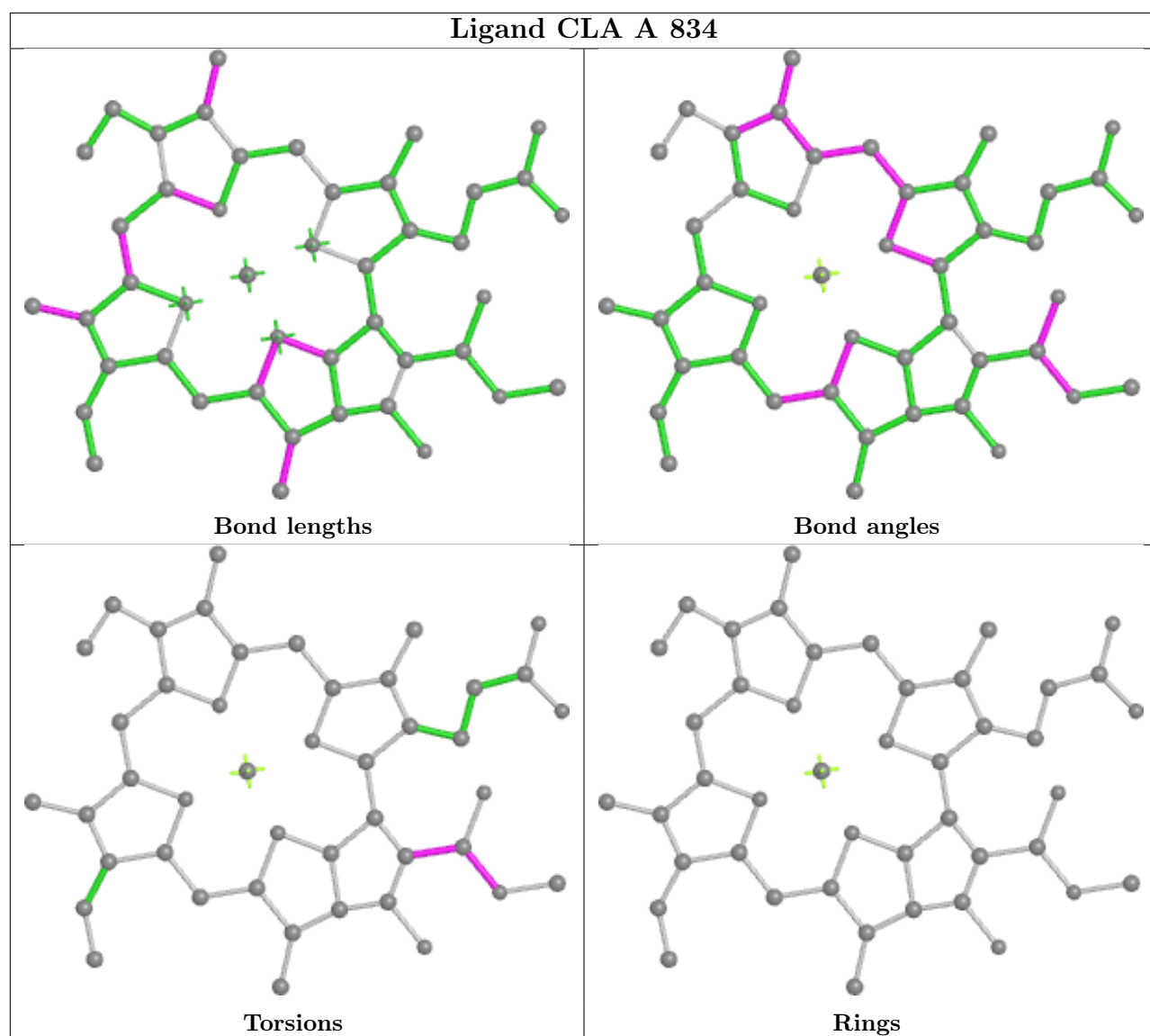
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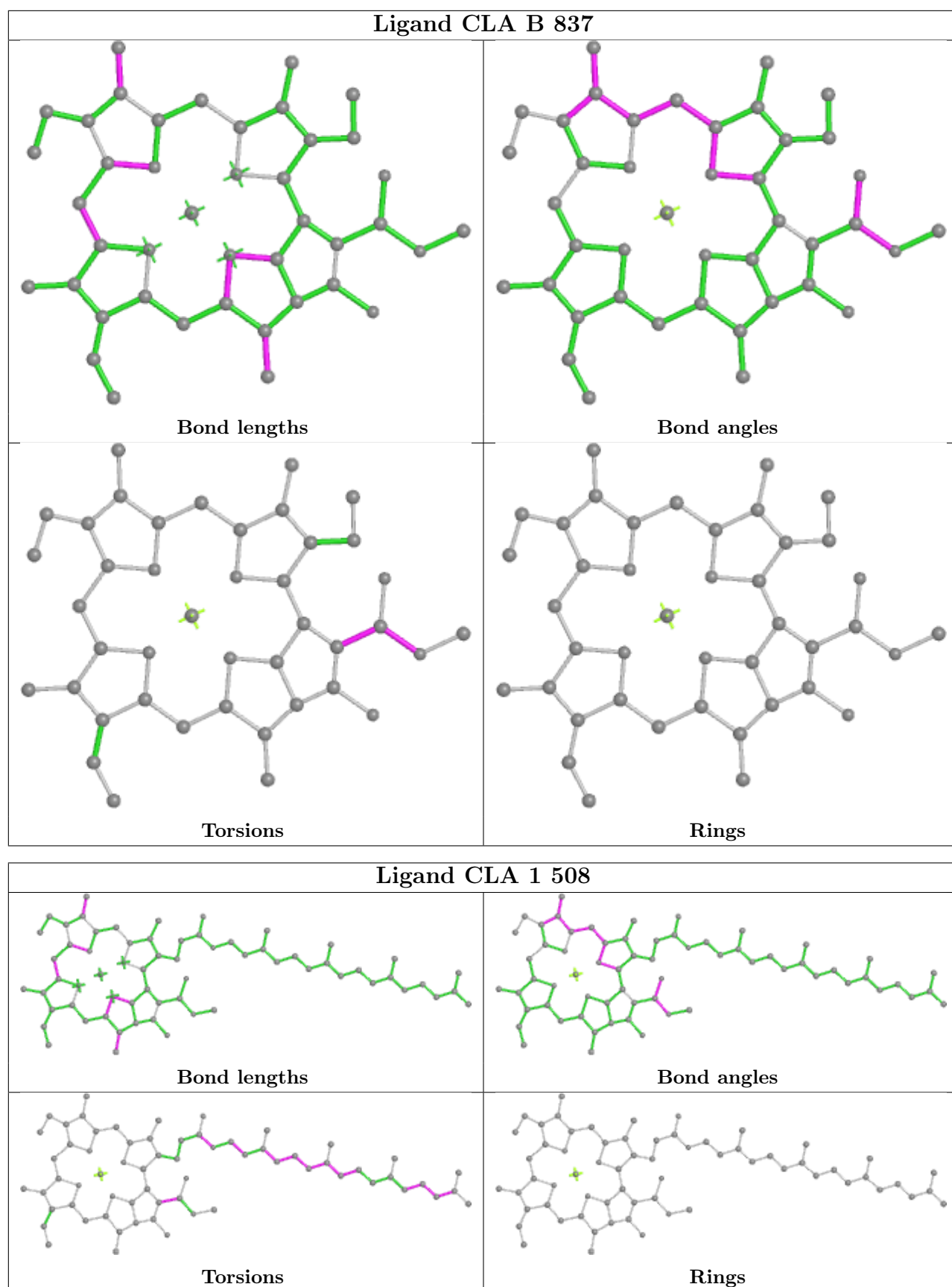
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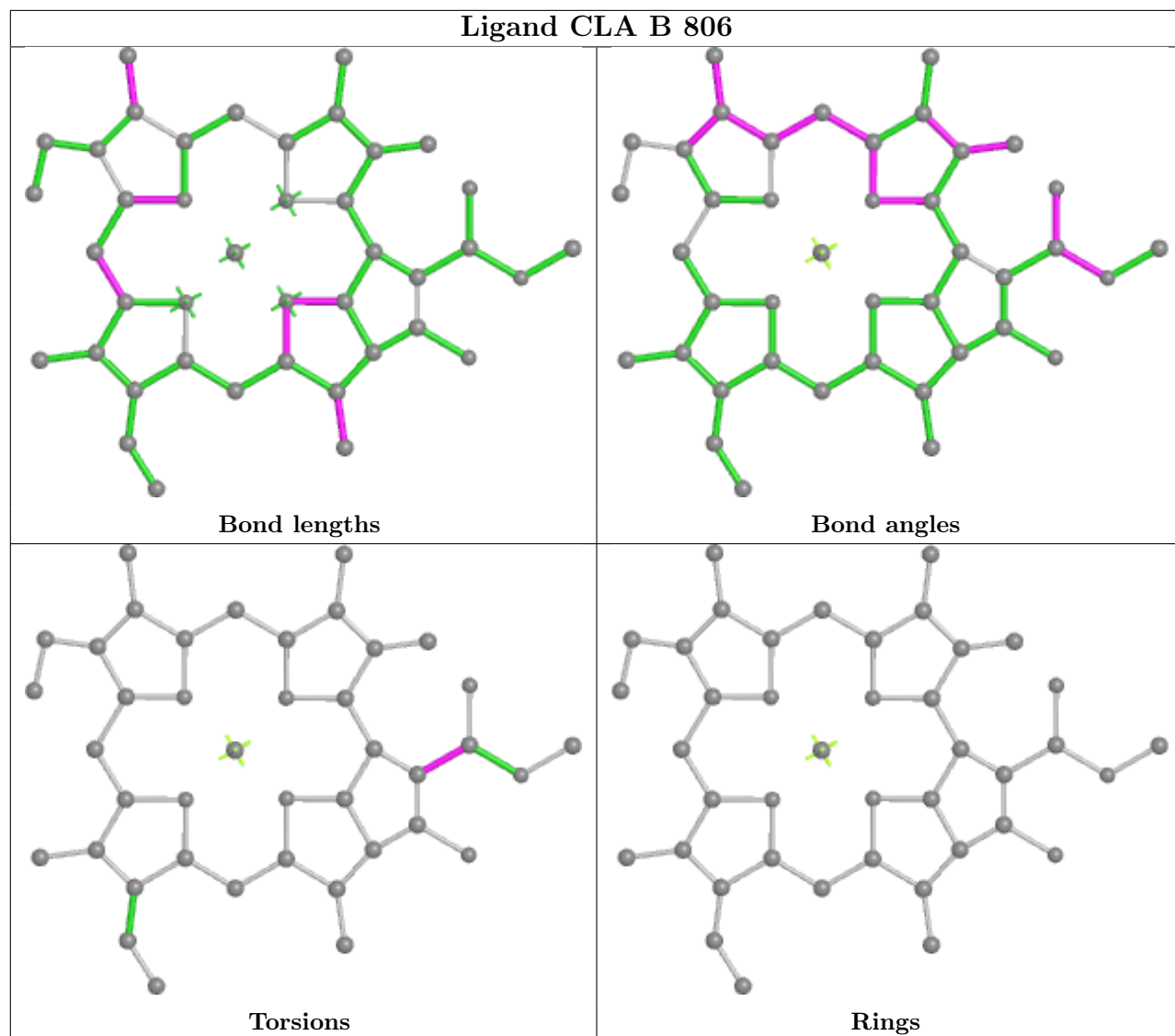
Mol	Chain	Res	Type	Clashes	Symm-Clashes
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25	6	513	CHL	2	0
17	A	839	CLA	5	0
17	A	818	CLA	4	0
17	L	302	CLA	1	0
20	L	301	BCR	1	0
17	B	805	CLA	2	0
17	B	801	CLA	4	0
20	3	303	BCR	3	0
17	A	819	CLA	1	0
17	B	825	CLA	1	0
17	J	101	CLA	1	0
17	4	309	CLA	1	0
20	A	847	BCR	2	0
25	6	515	CHL	1	0
17	3	312	CLA	4	0
17	B	803	CLA	4	0
17	3	314	CLA	4	0
17	A	806	CLA	1	0
25	1	517	CHL	2	0
25	4	316	CHL	2	0
17	A	830	CLA	3	0
17	A	820	CLA	2	0
17	6	505	CLA	3	0
18	B	844	PQN	7	0
17	1	510	CLA	2	0
17	1	507	CLA	2	0

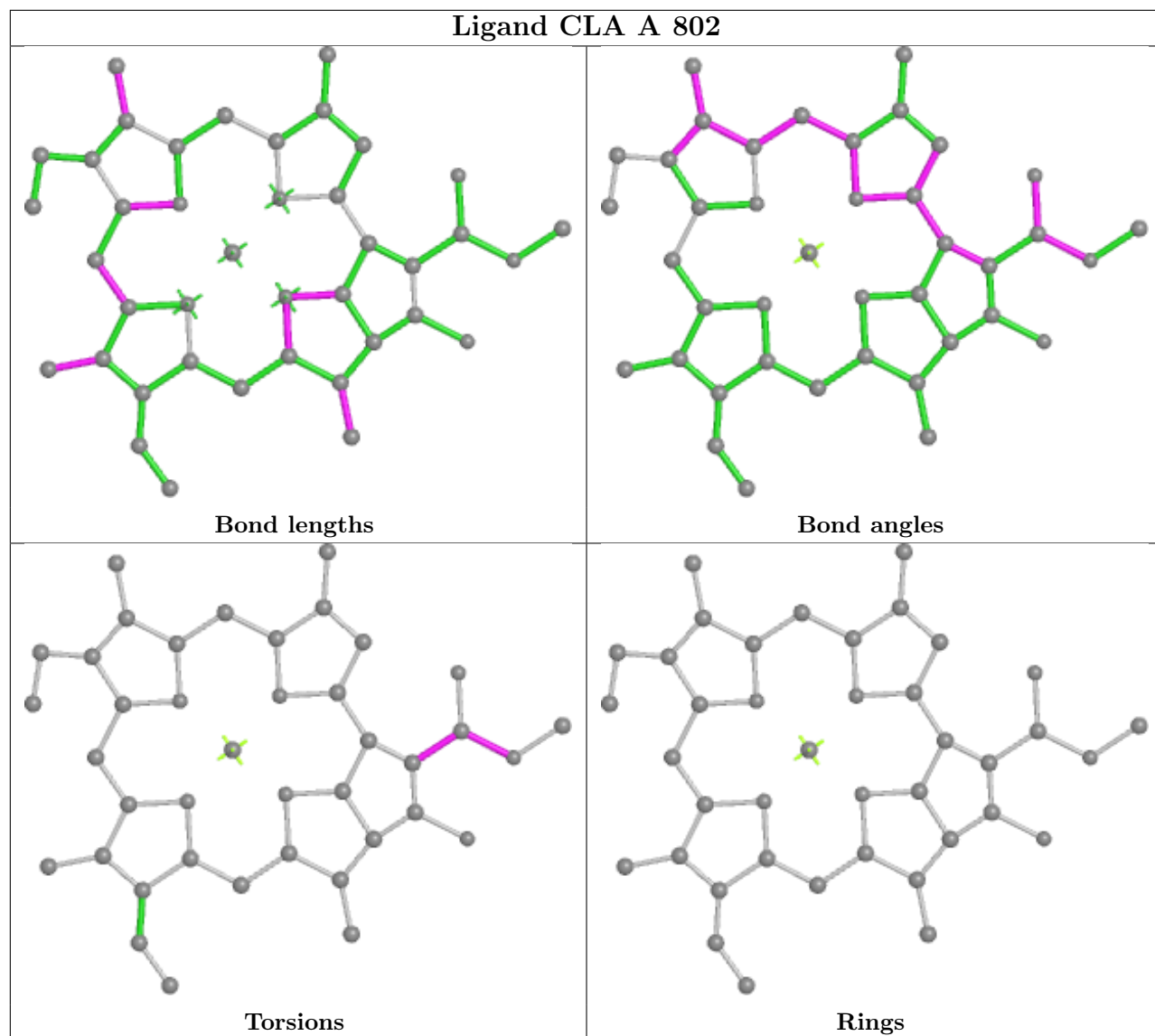
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

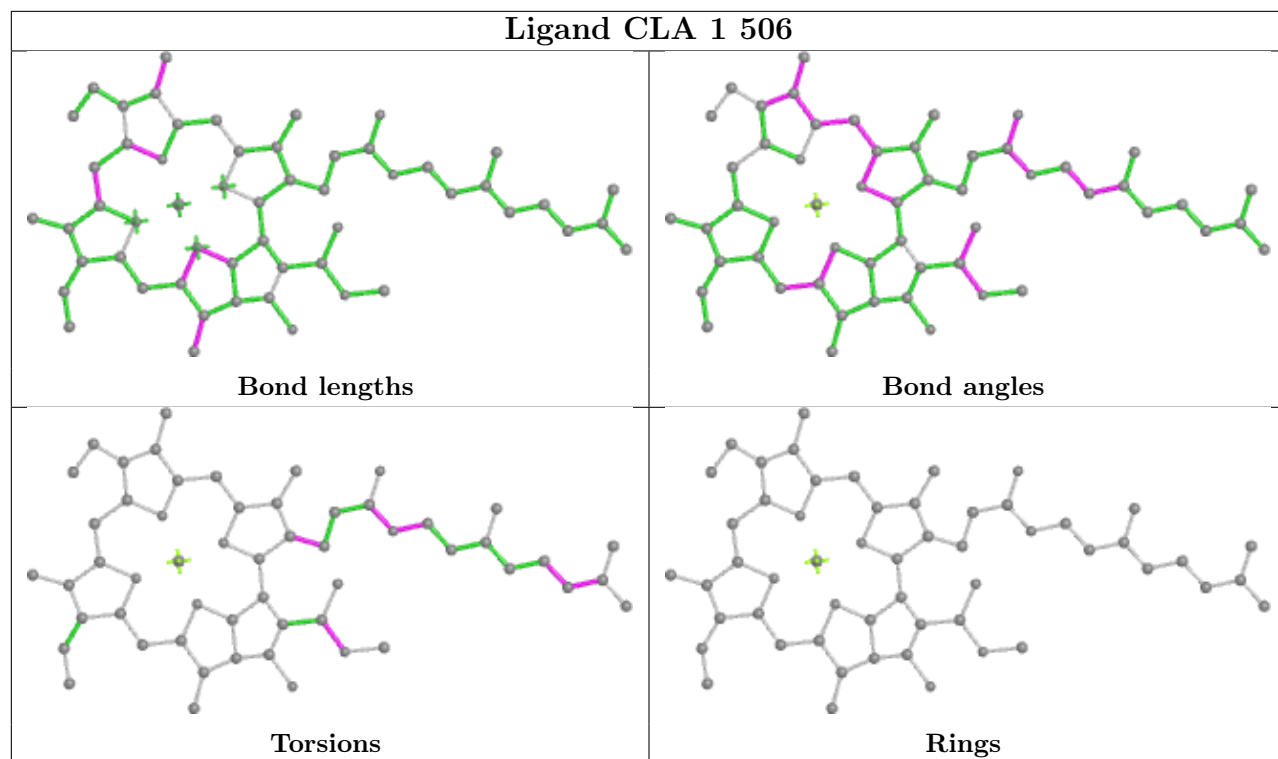


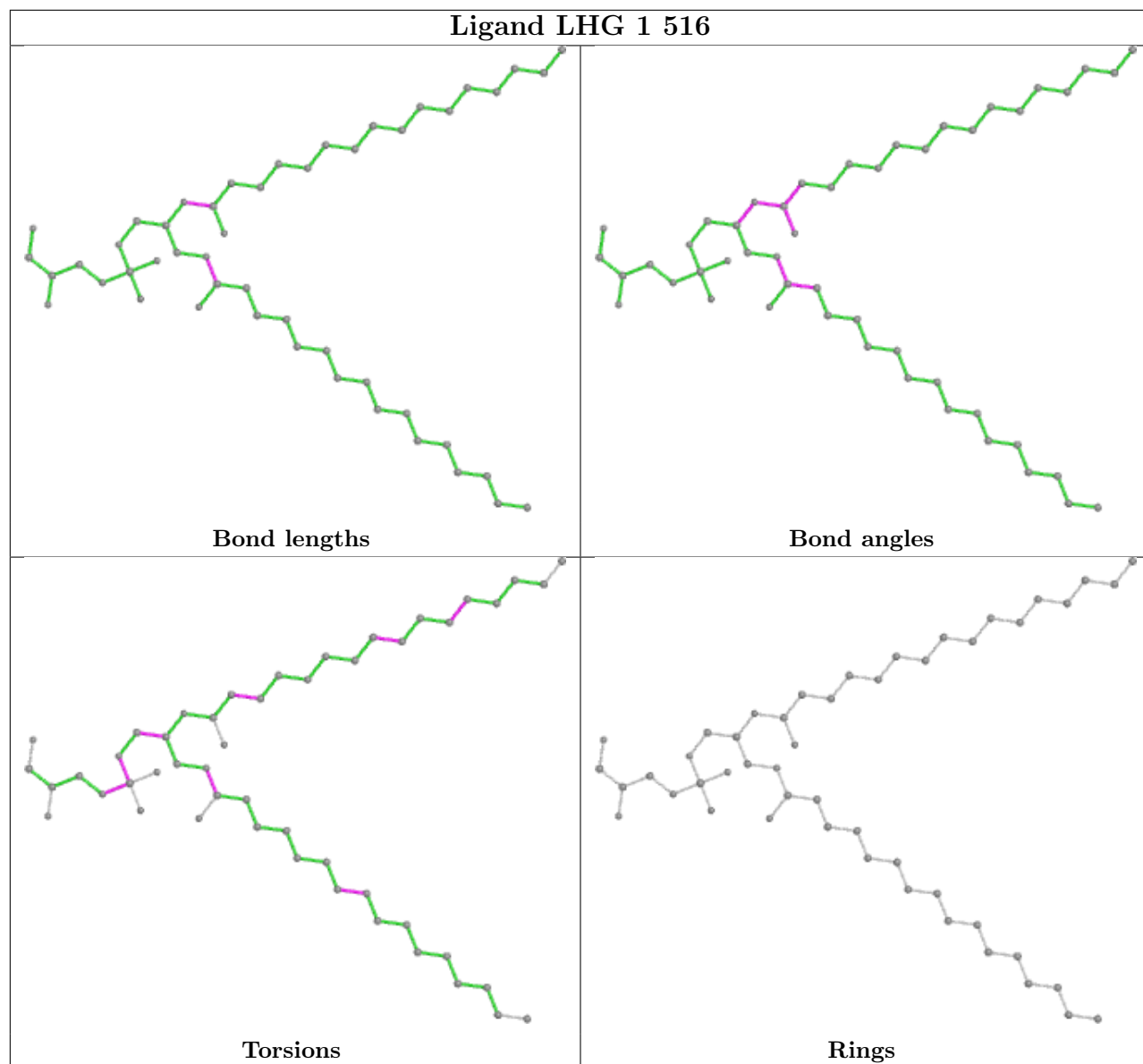


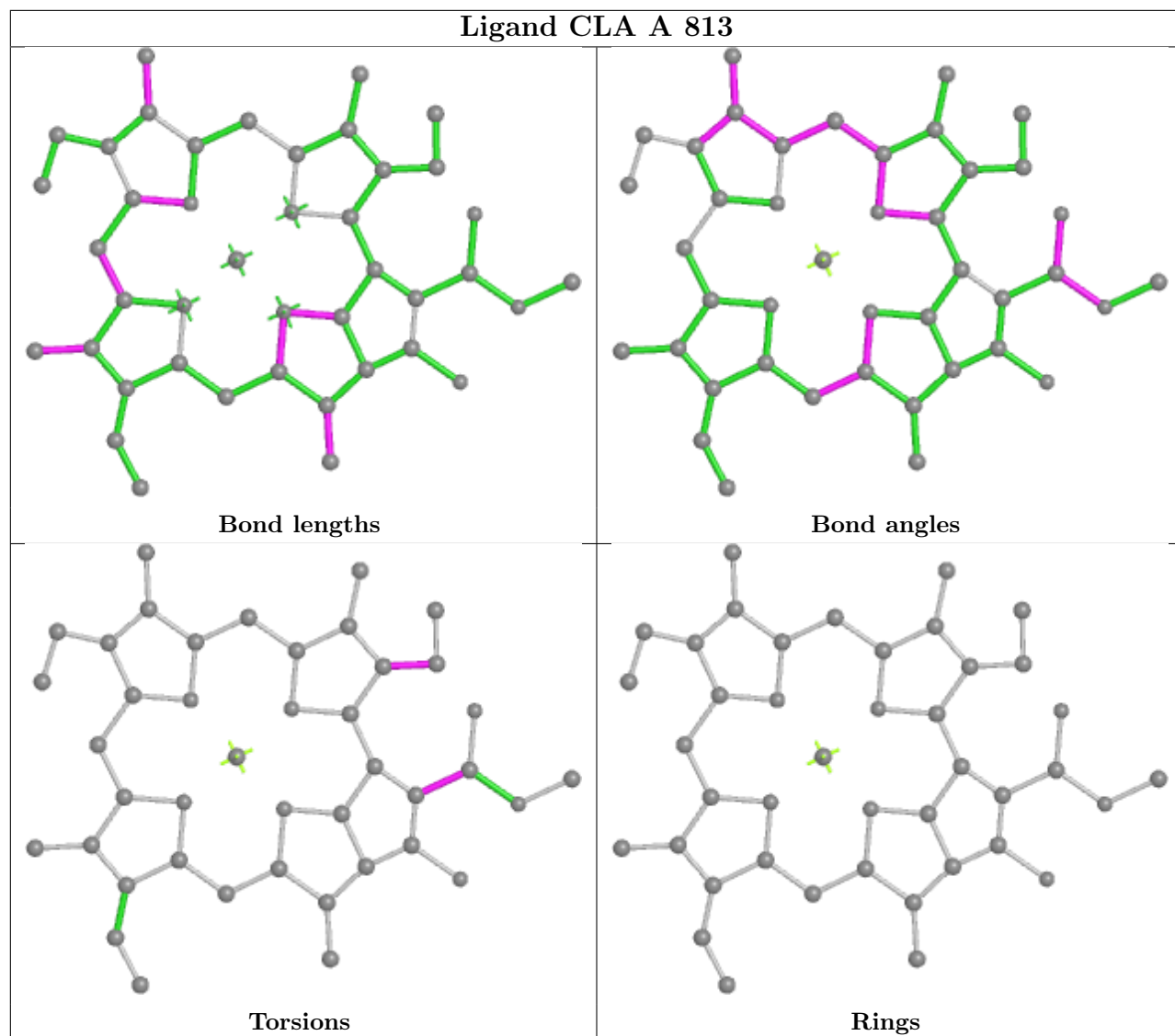


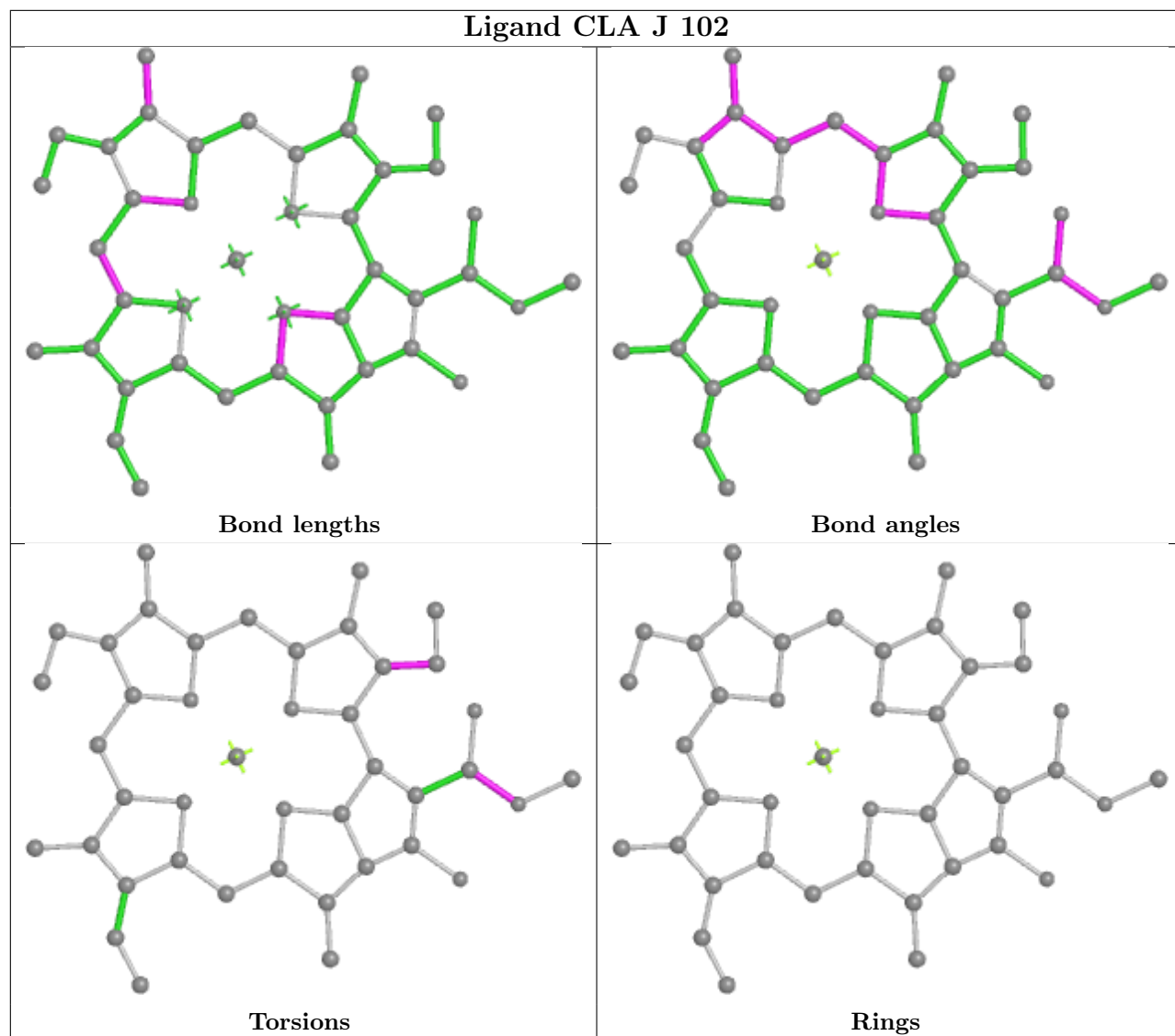


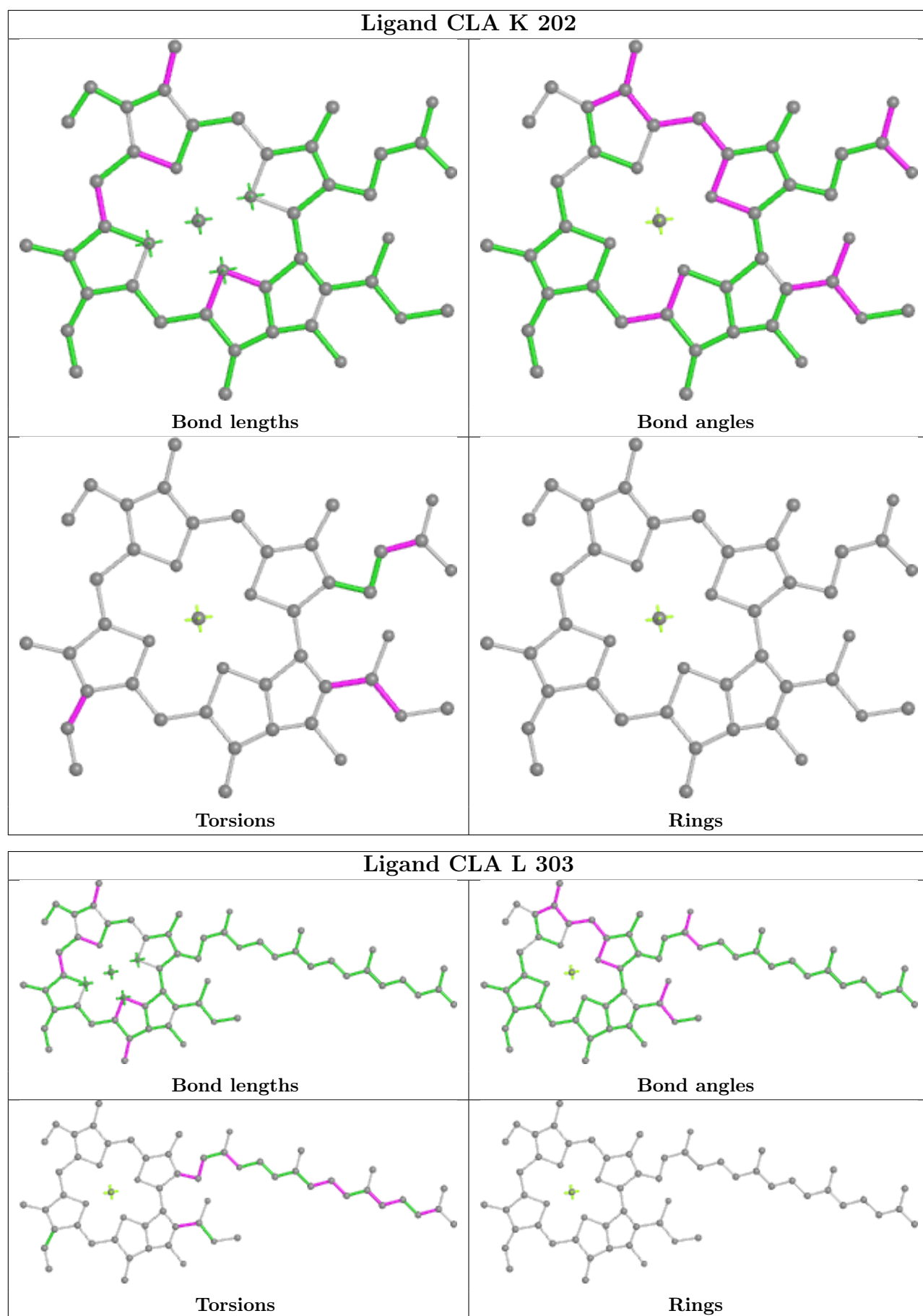


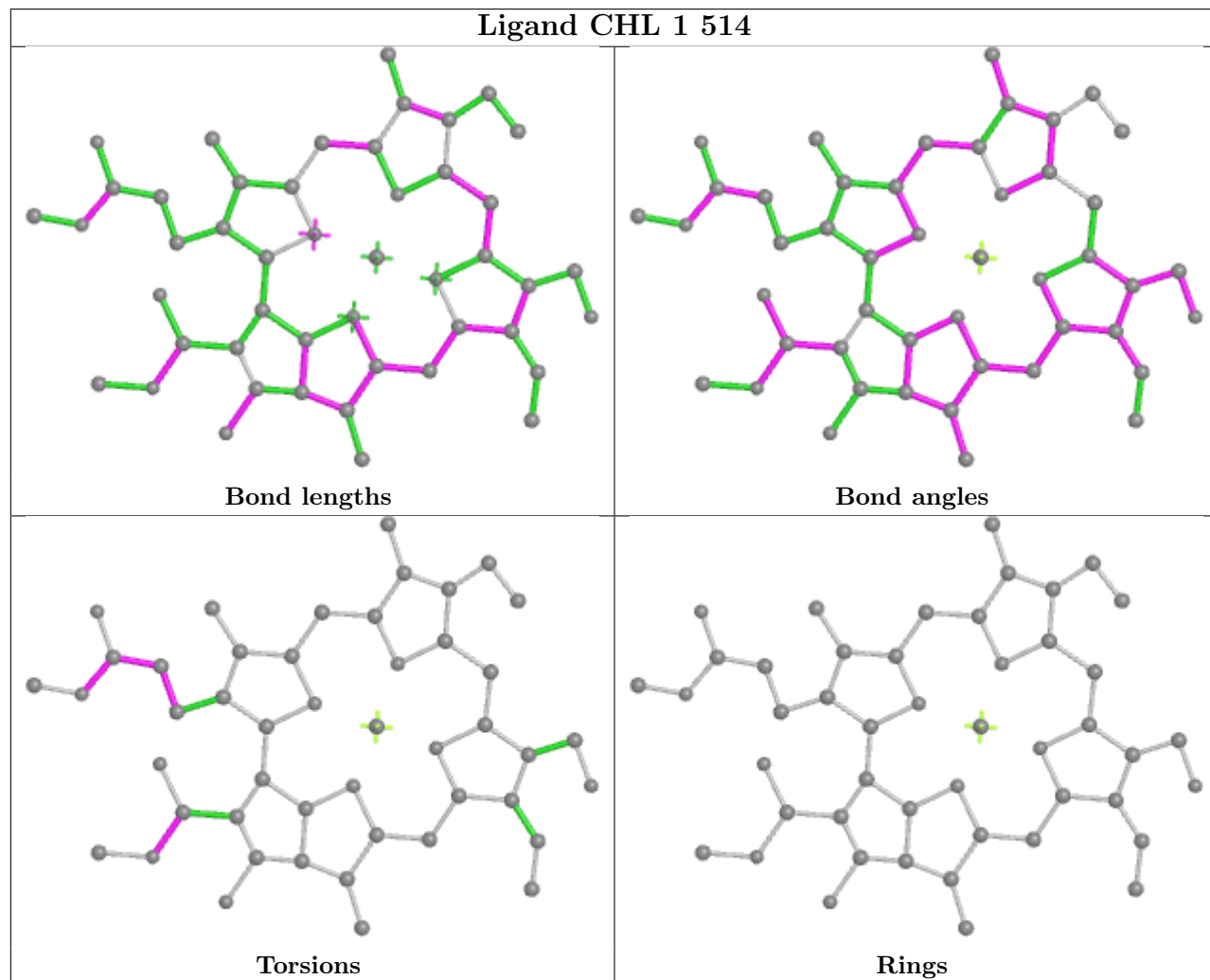
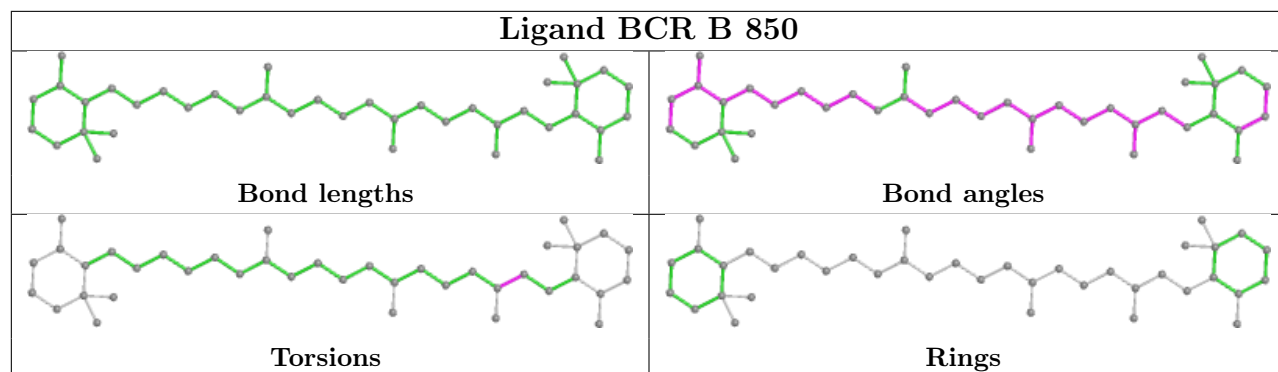


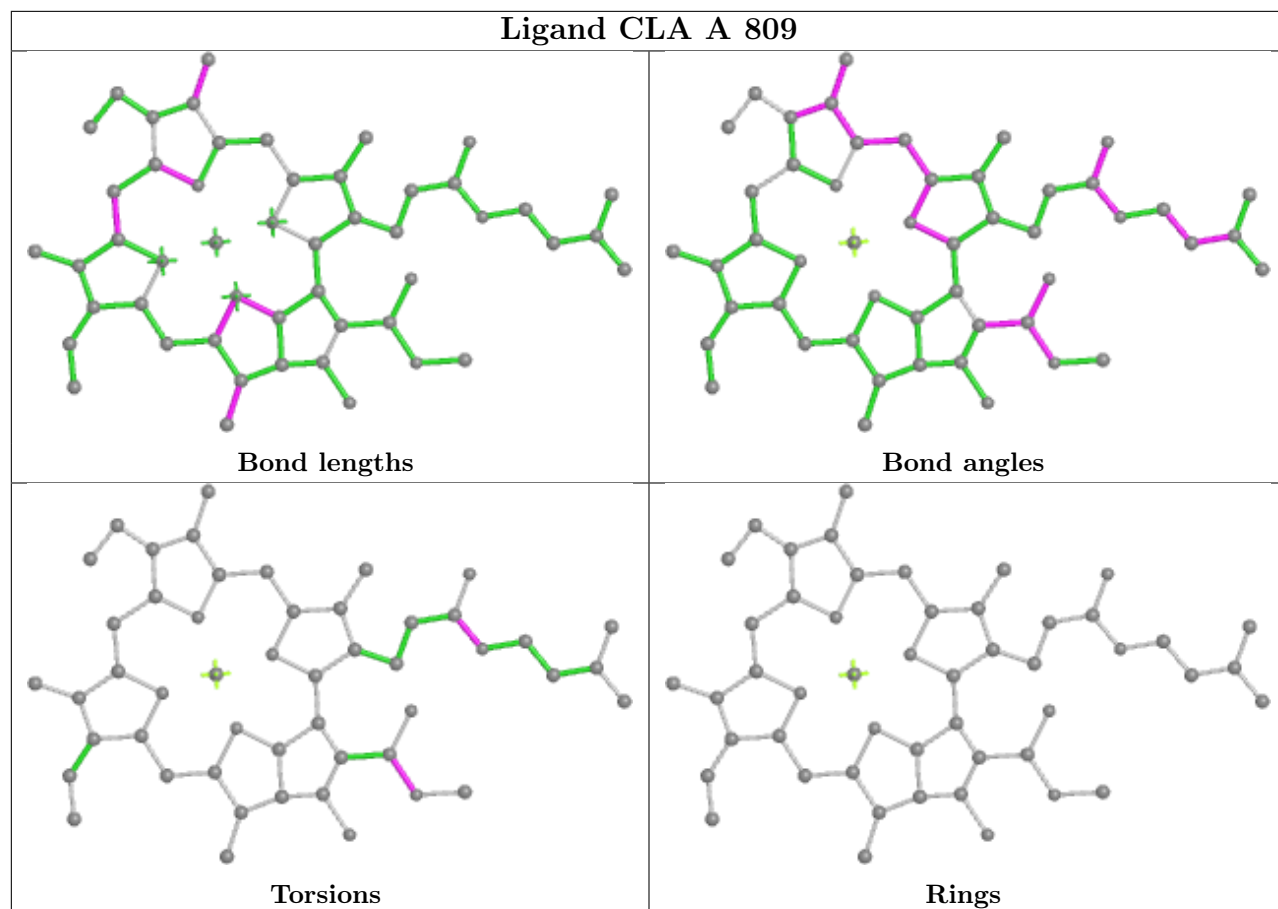


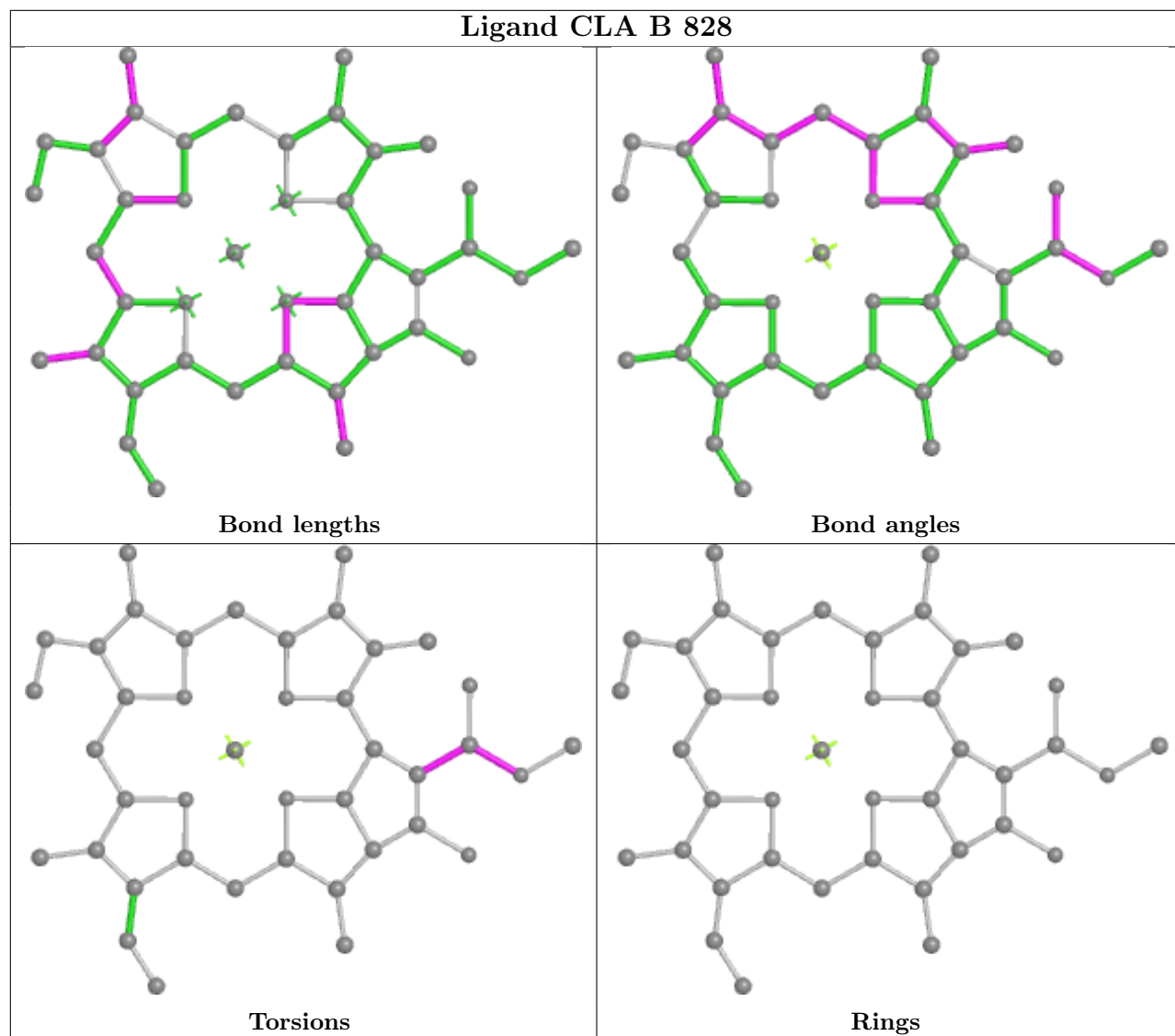


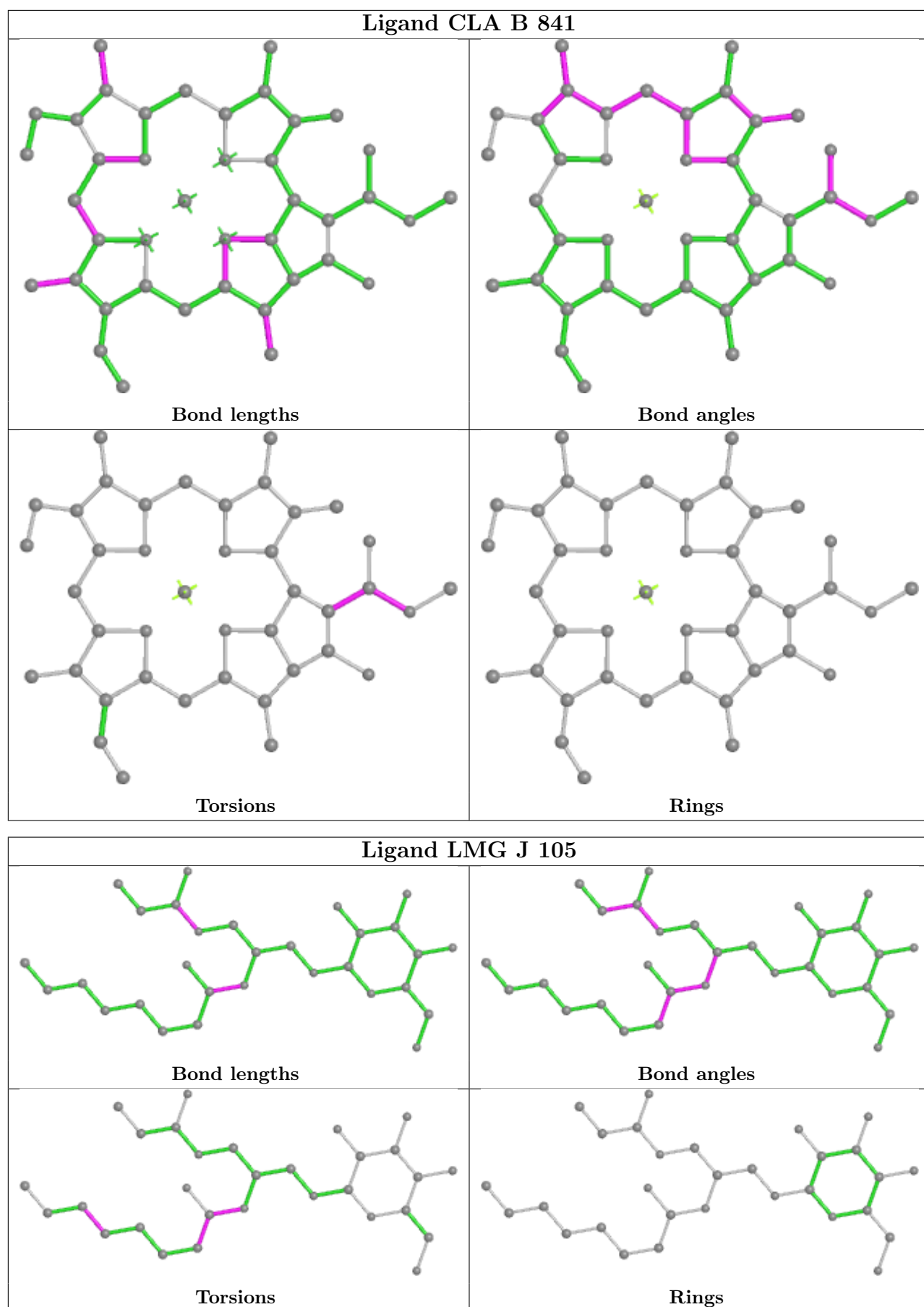


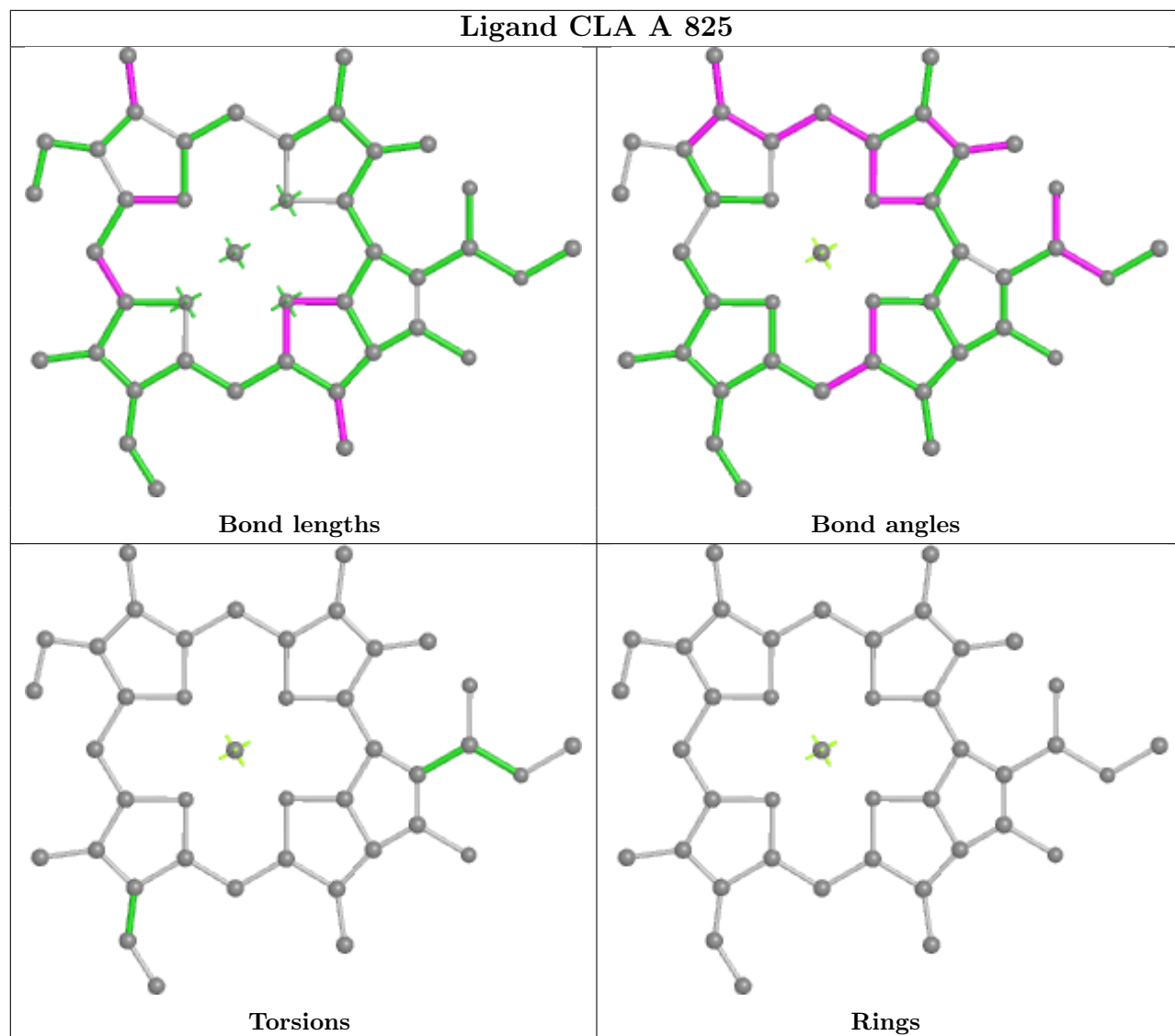


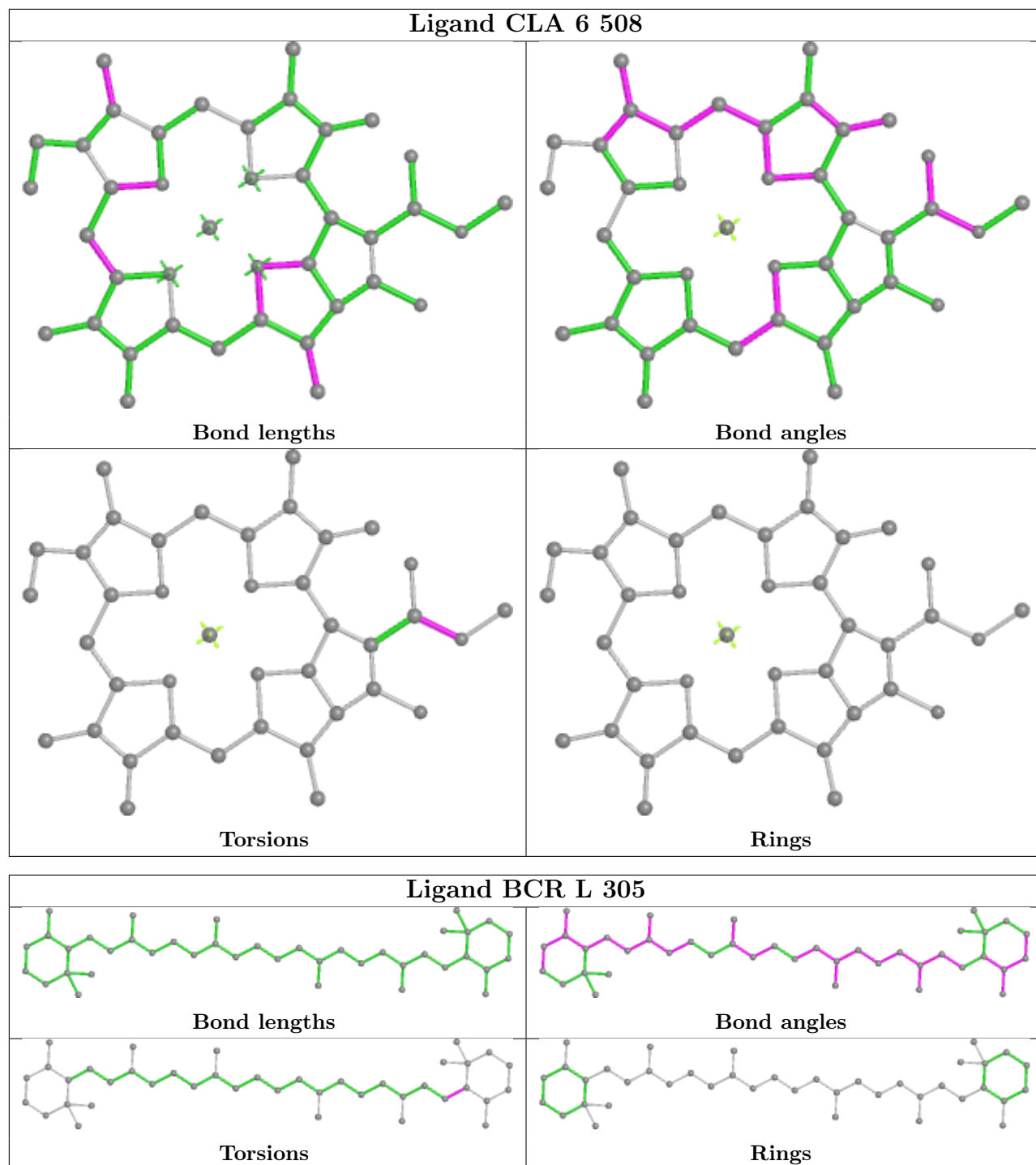


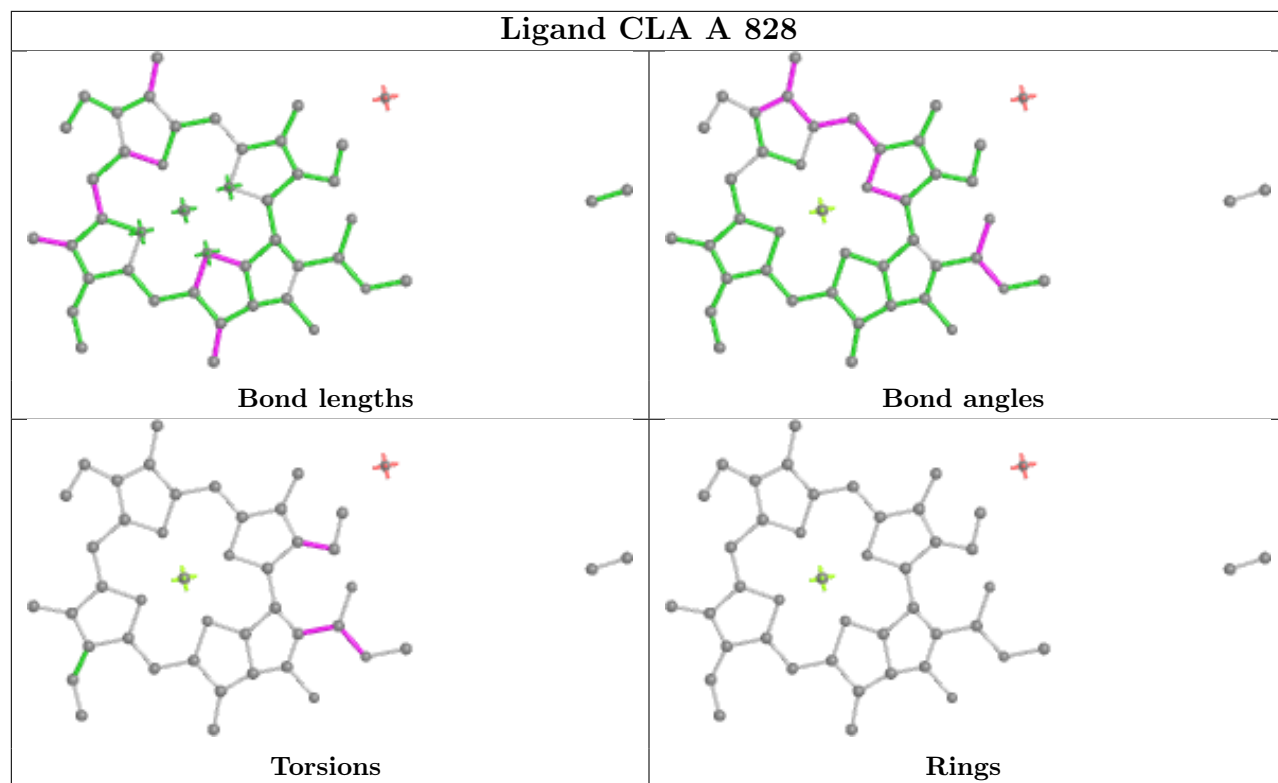


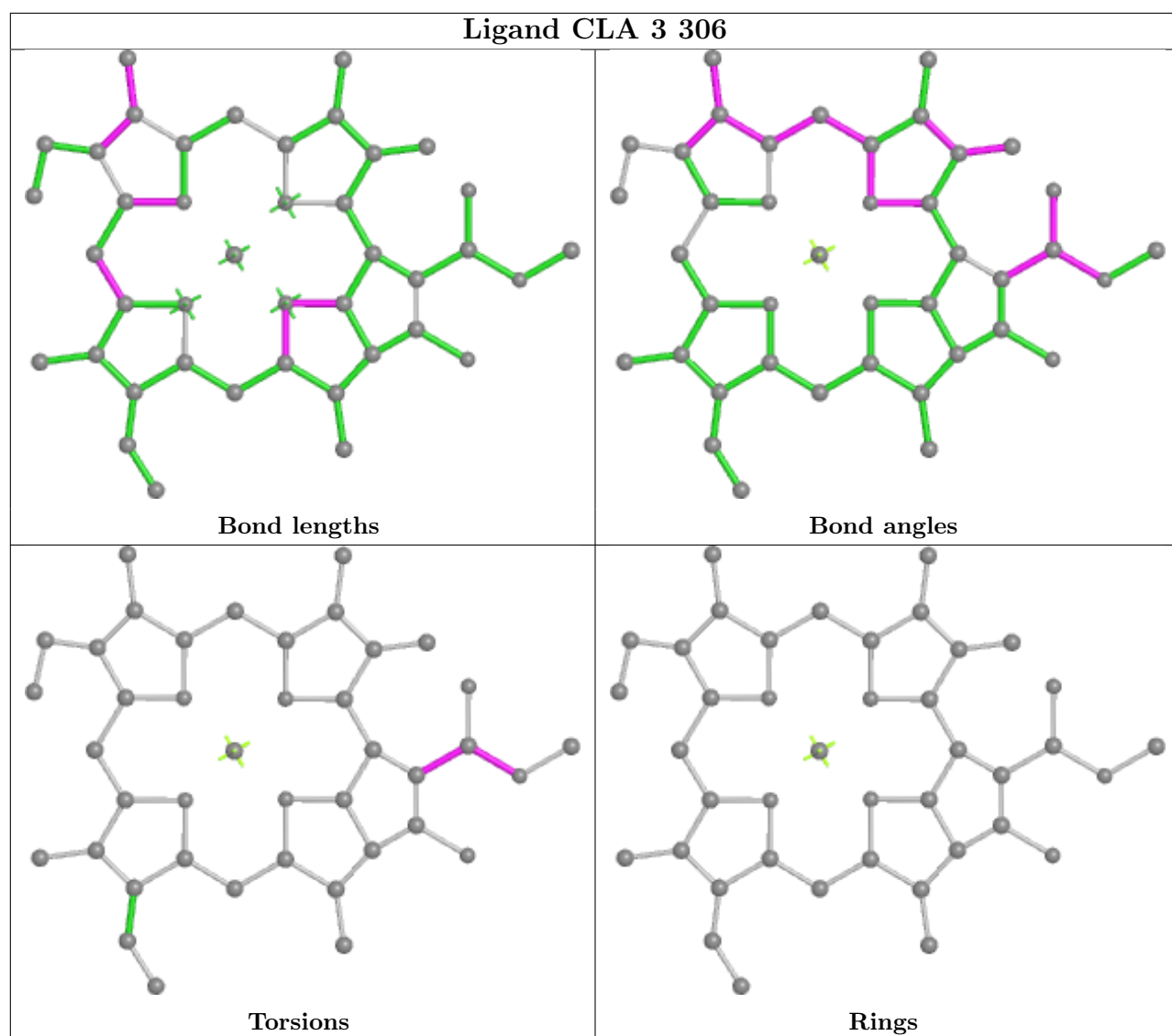


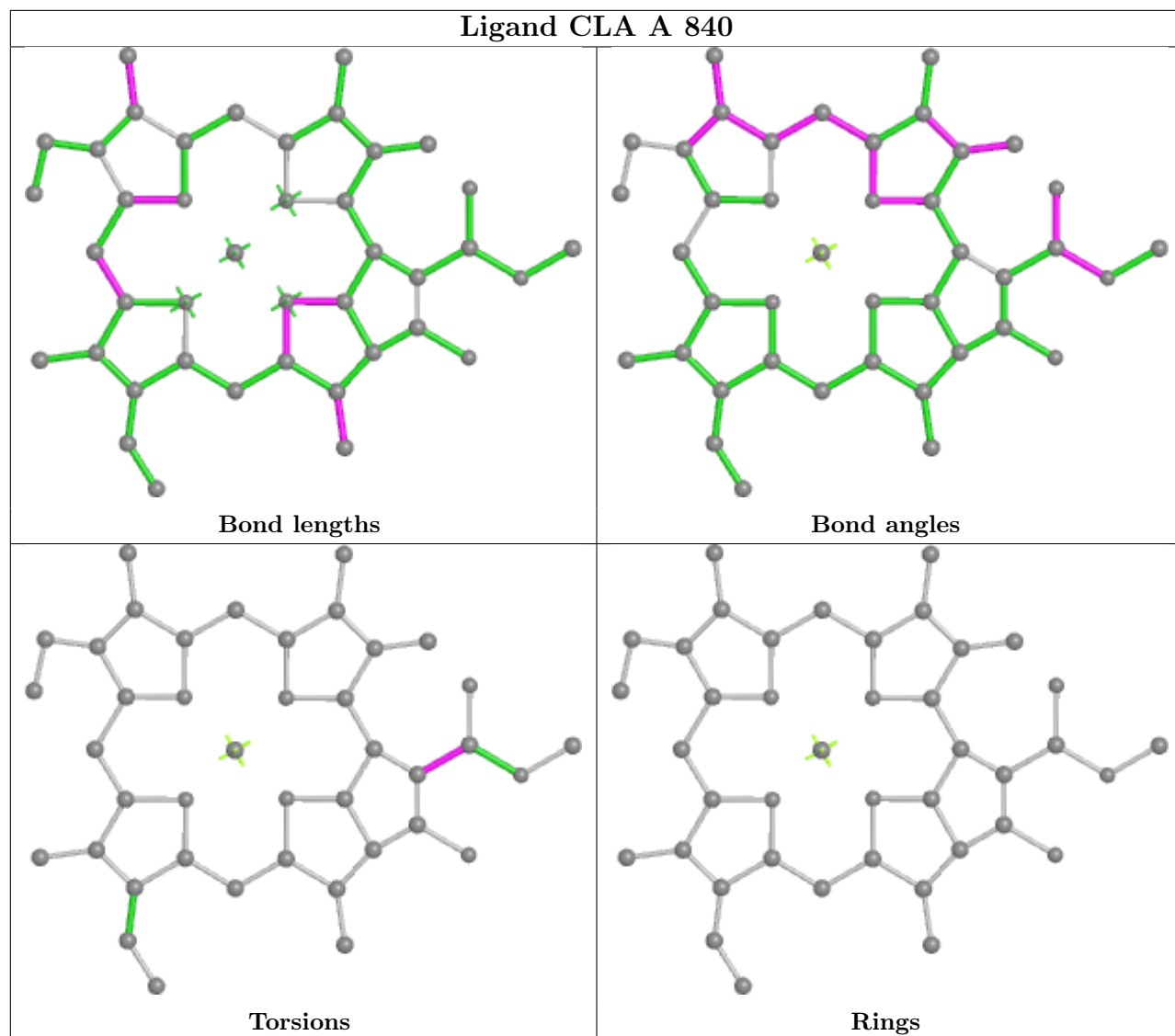


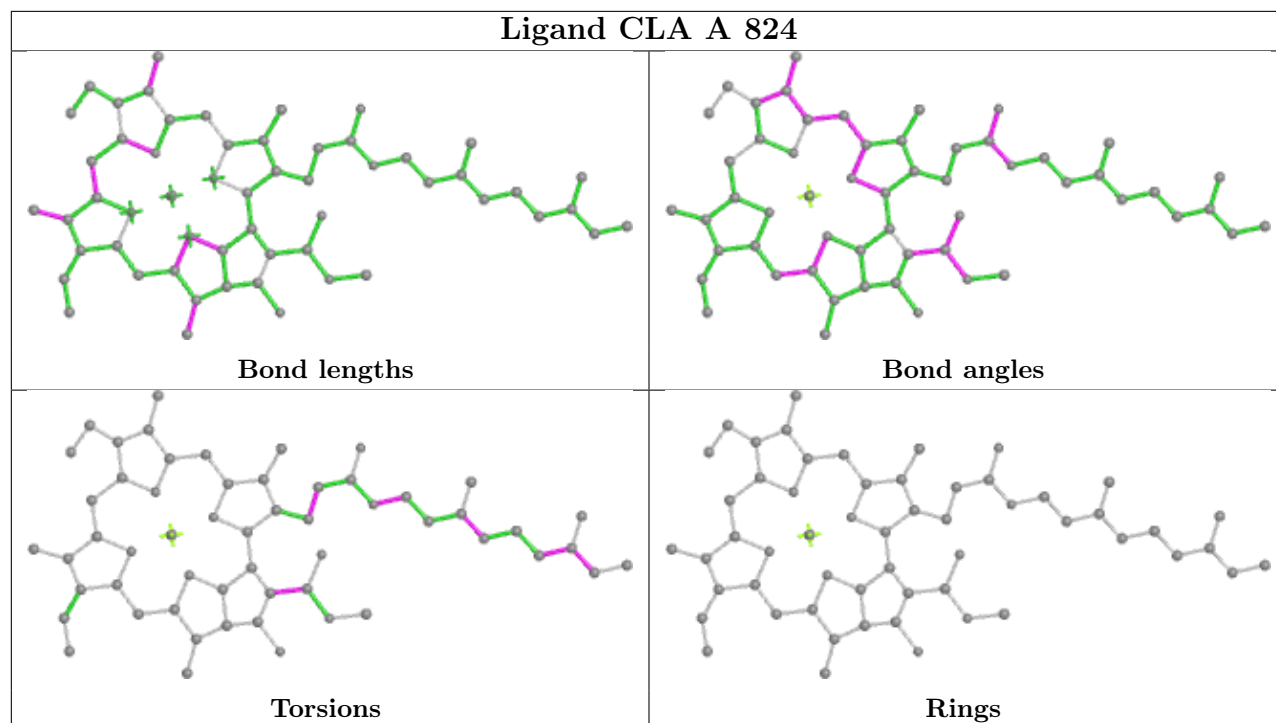


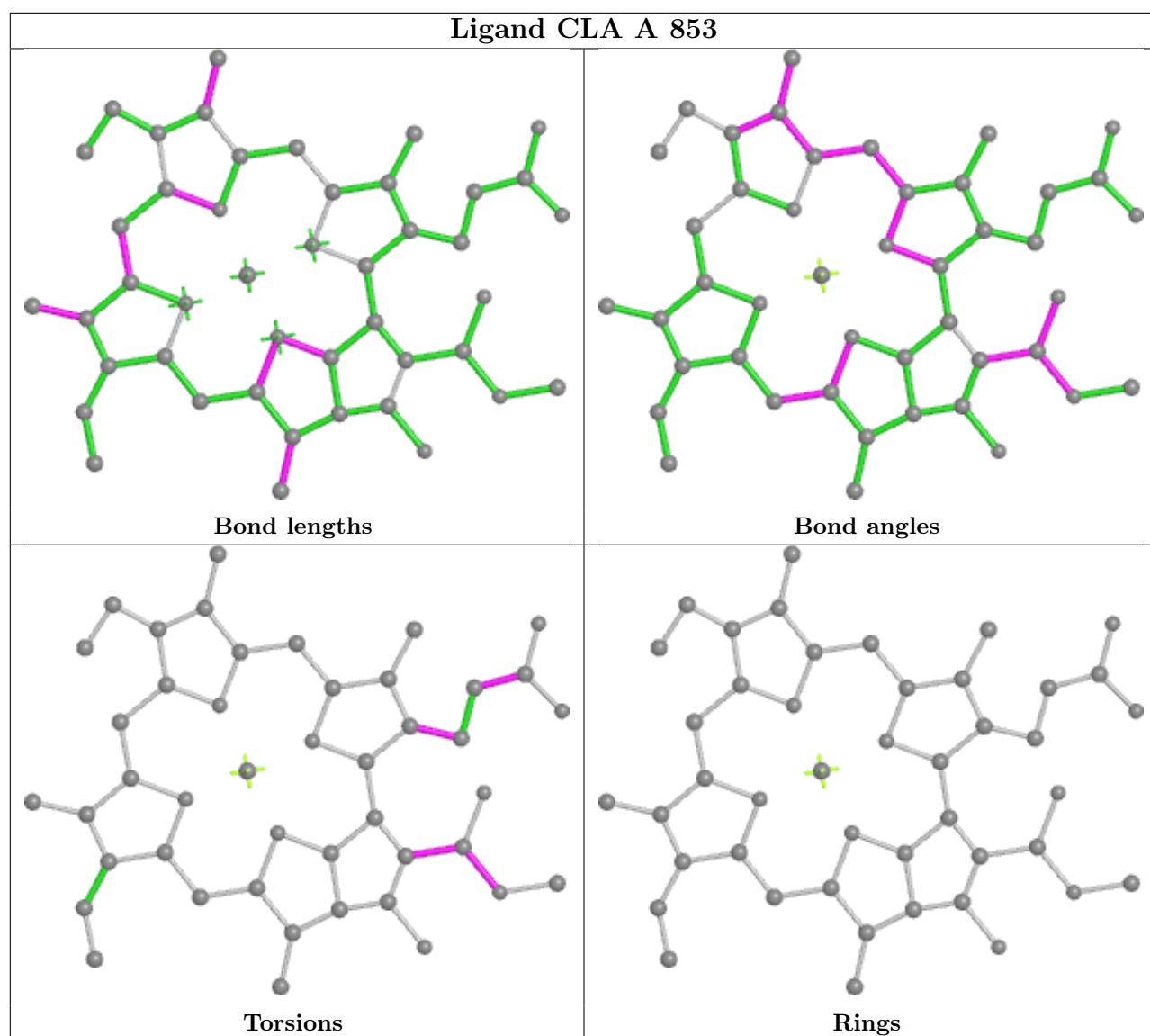


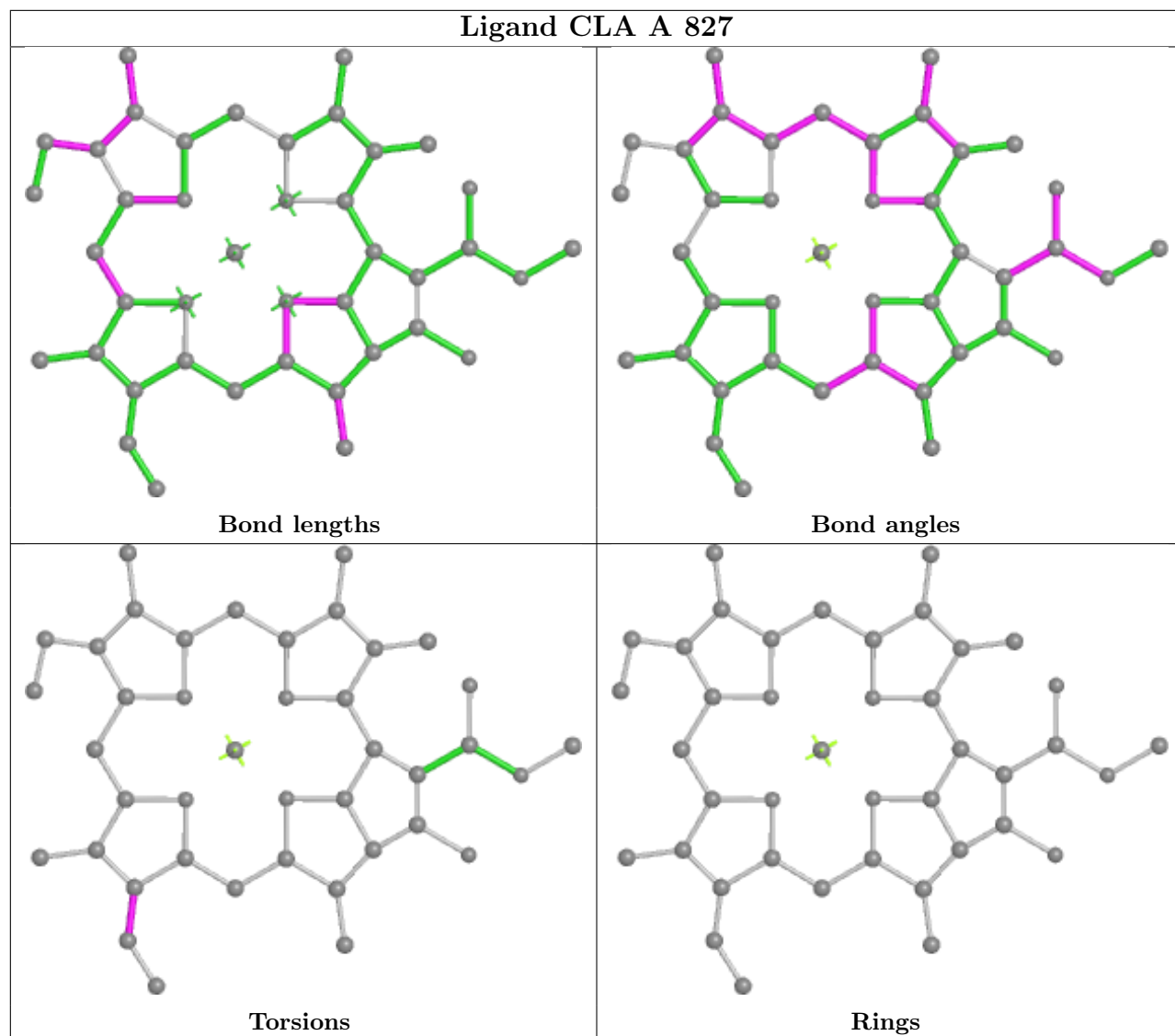


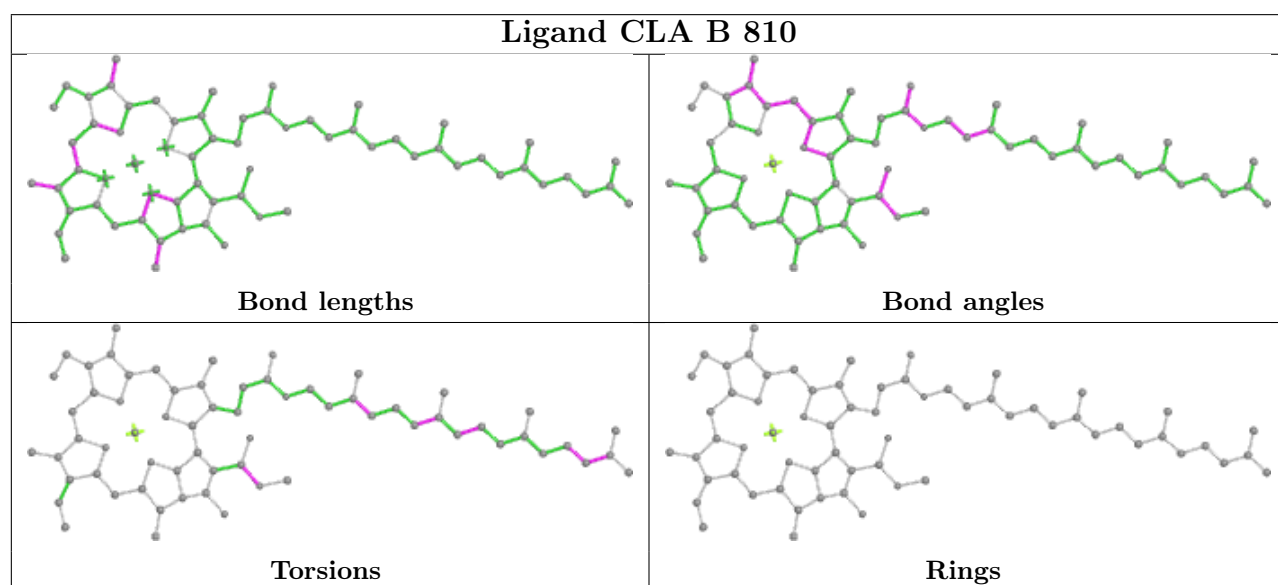
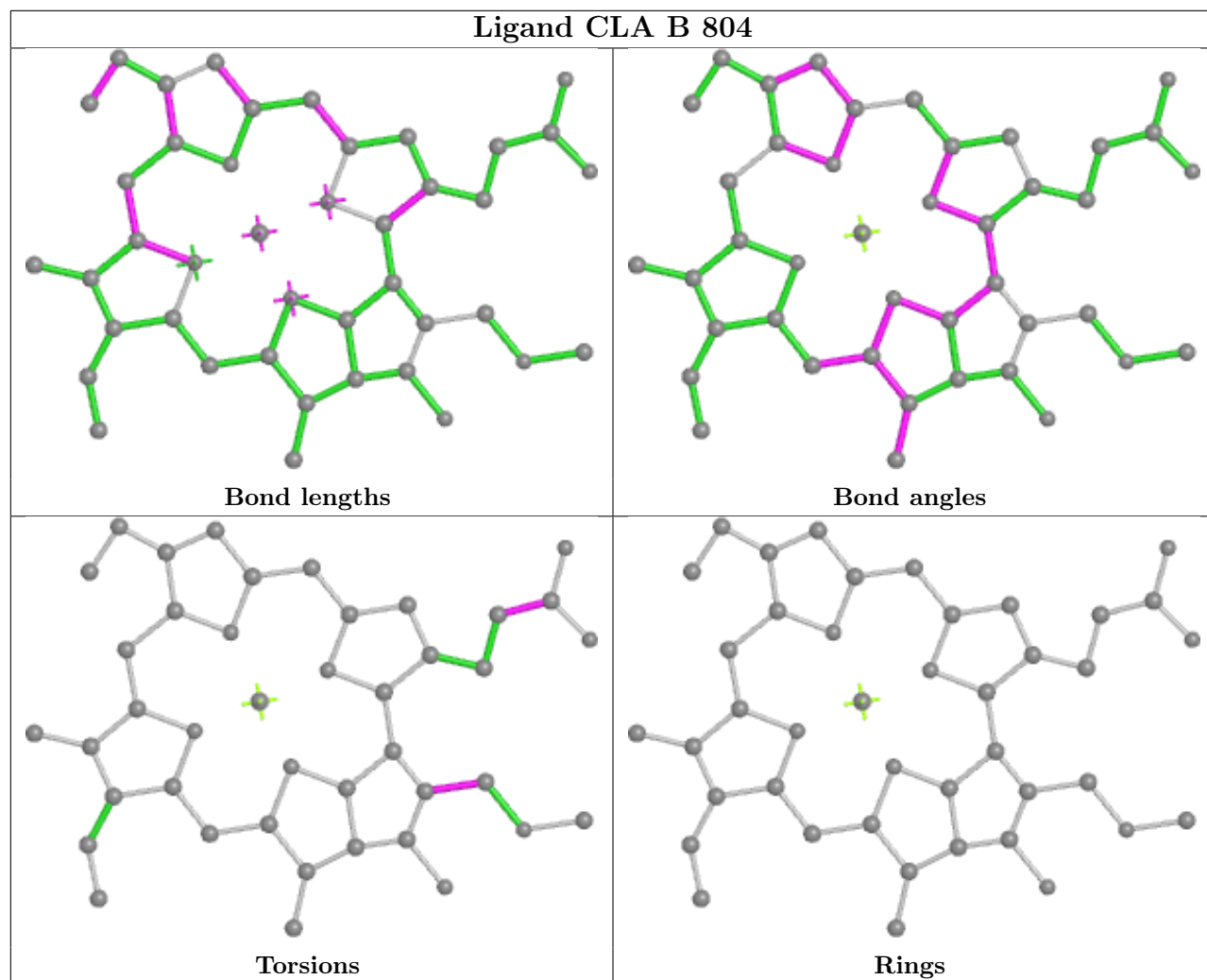


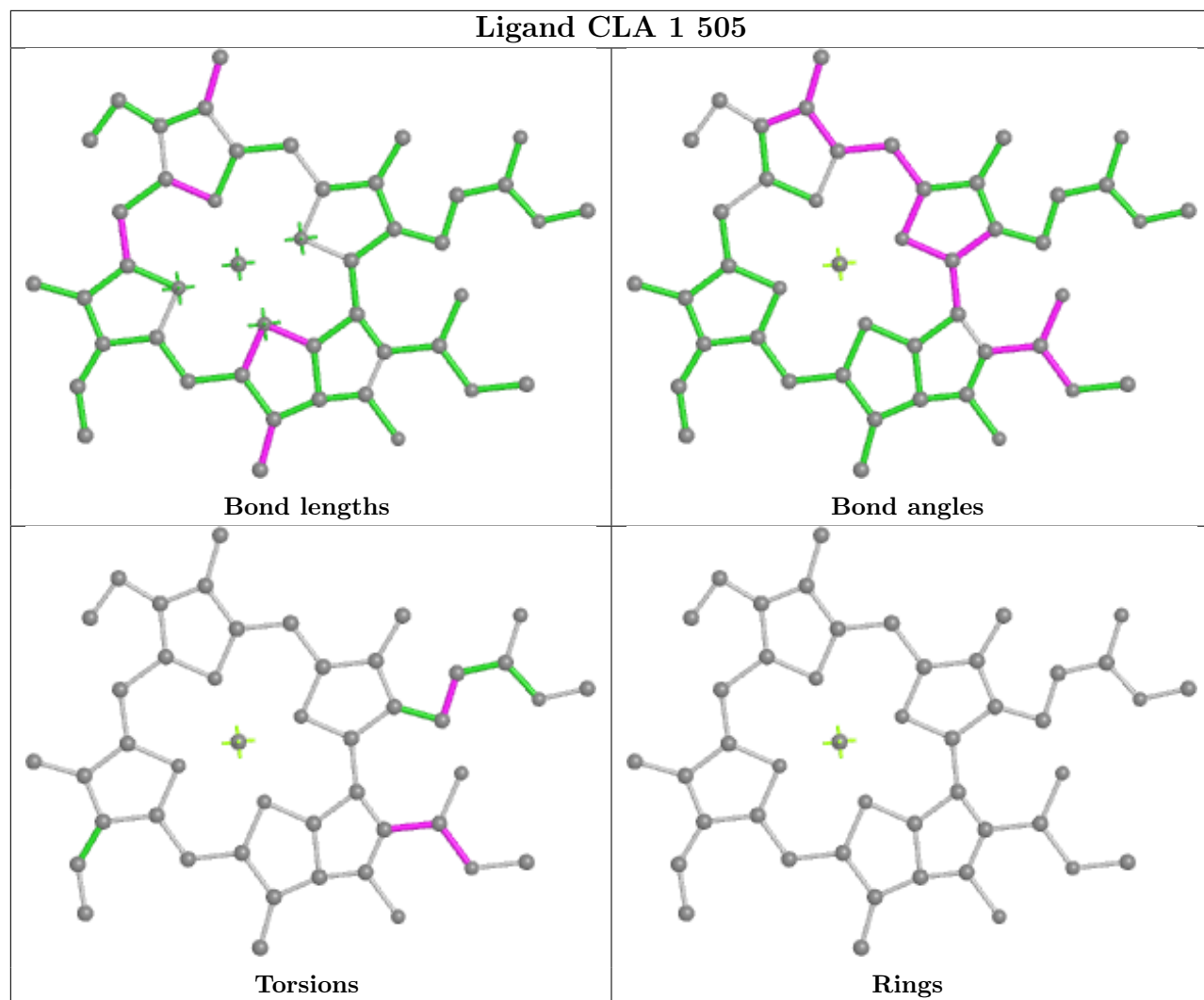


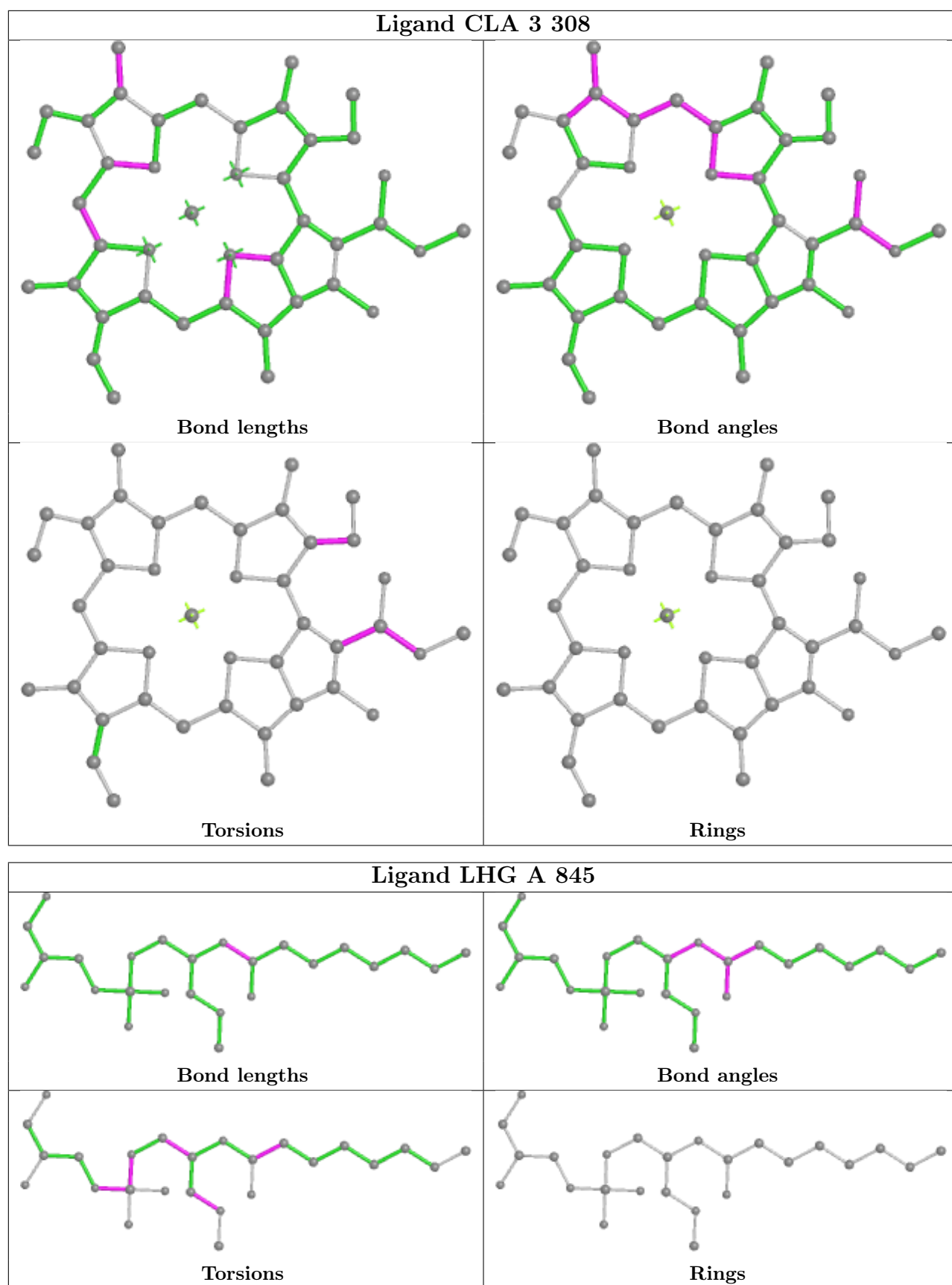


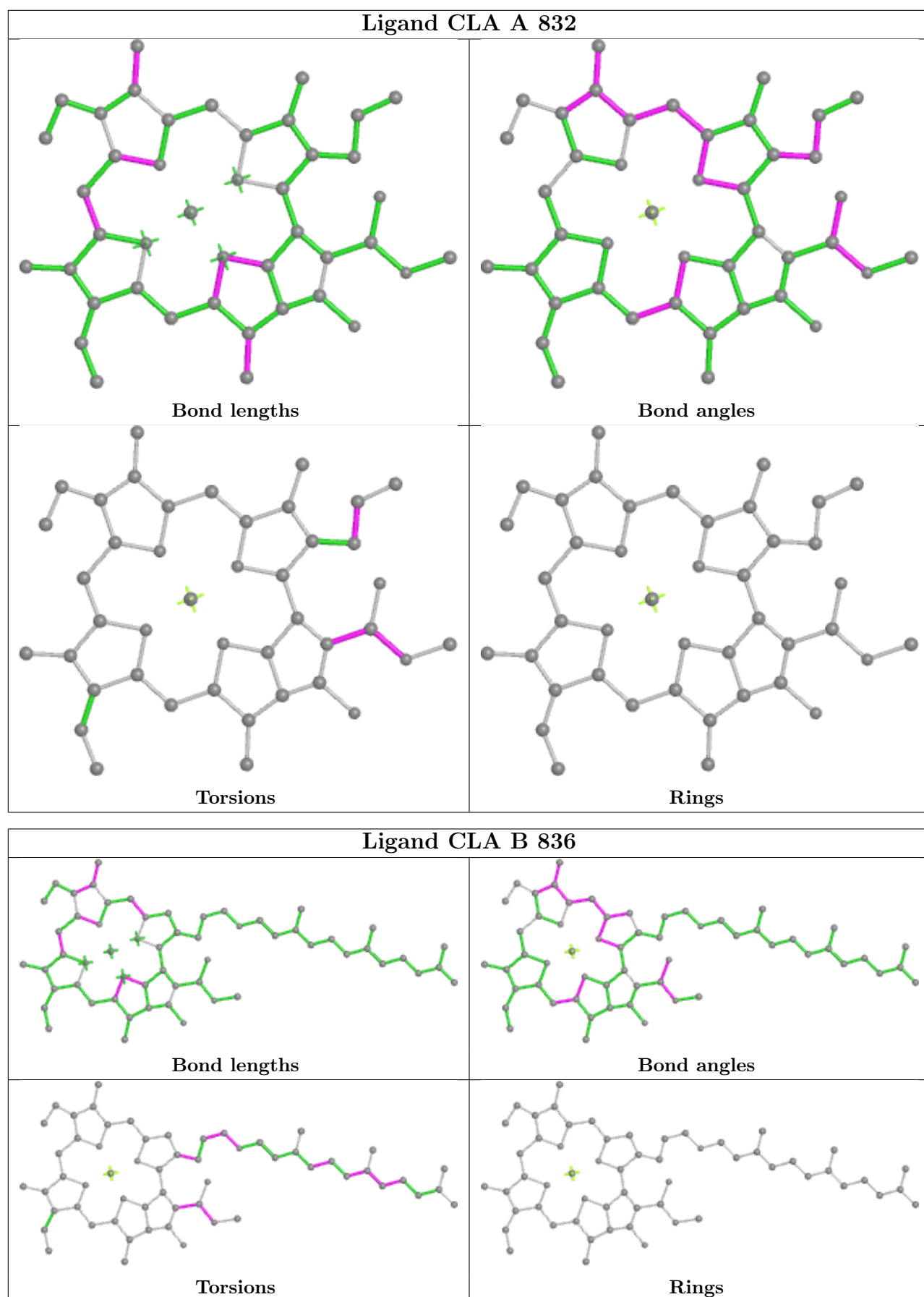


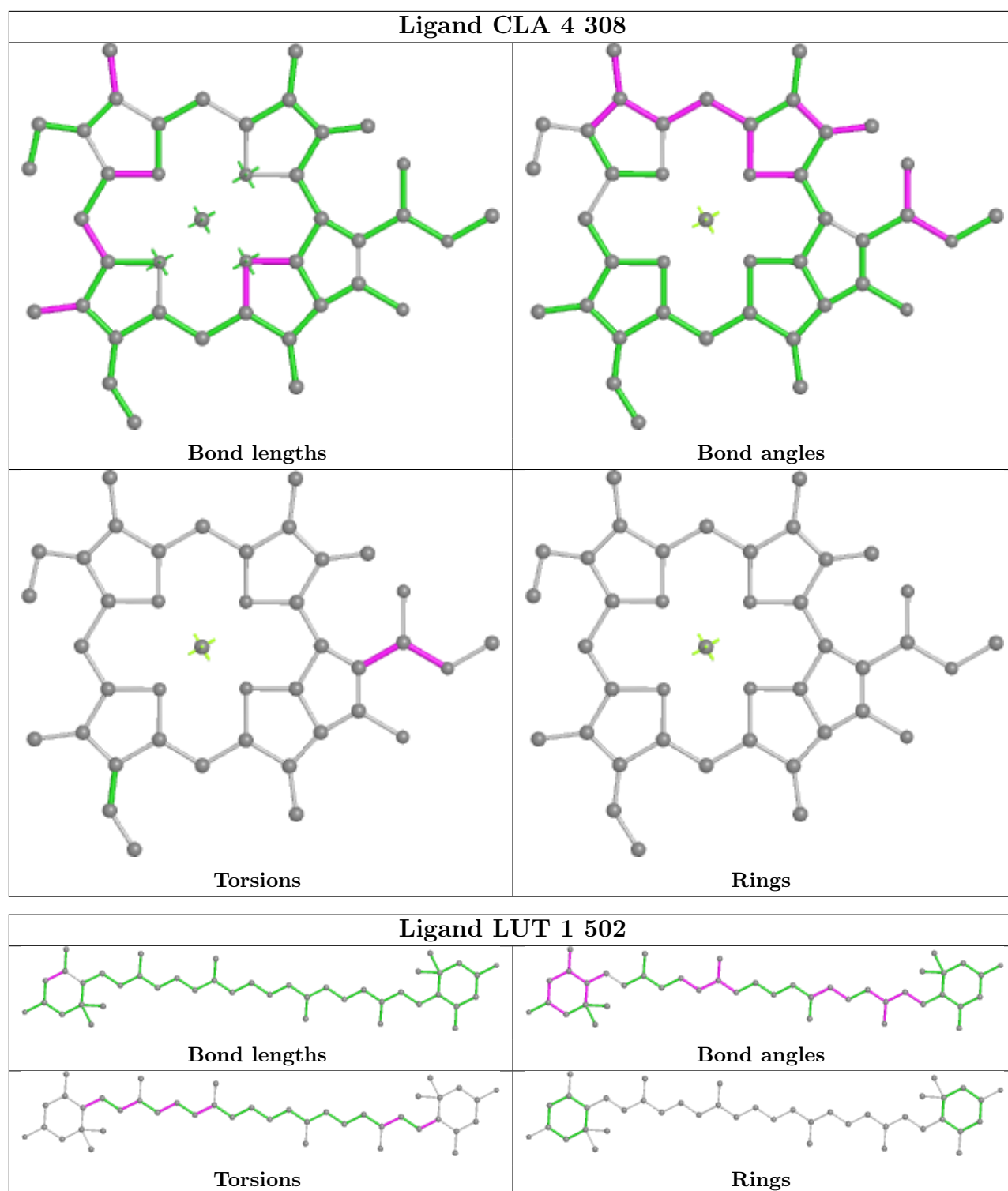


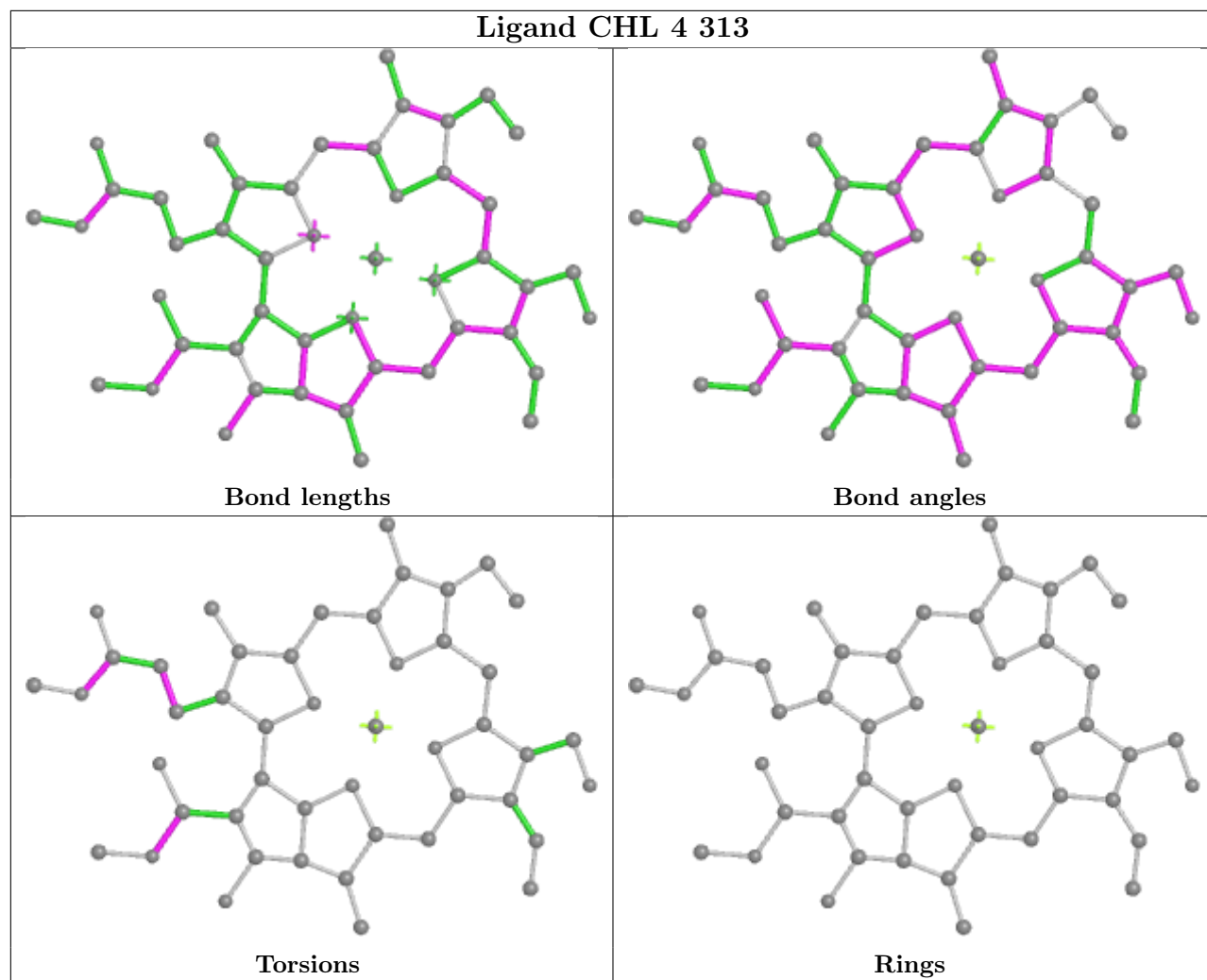


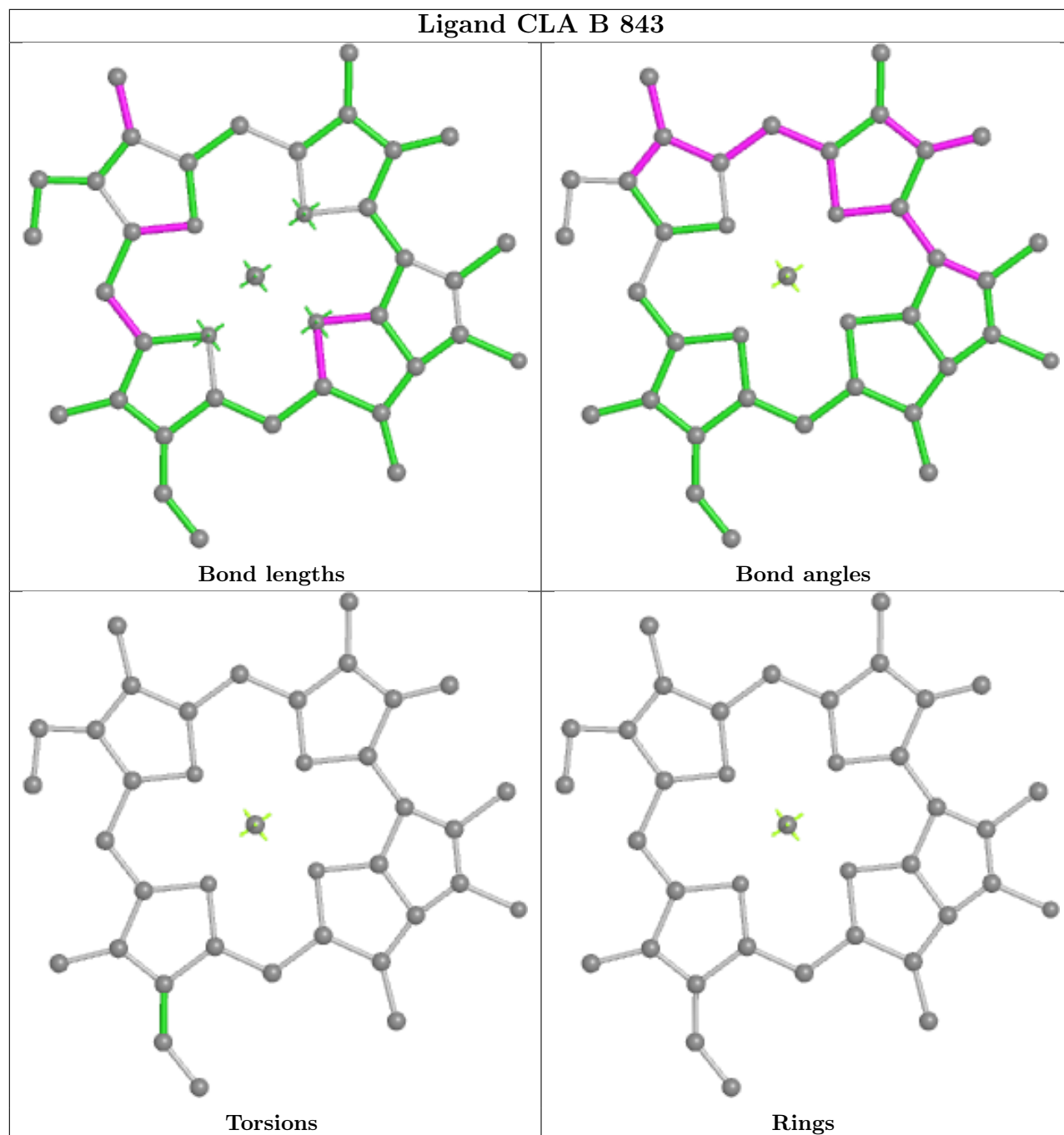


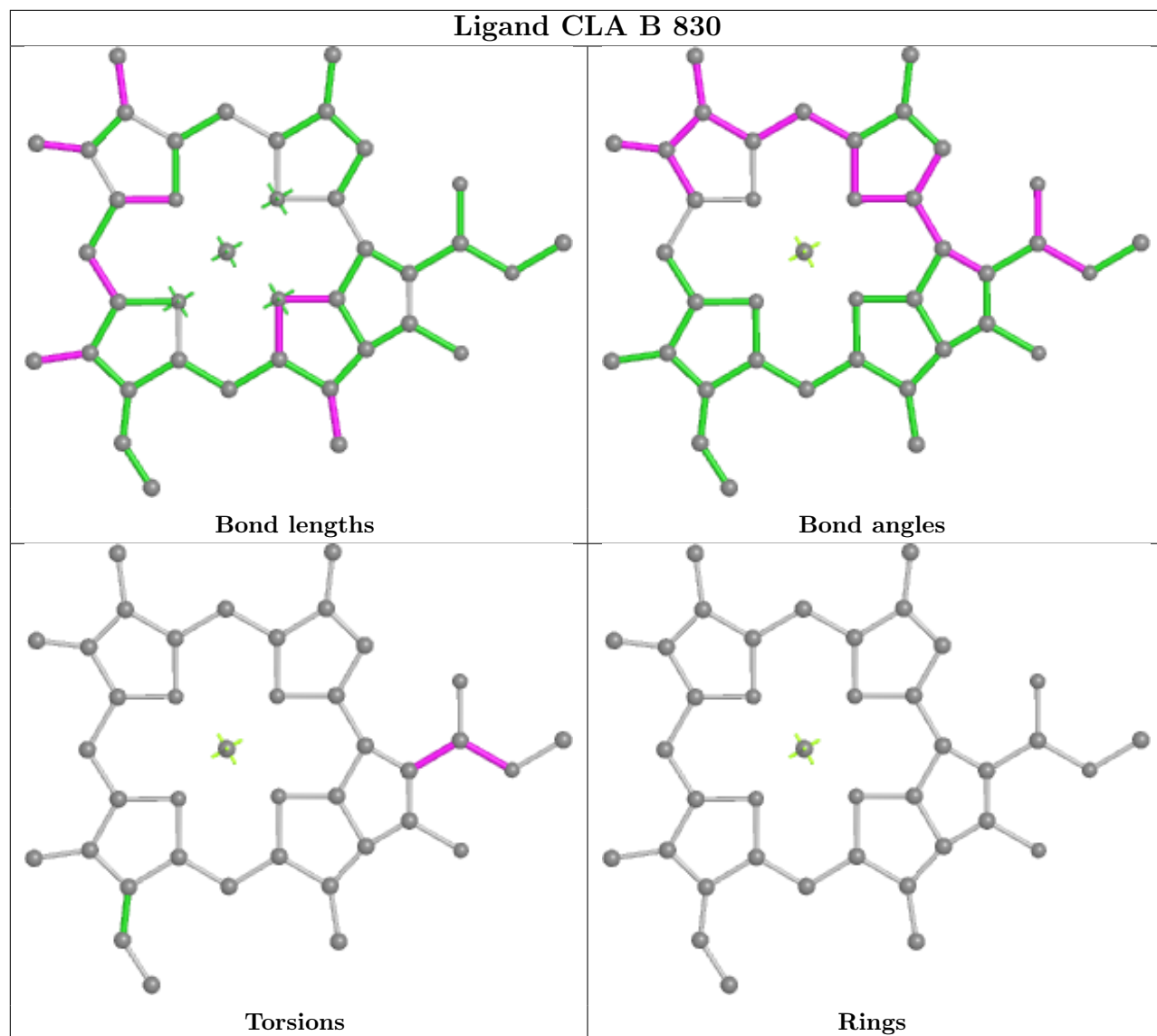


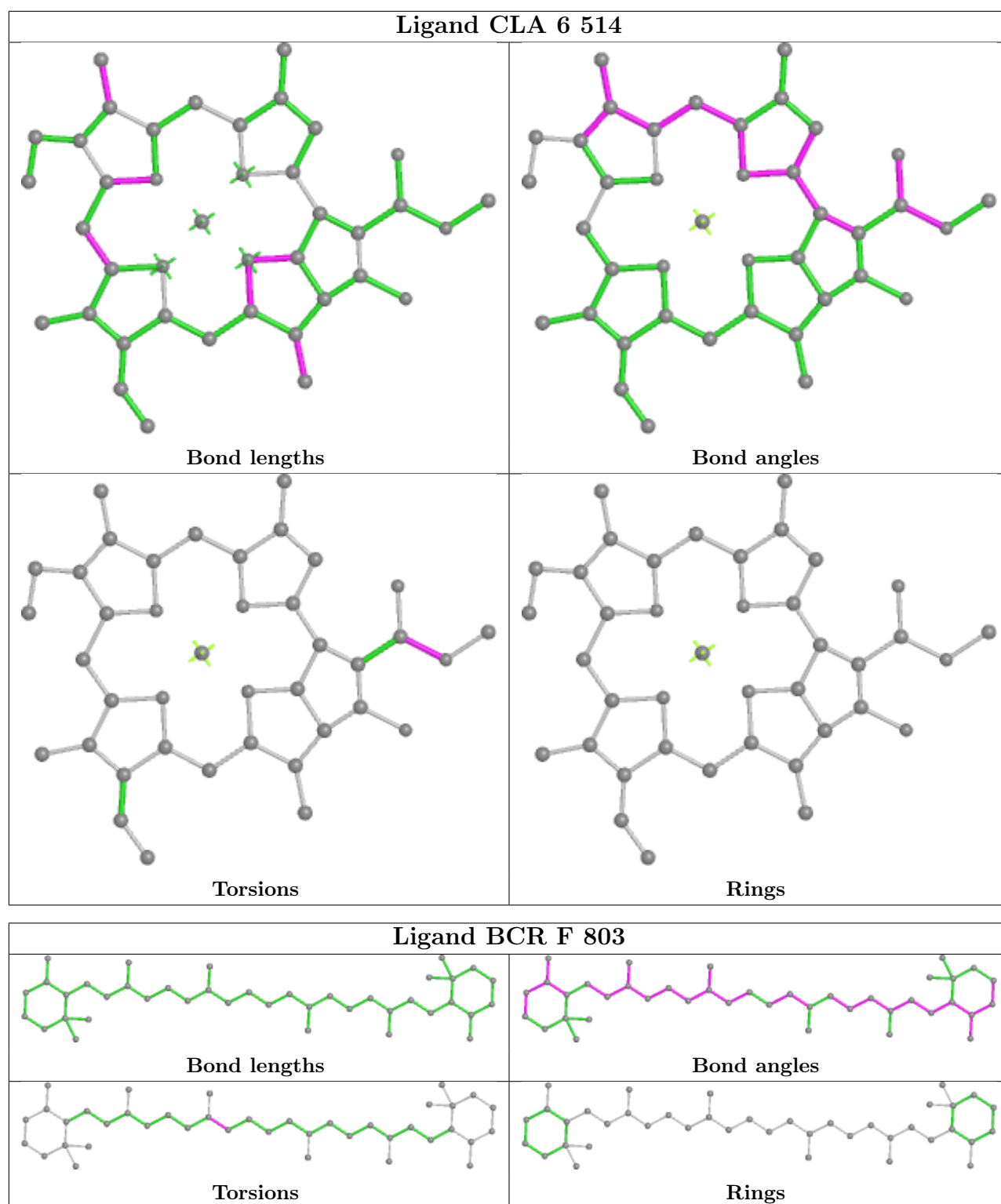


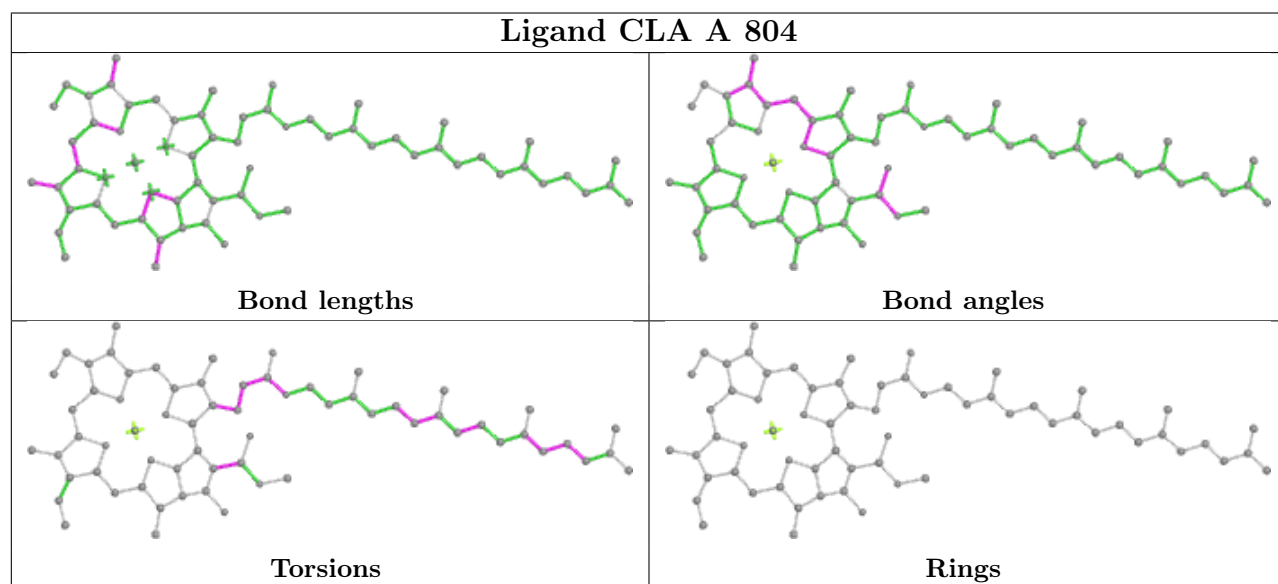
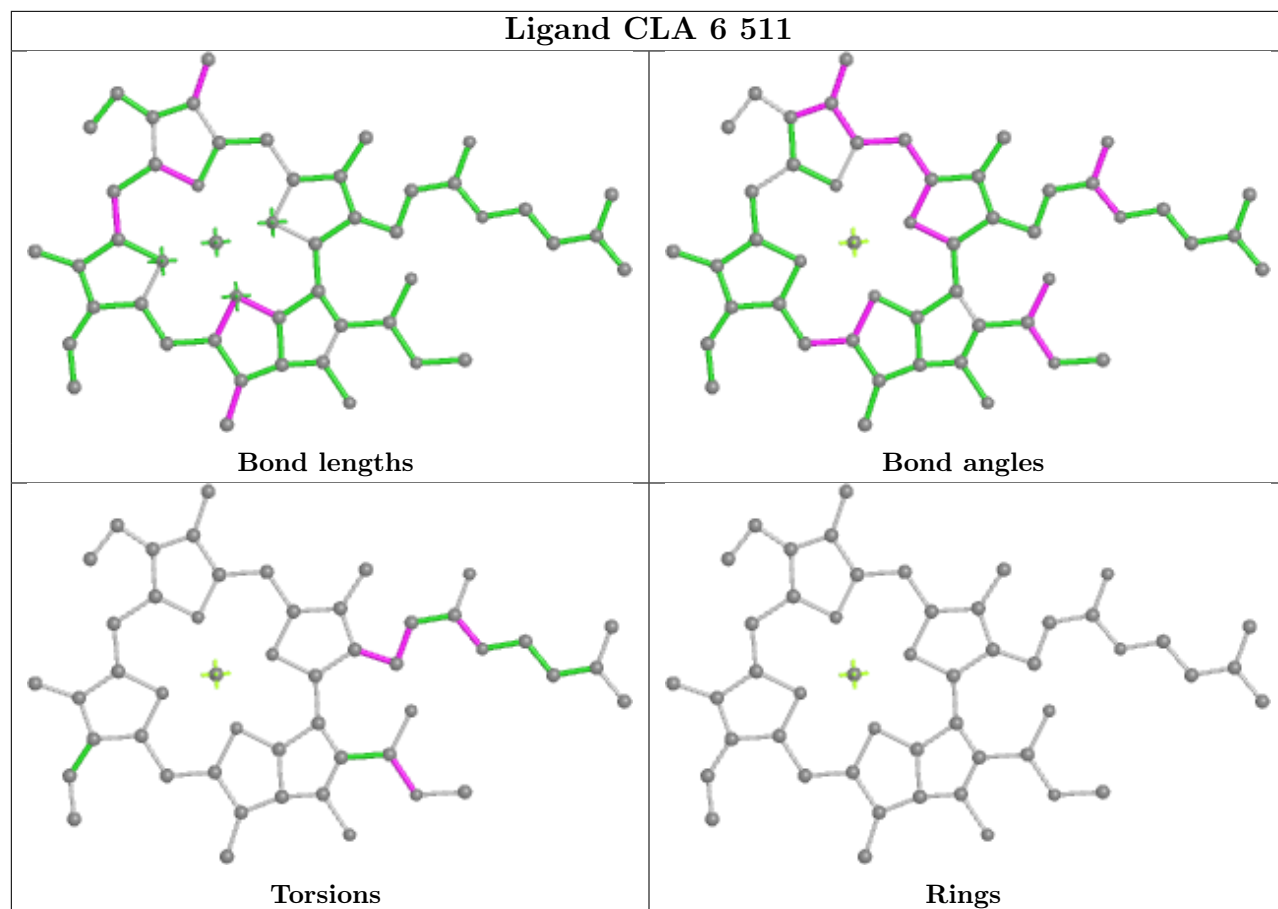


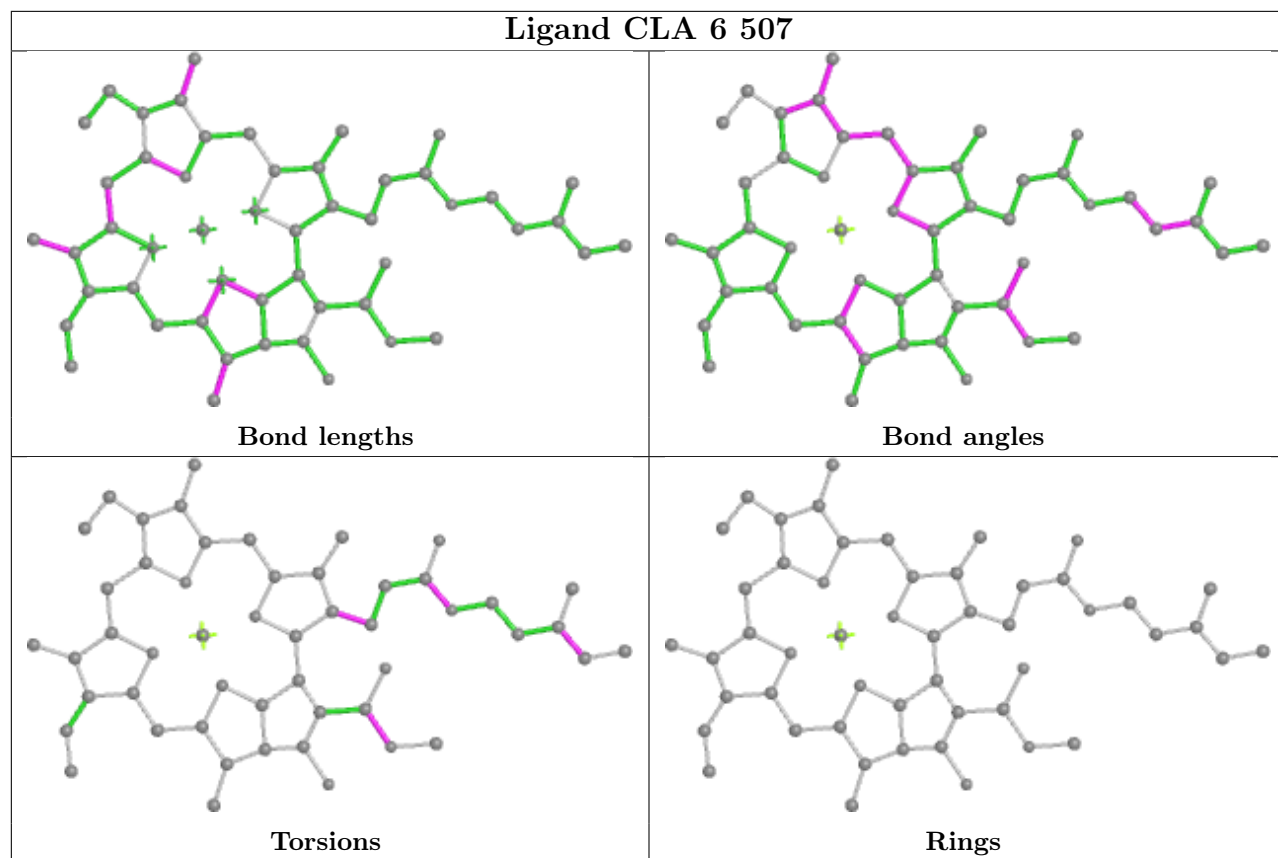


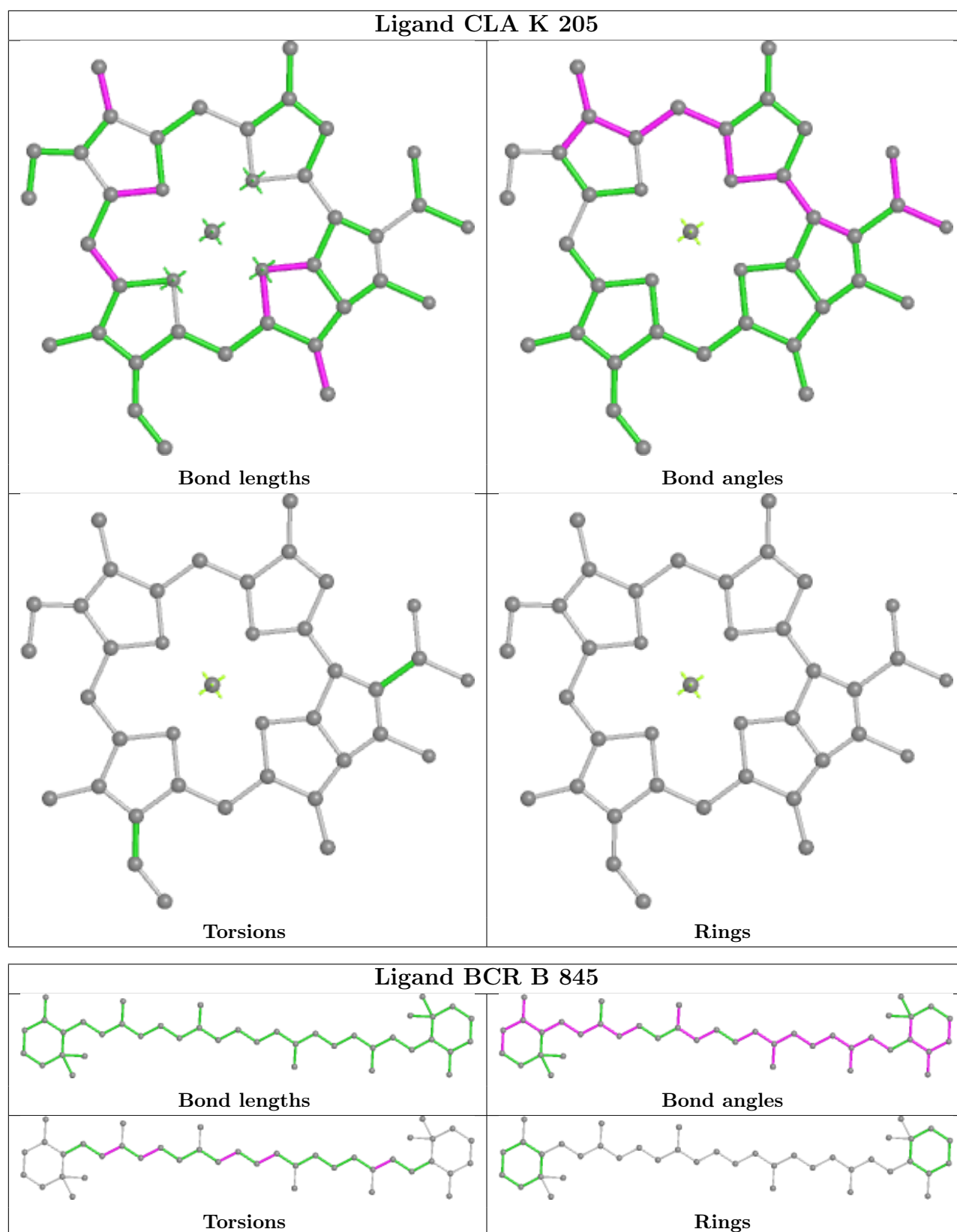


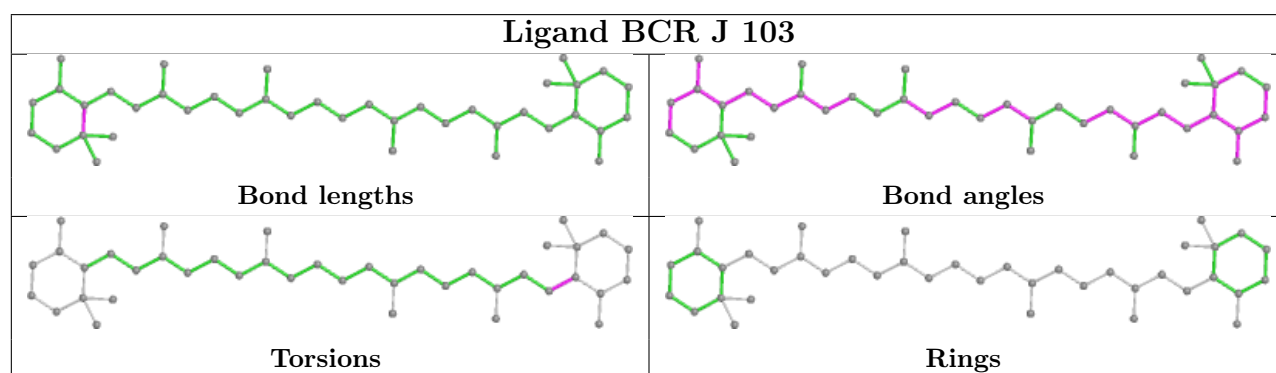
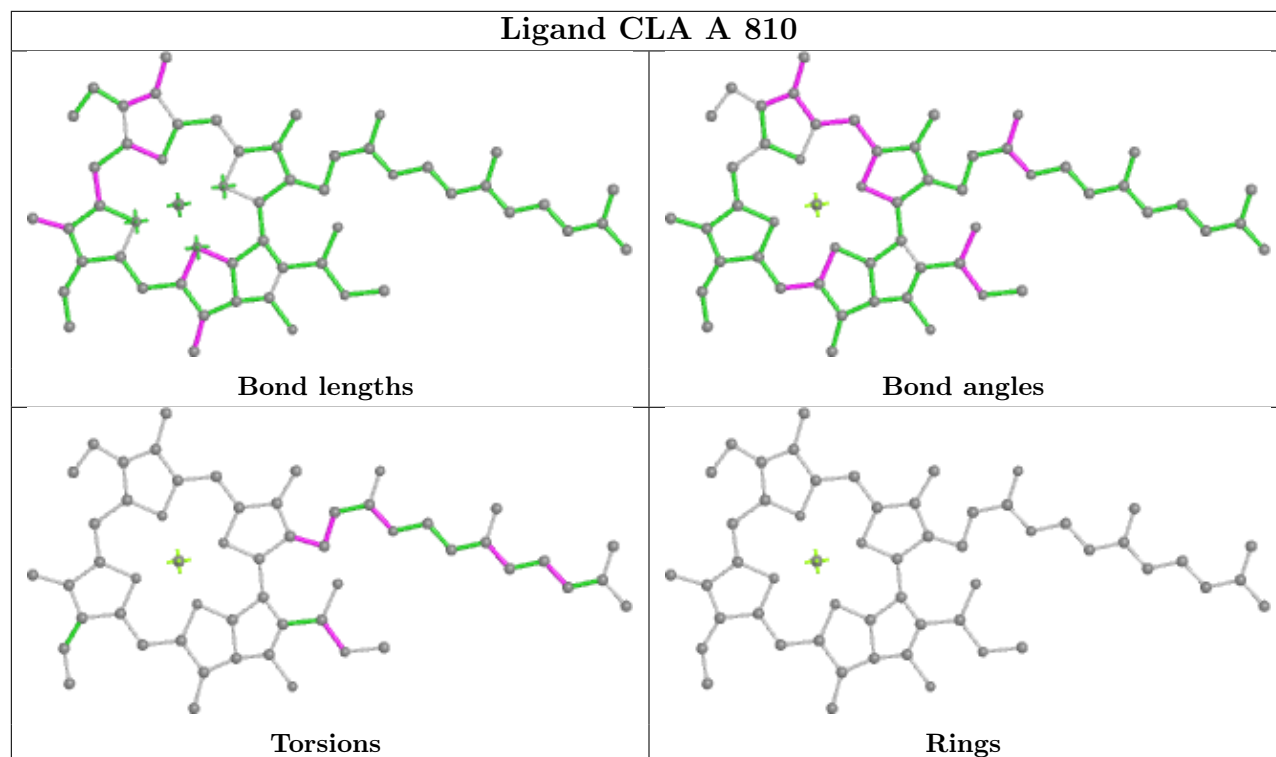
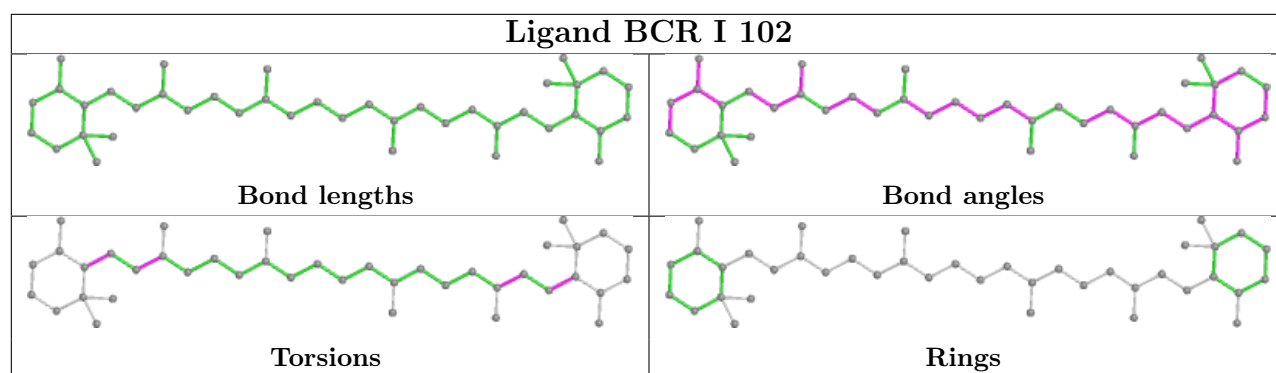


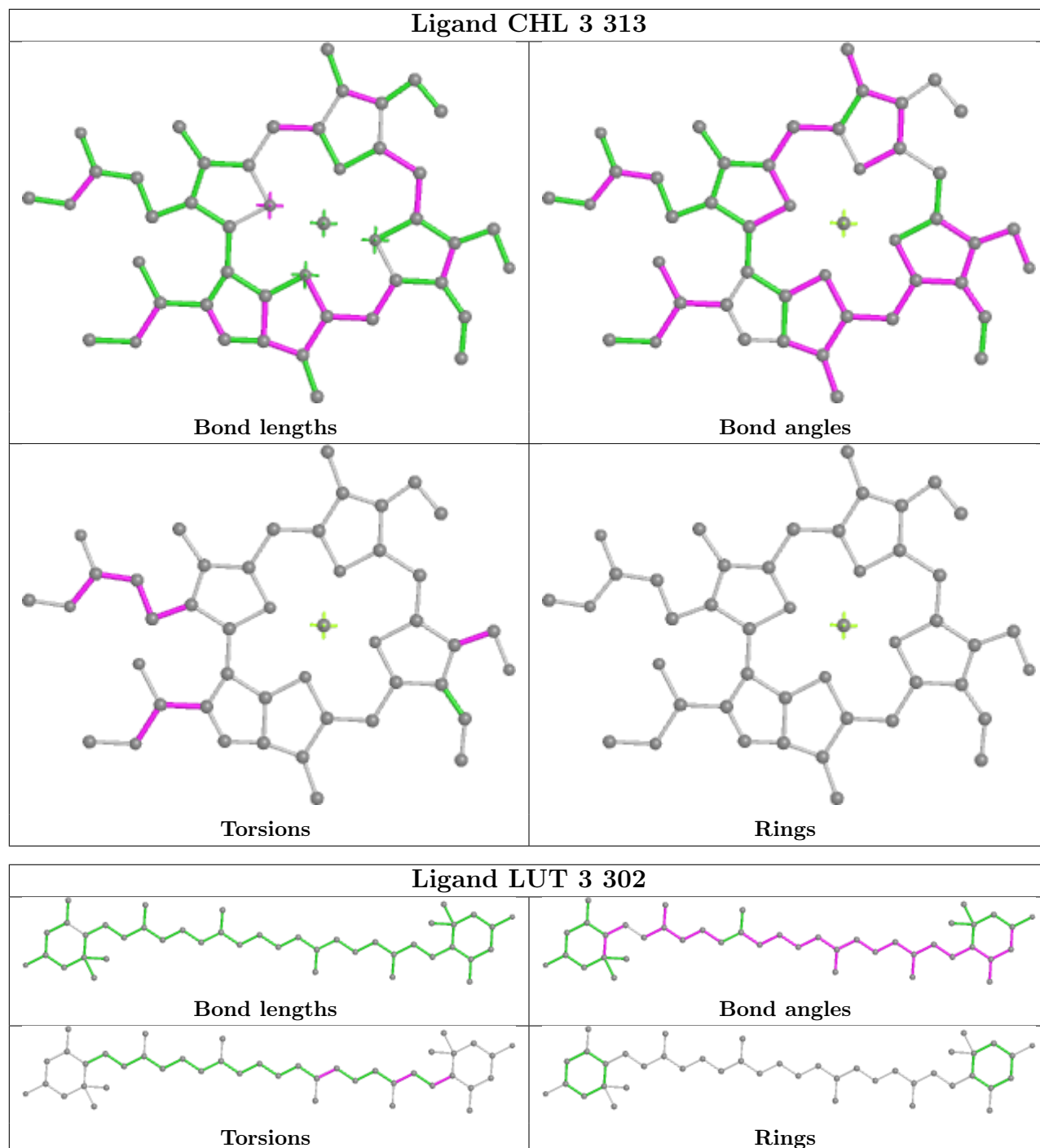


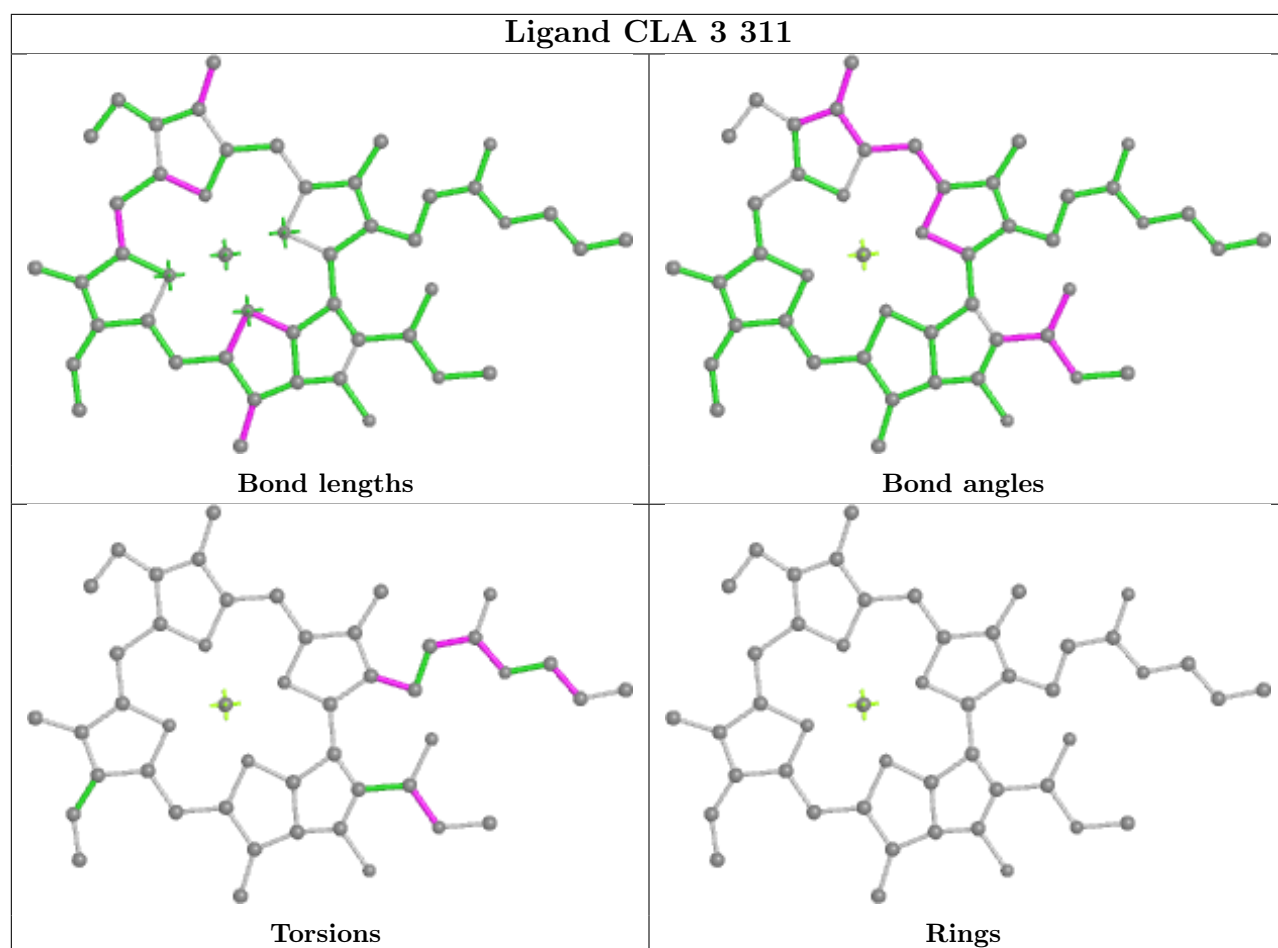


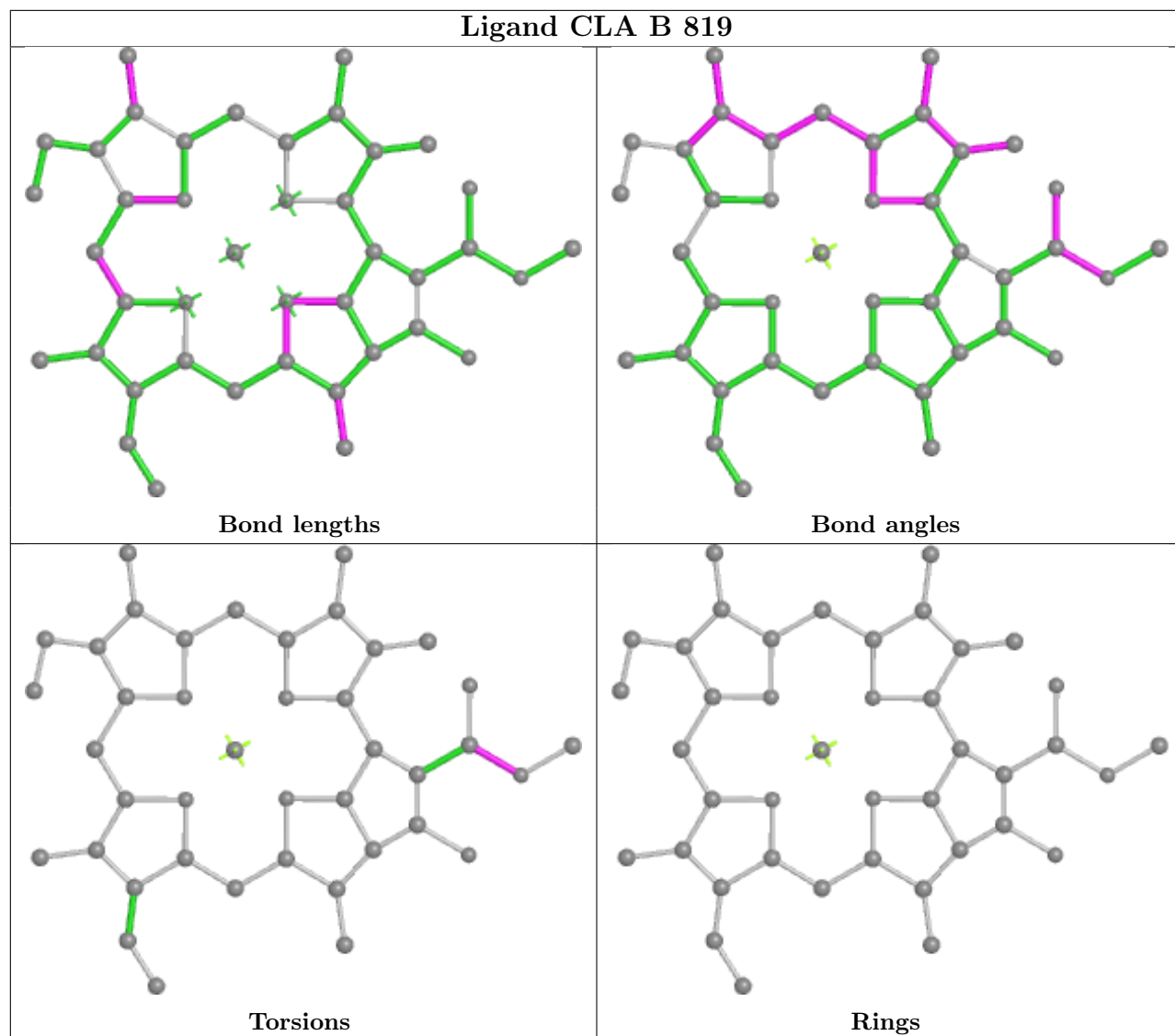


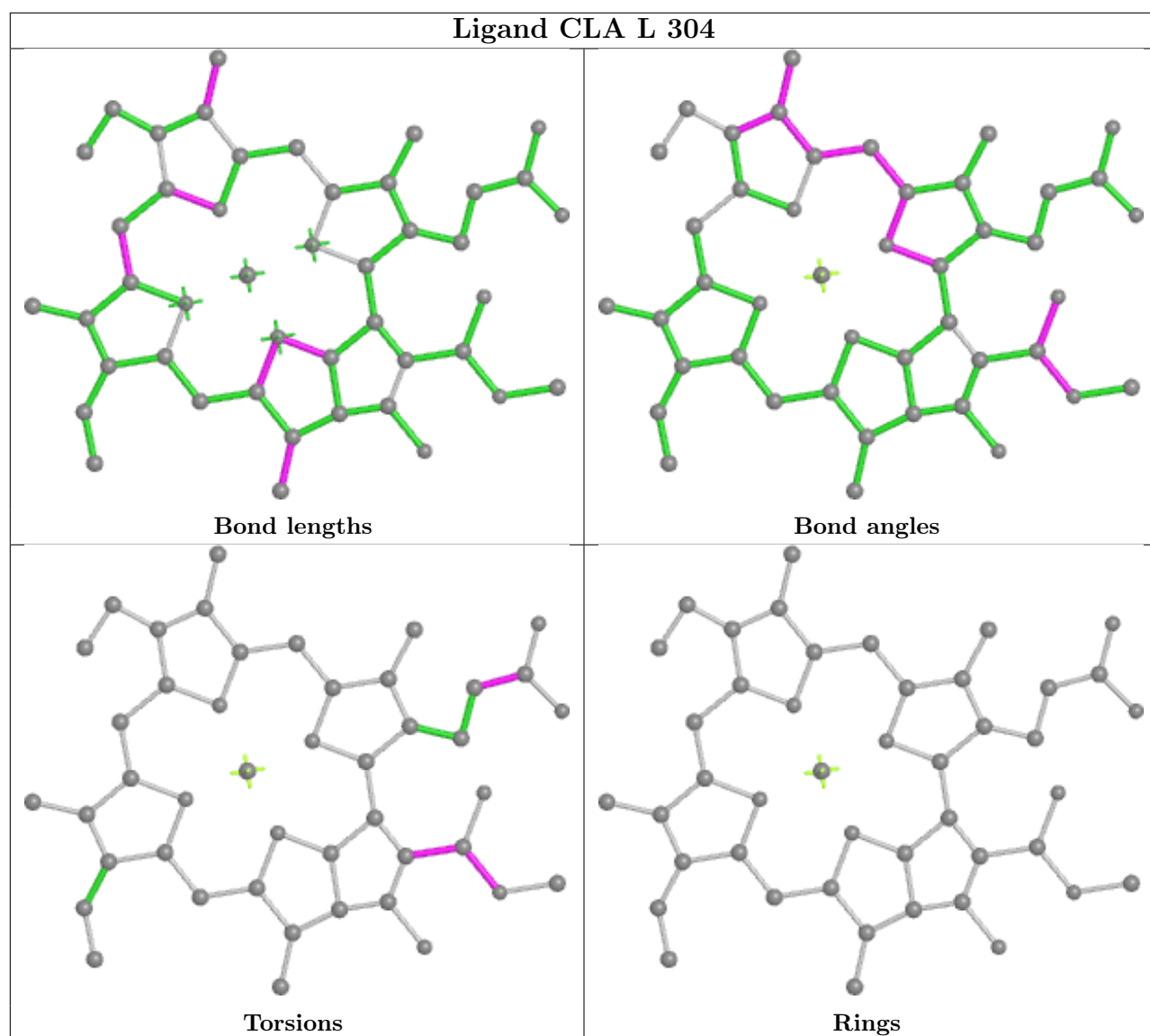


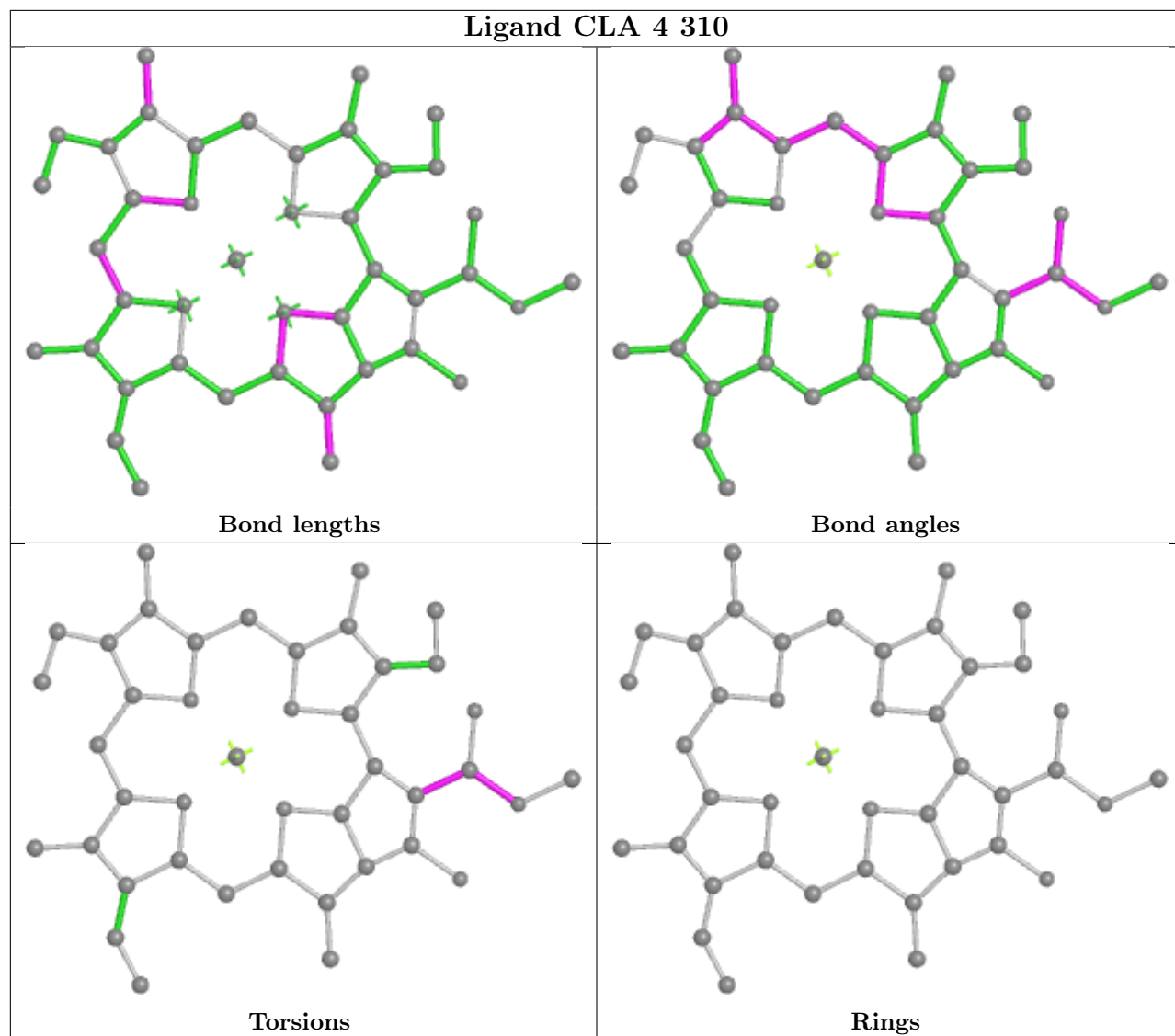


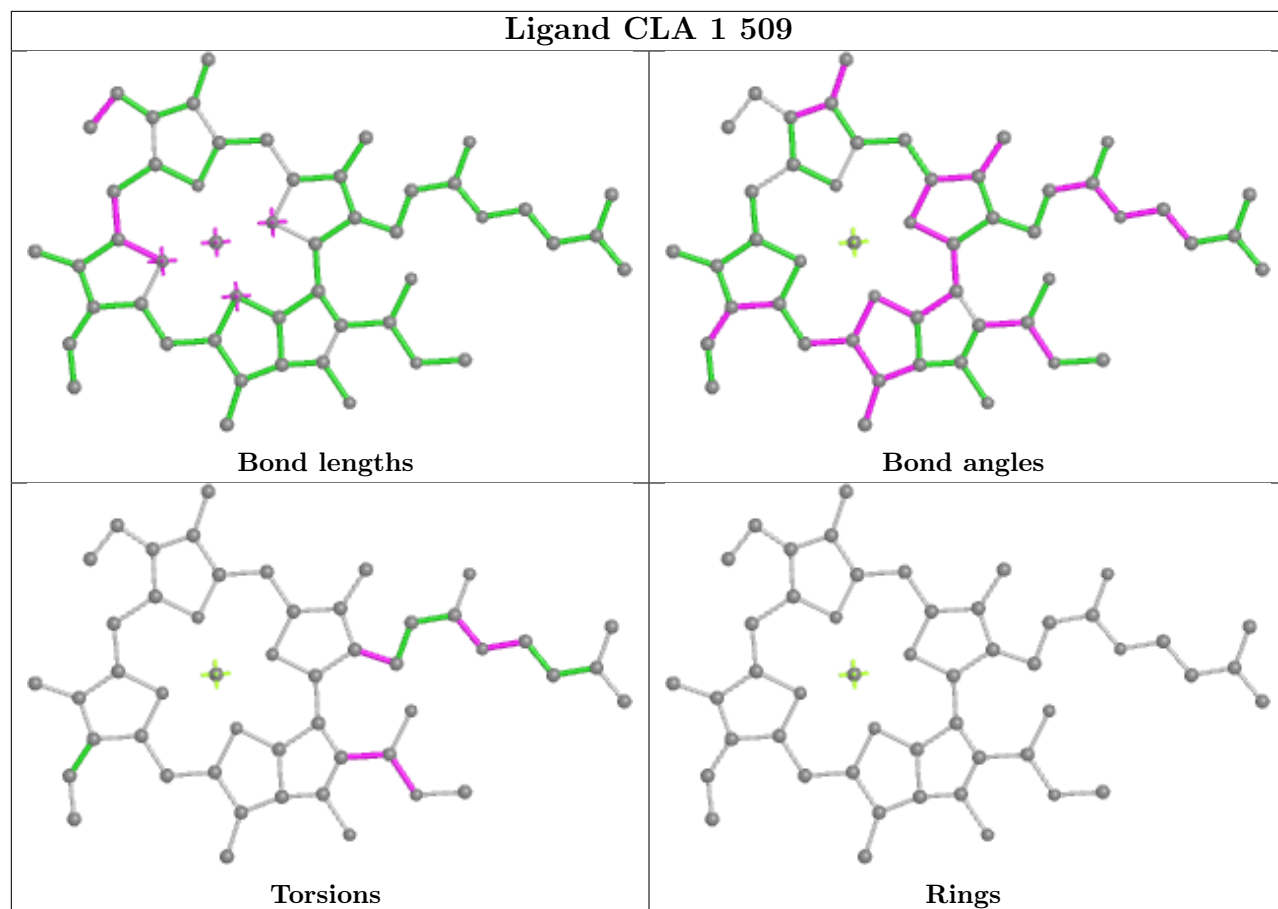


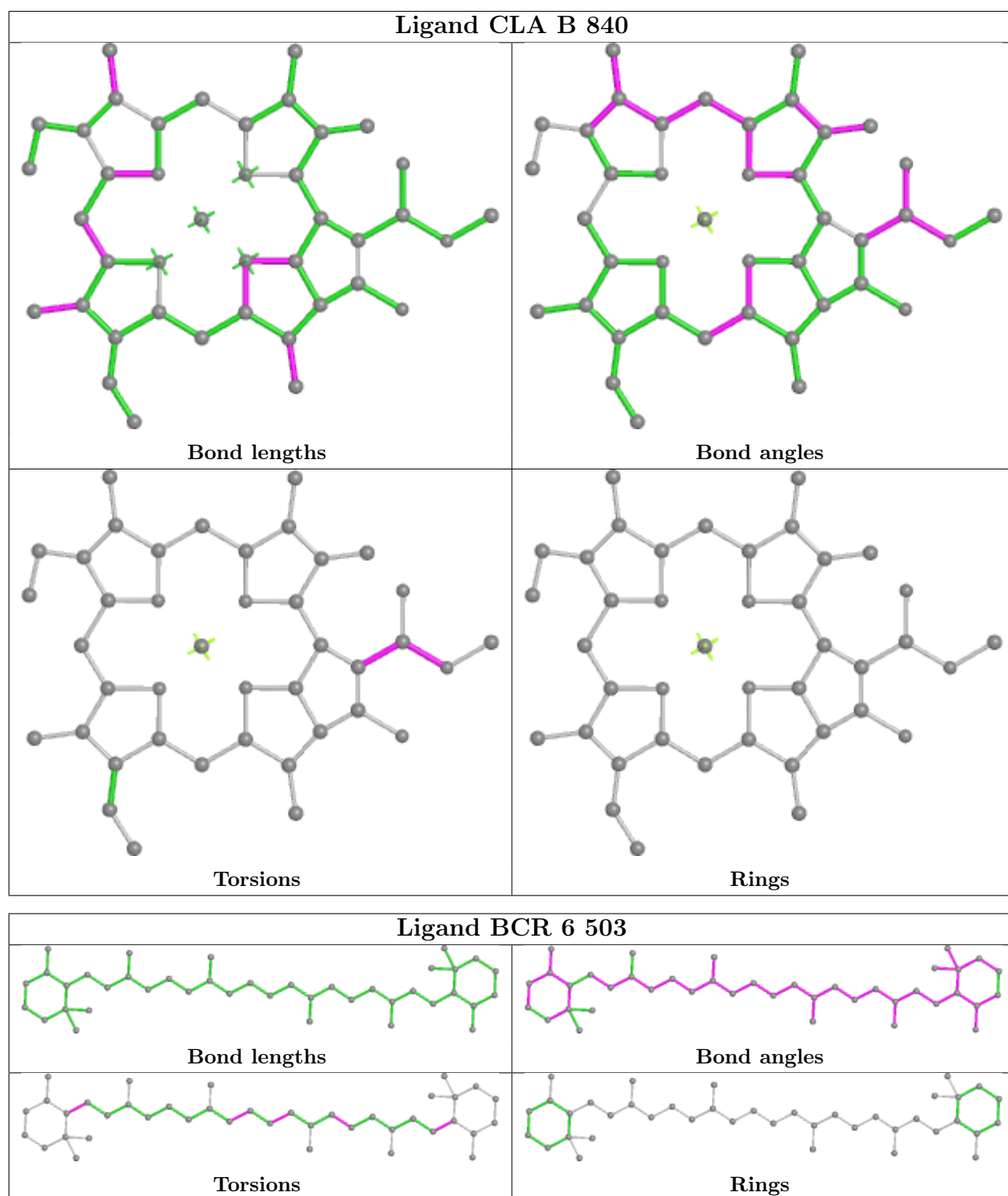


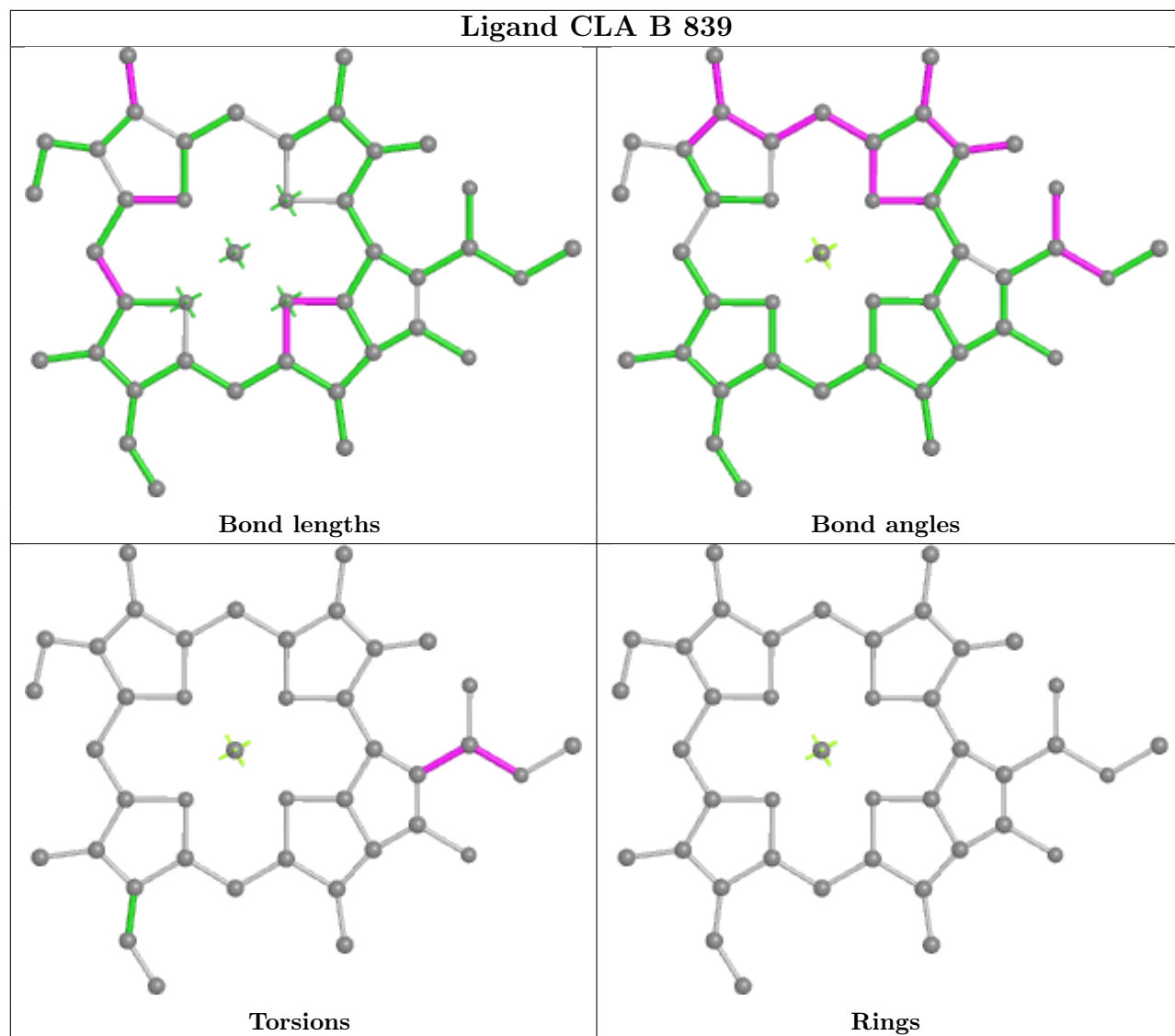


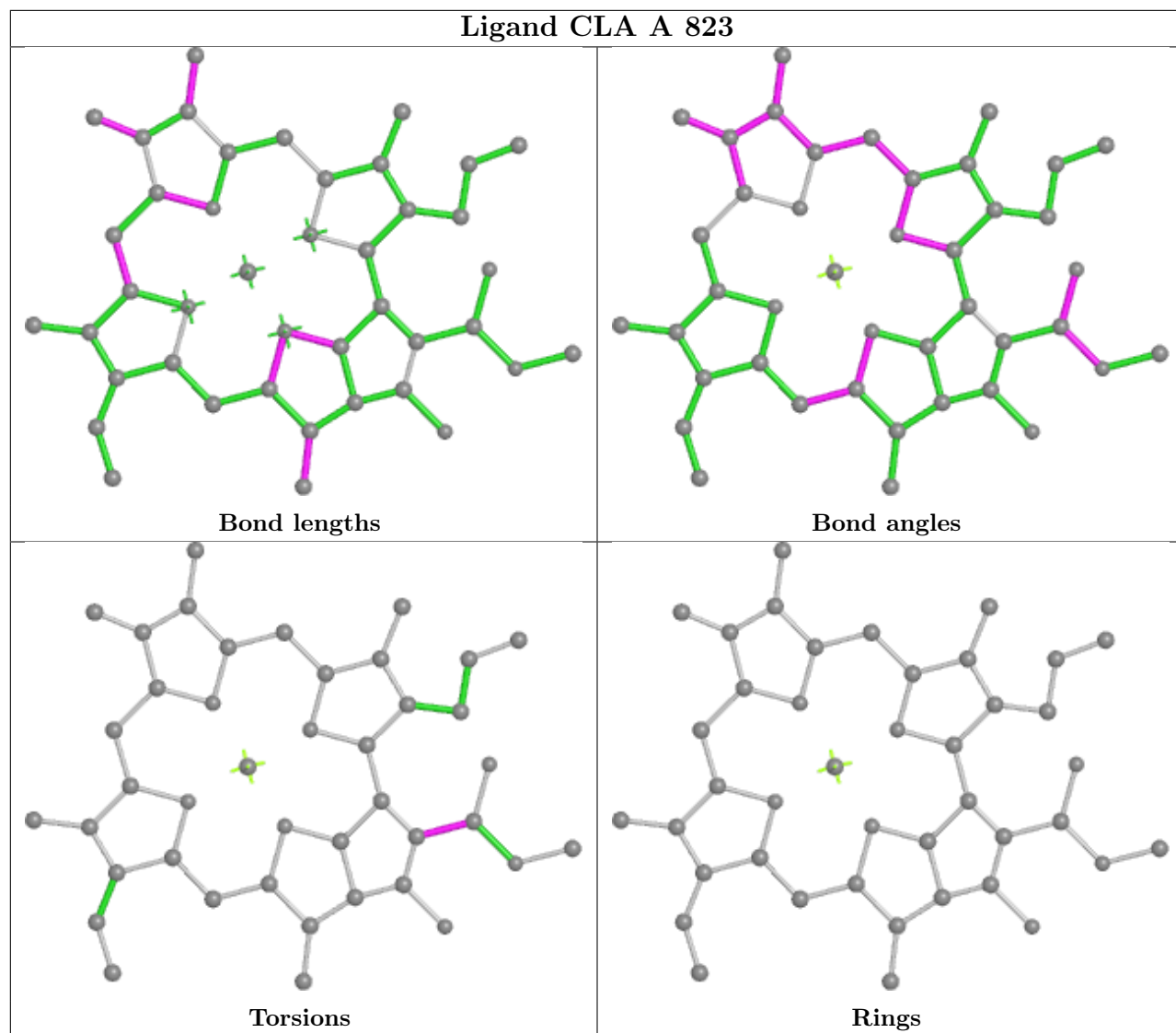


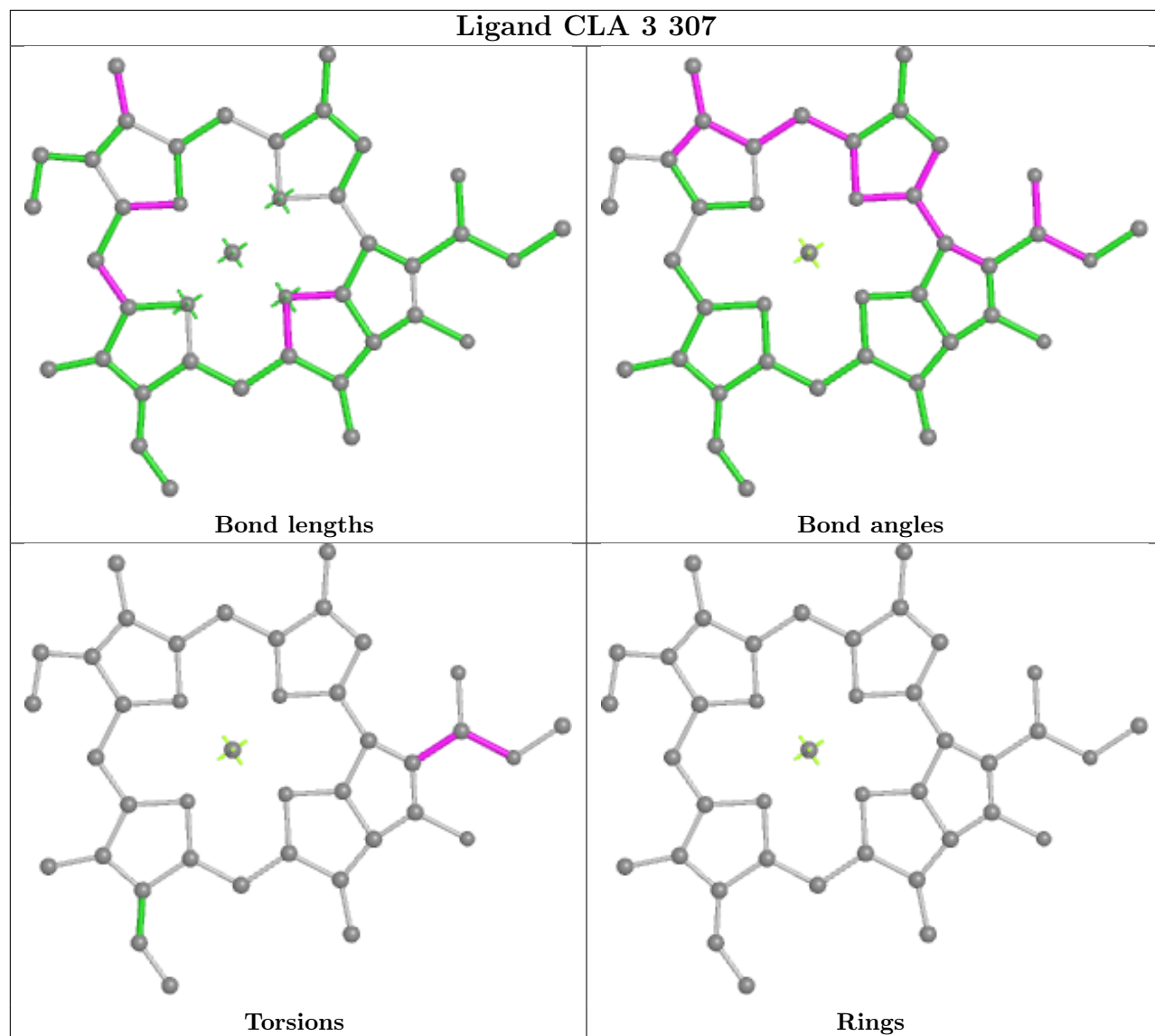


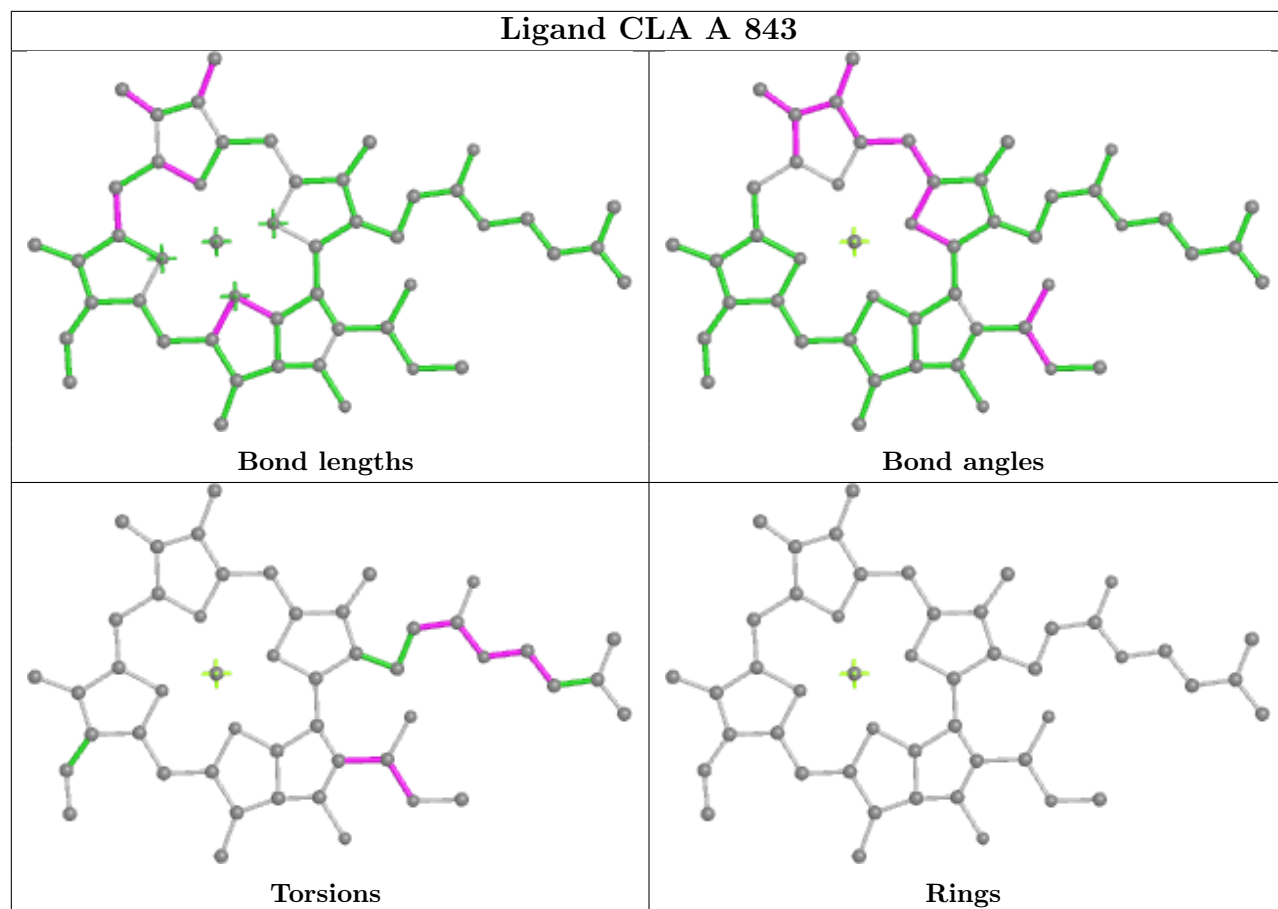


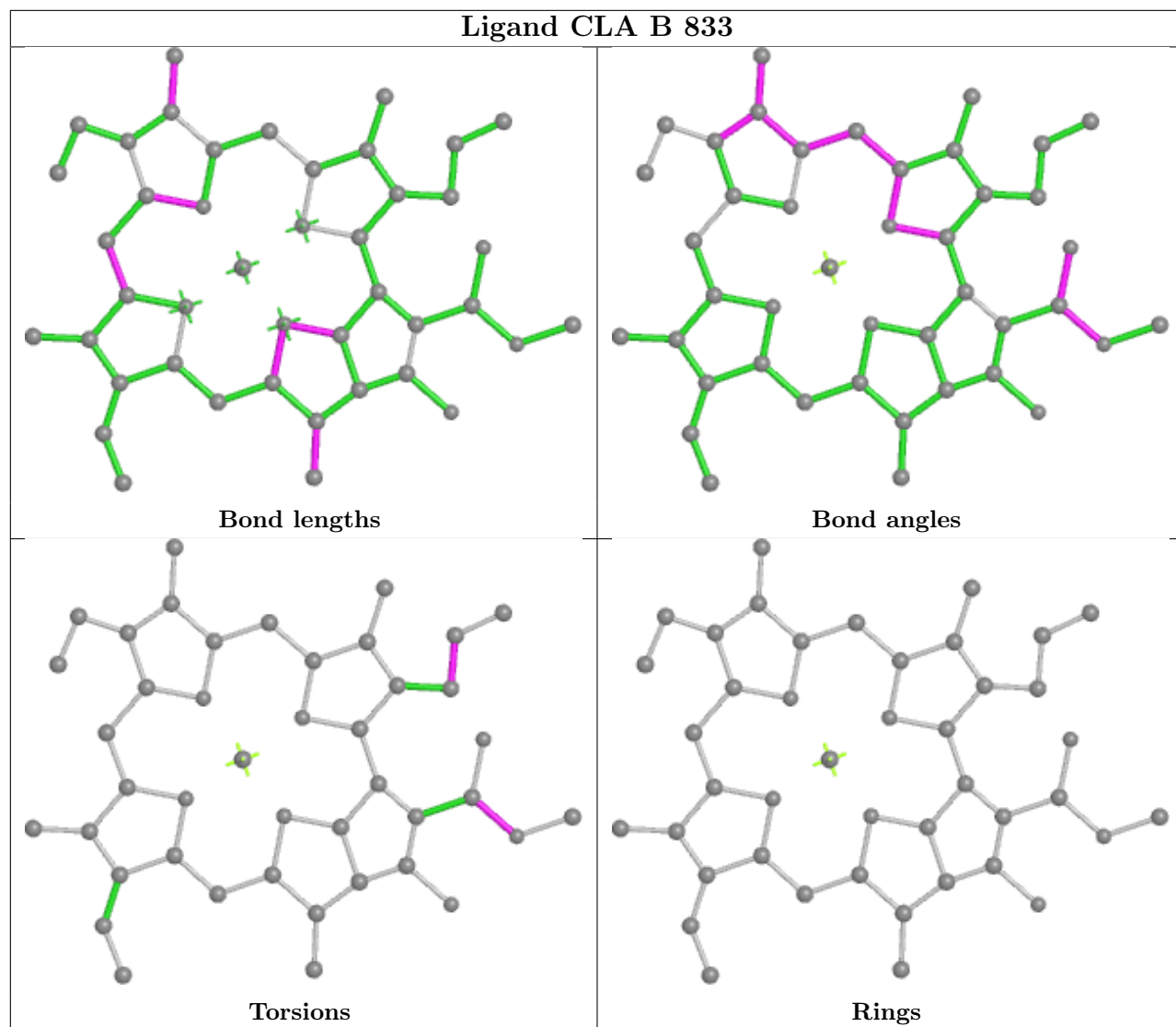


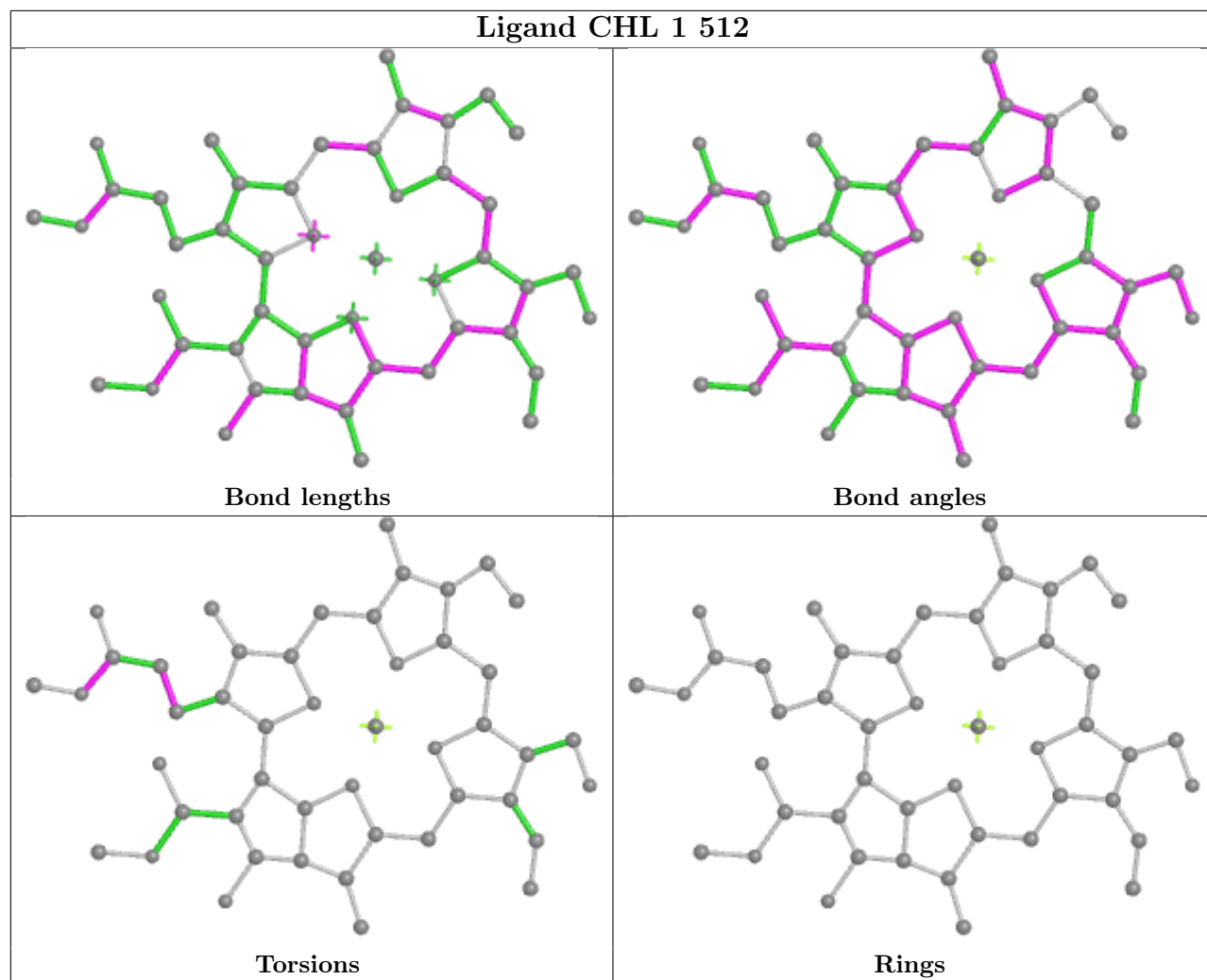


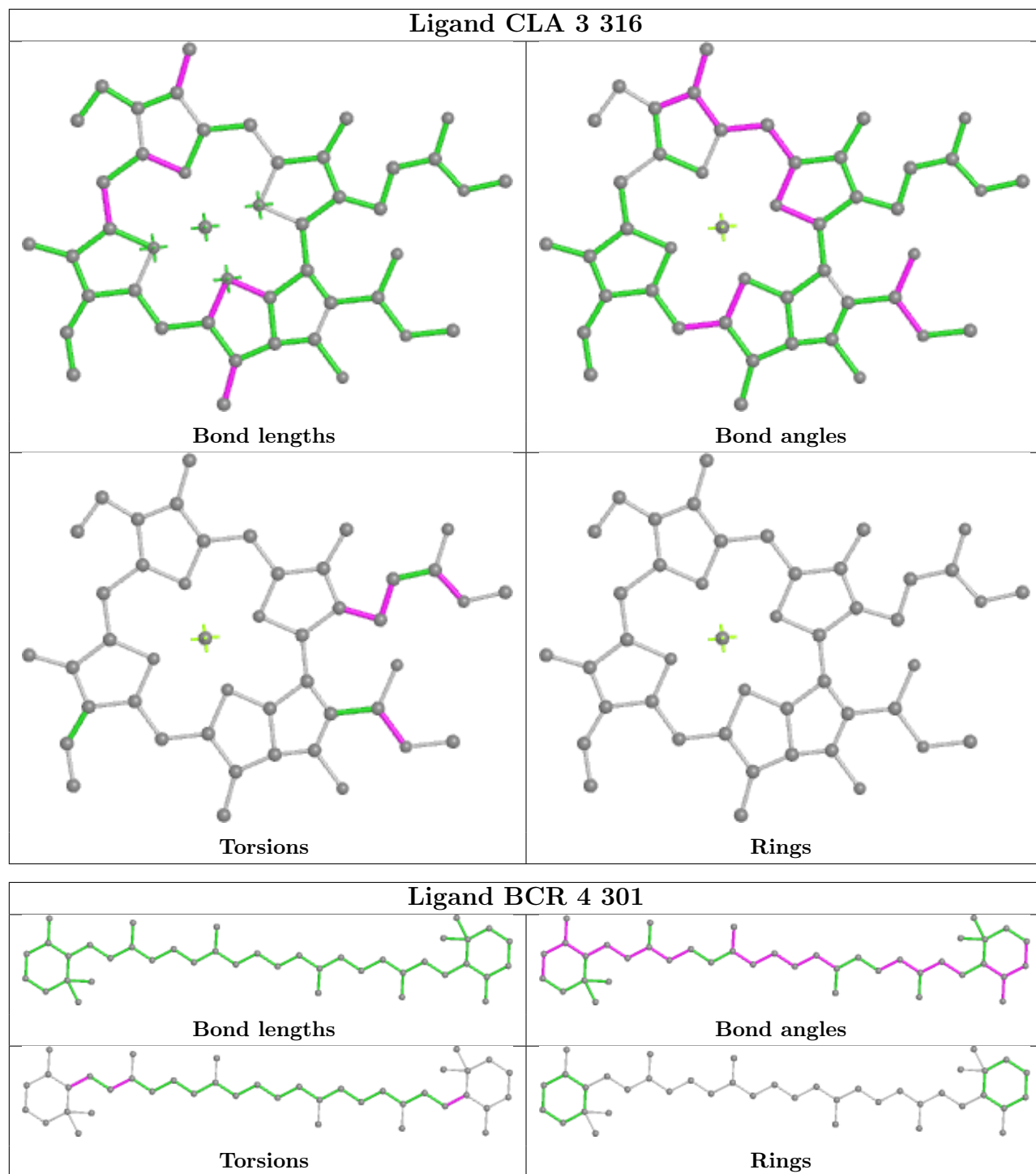


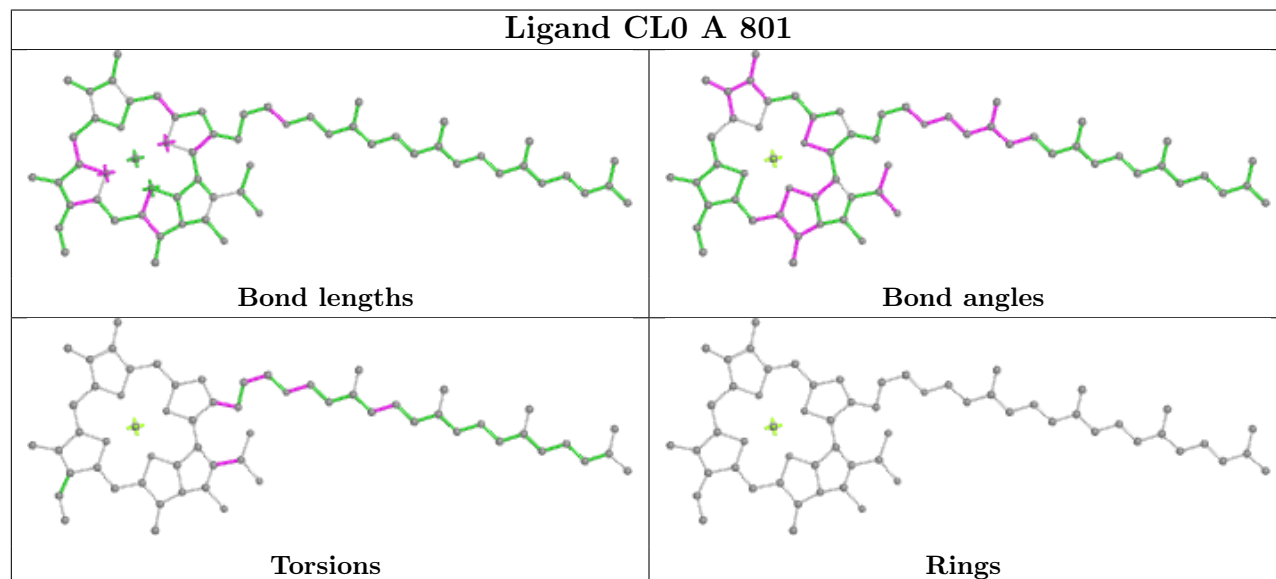
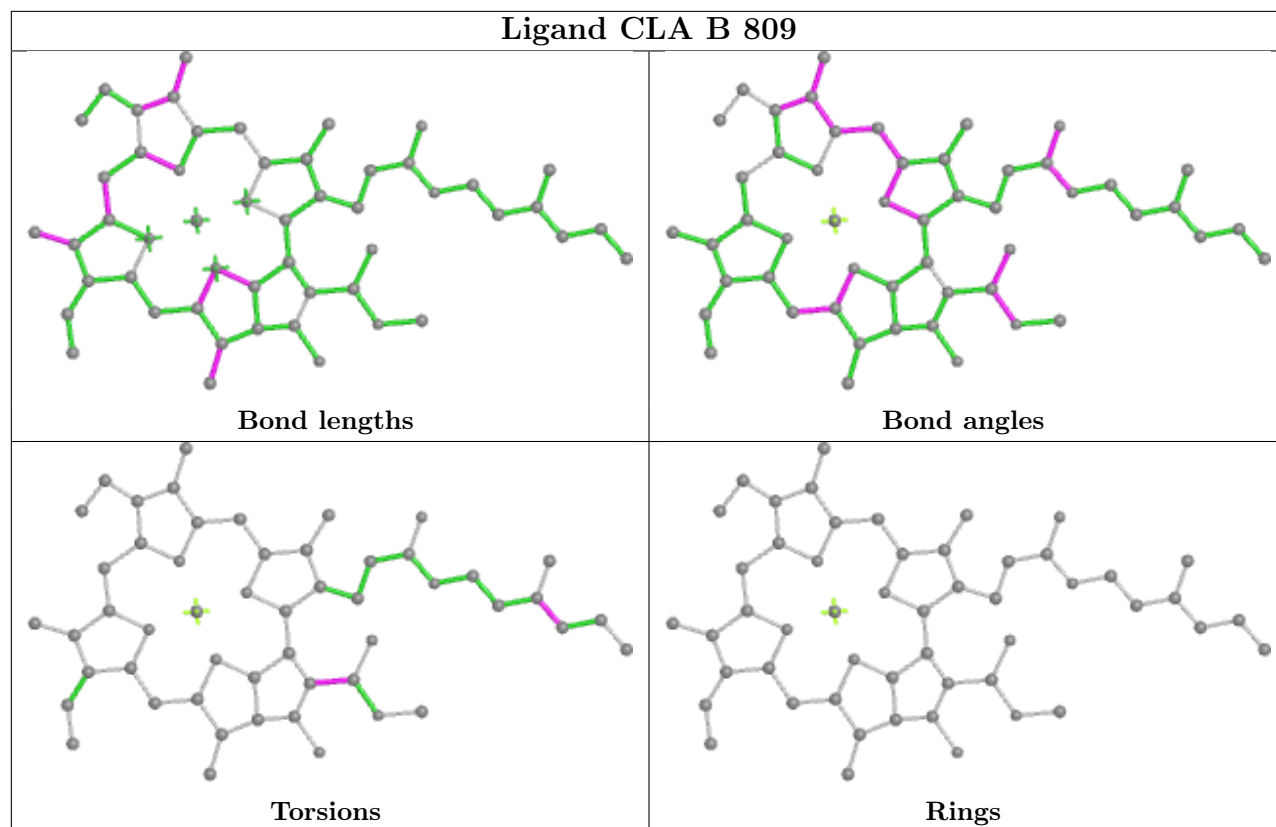


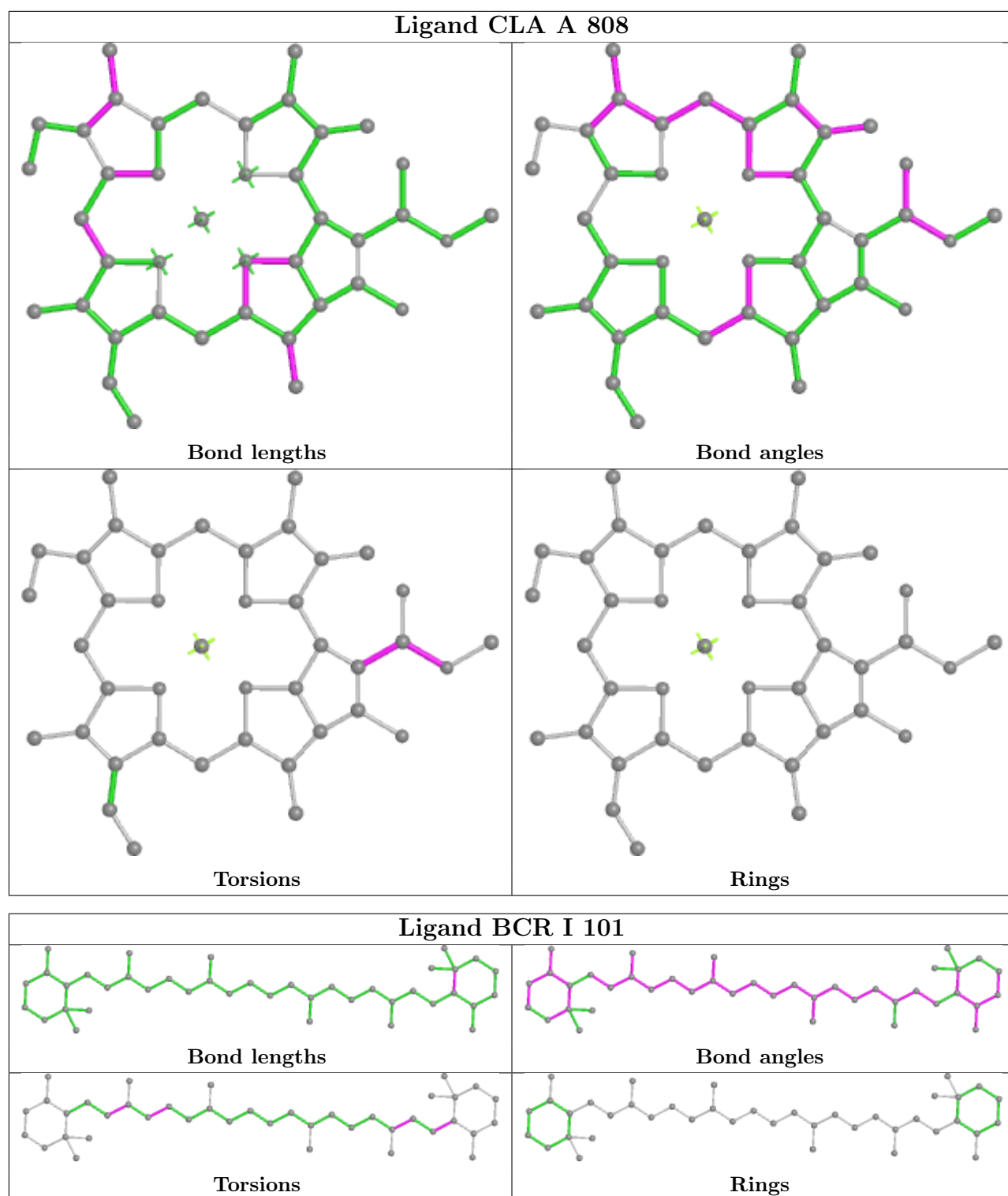


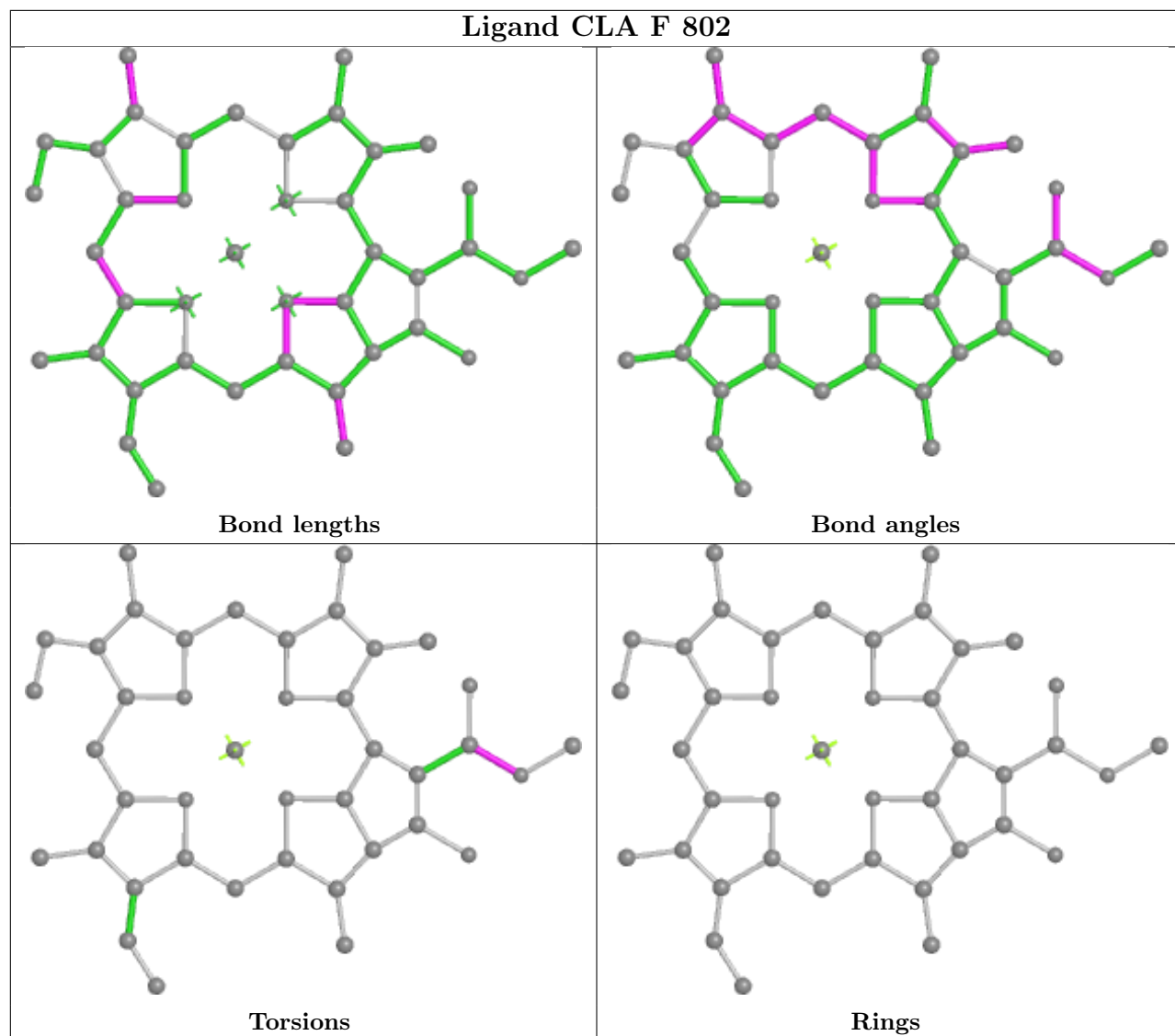


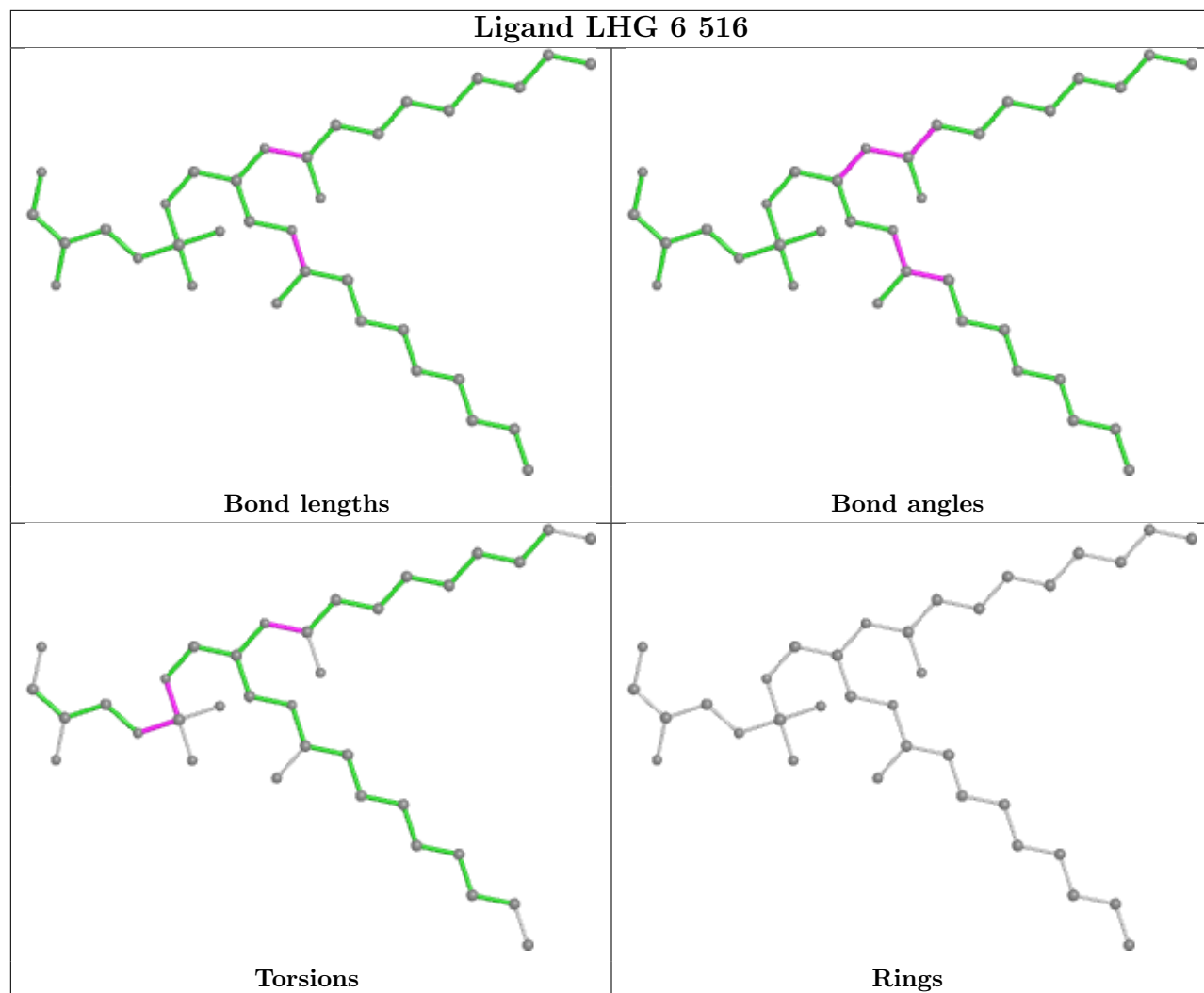


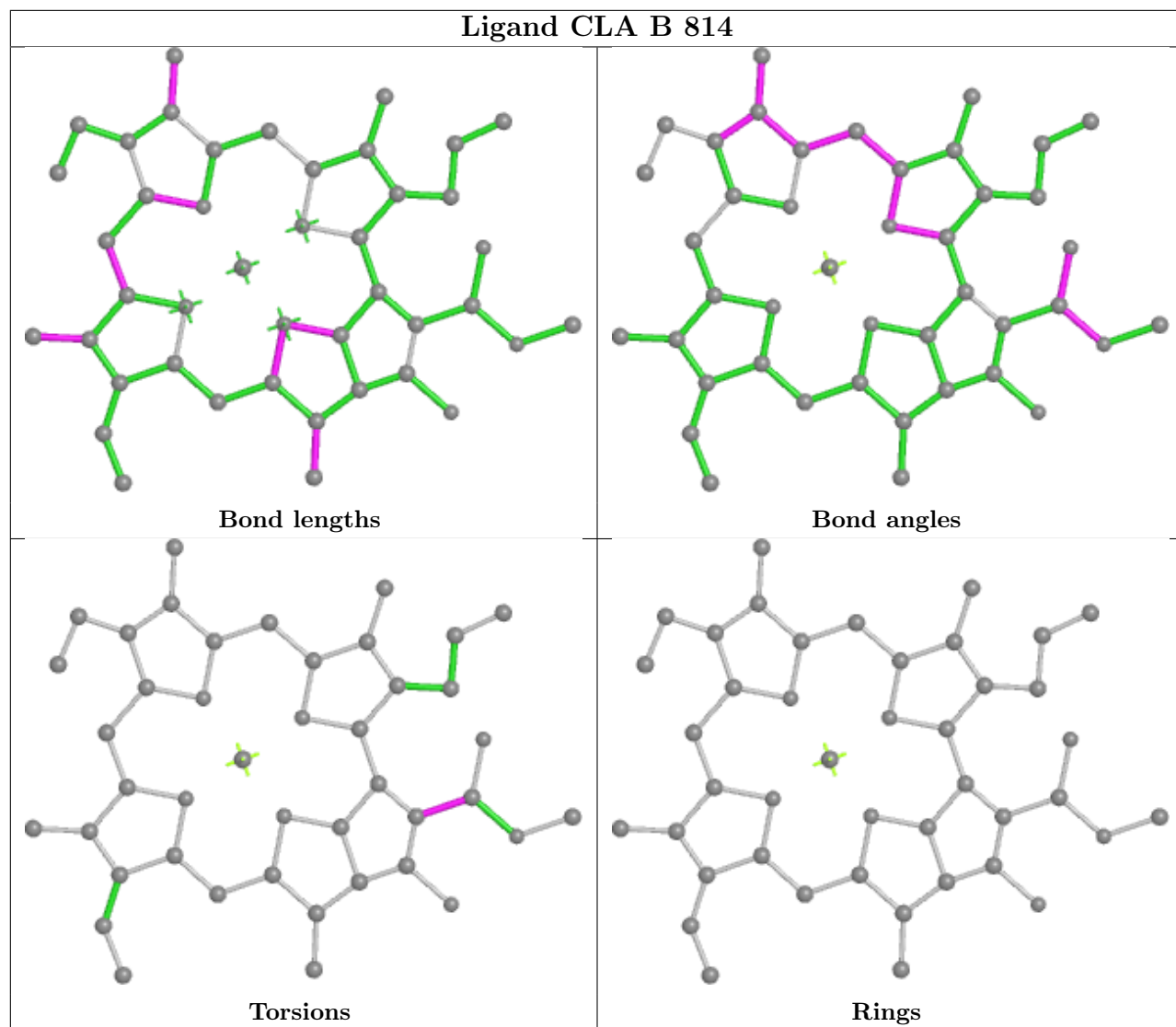


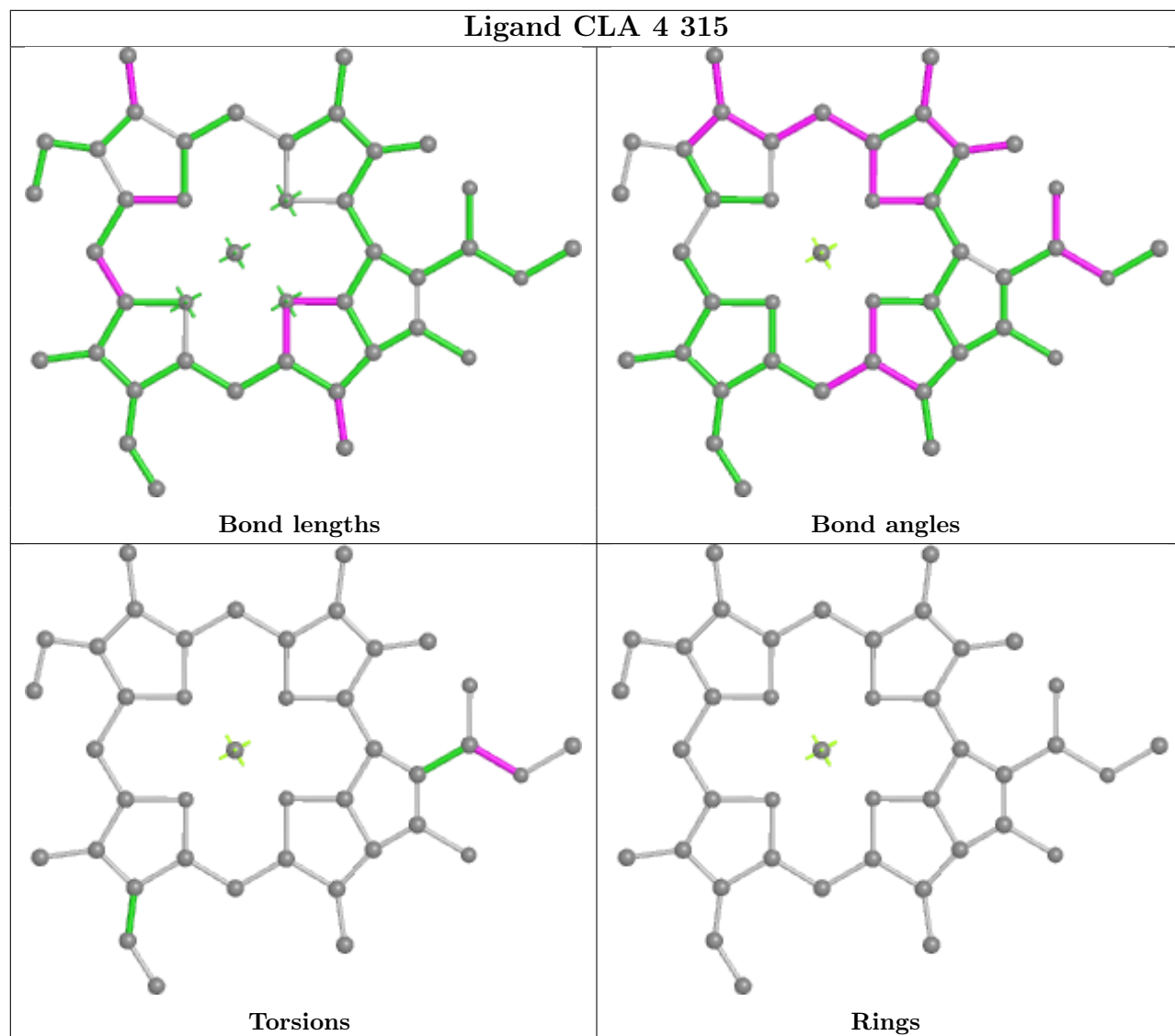


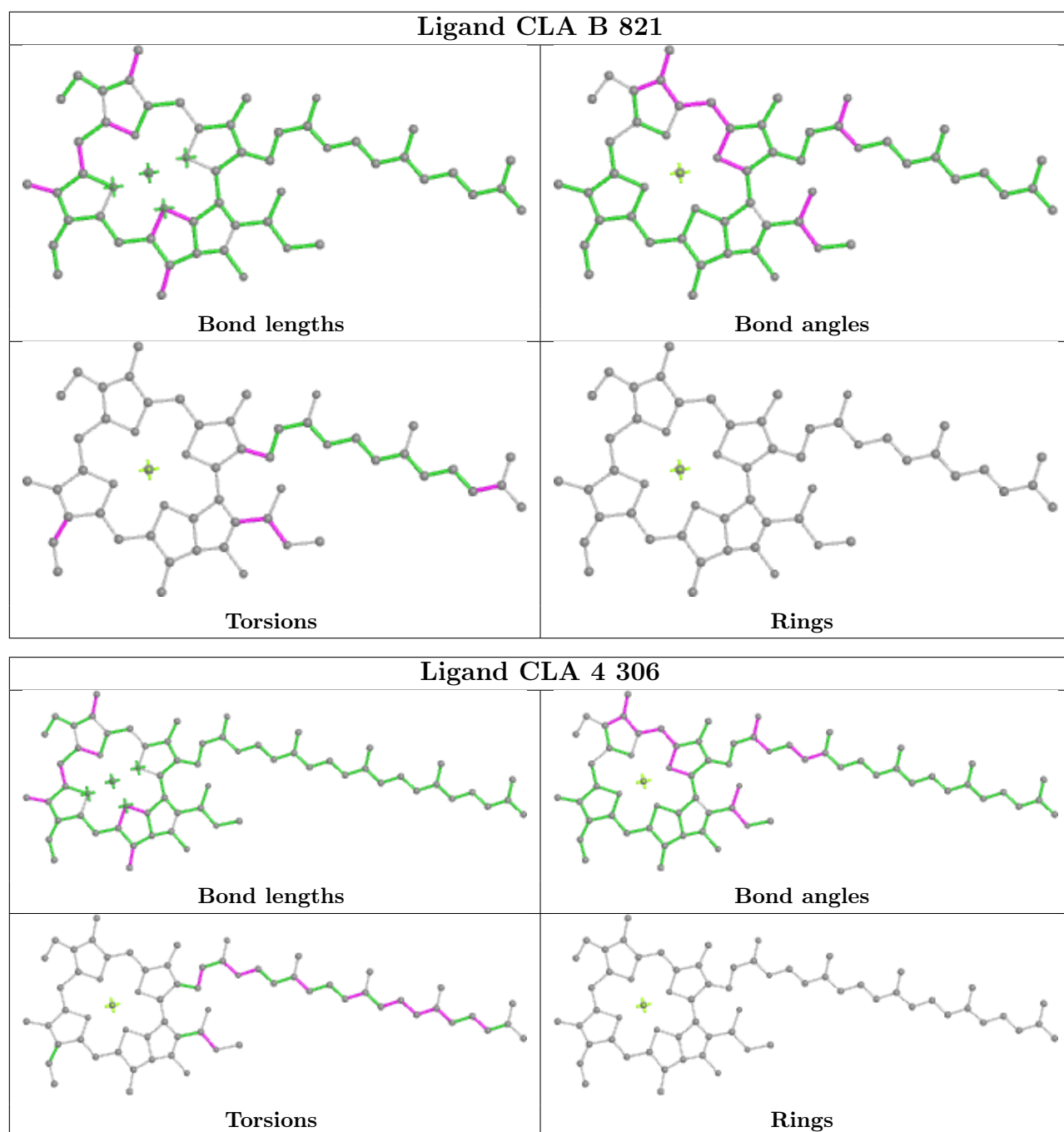


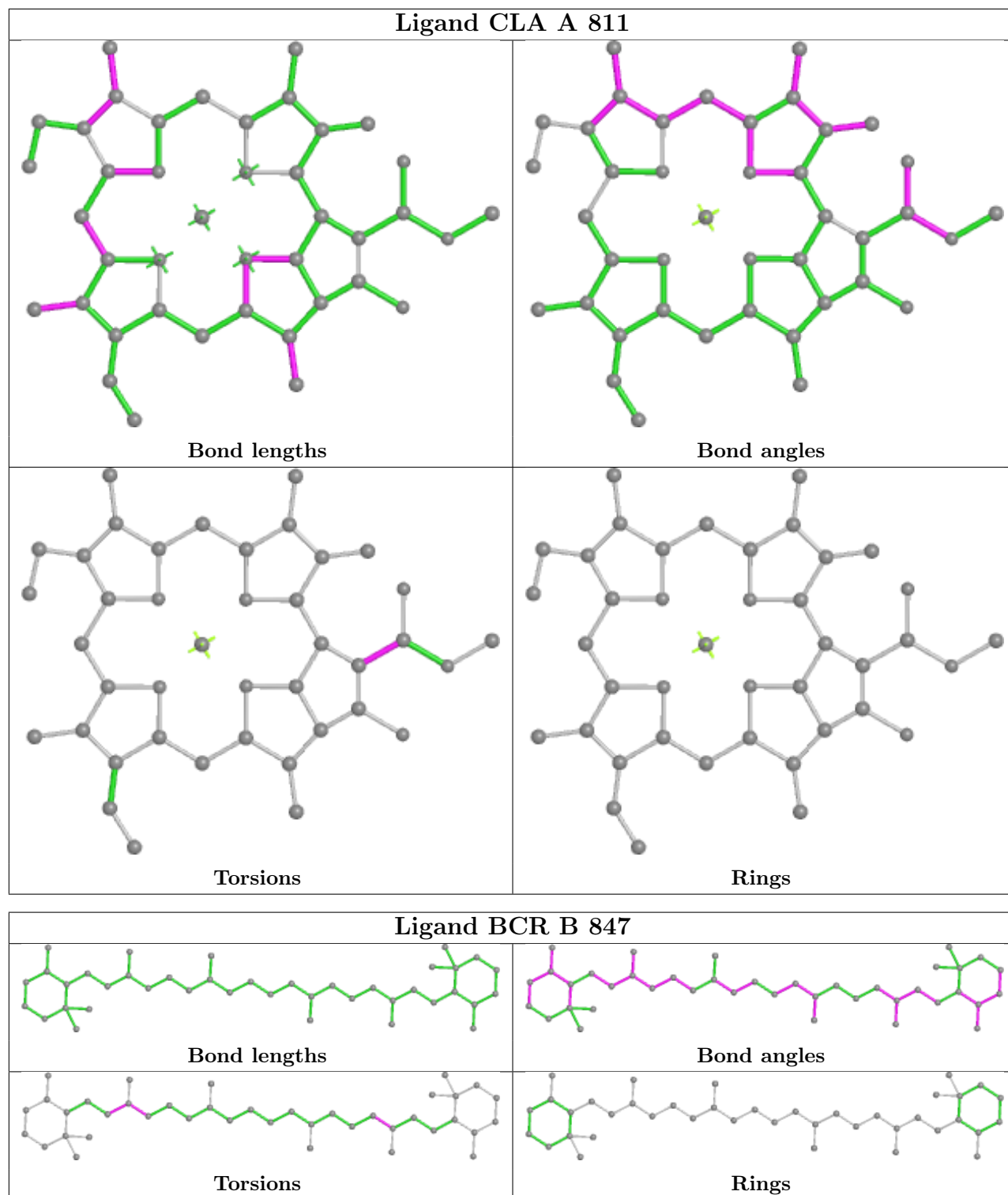


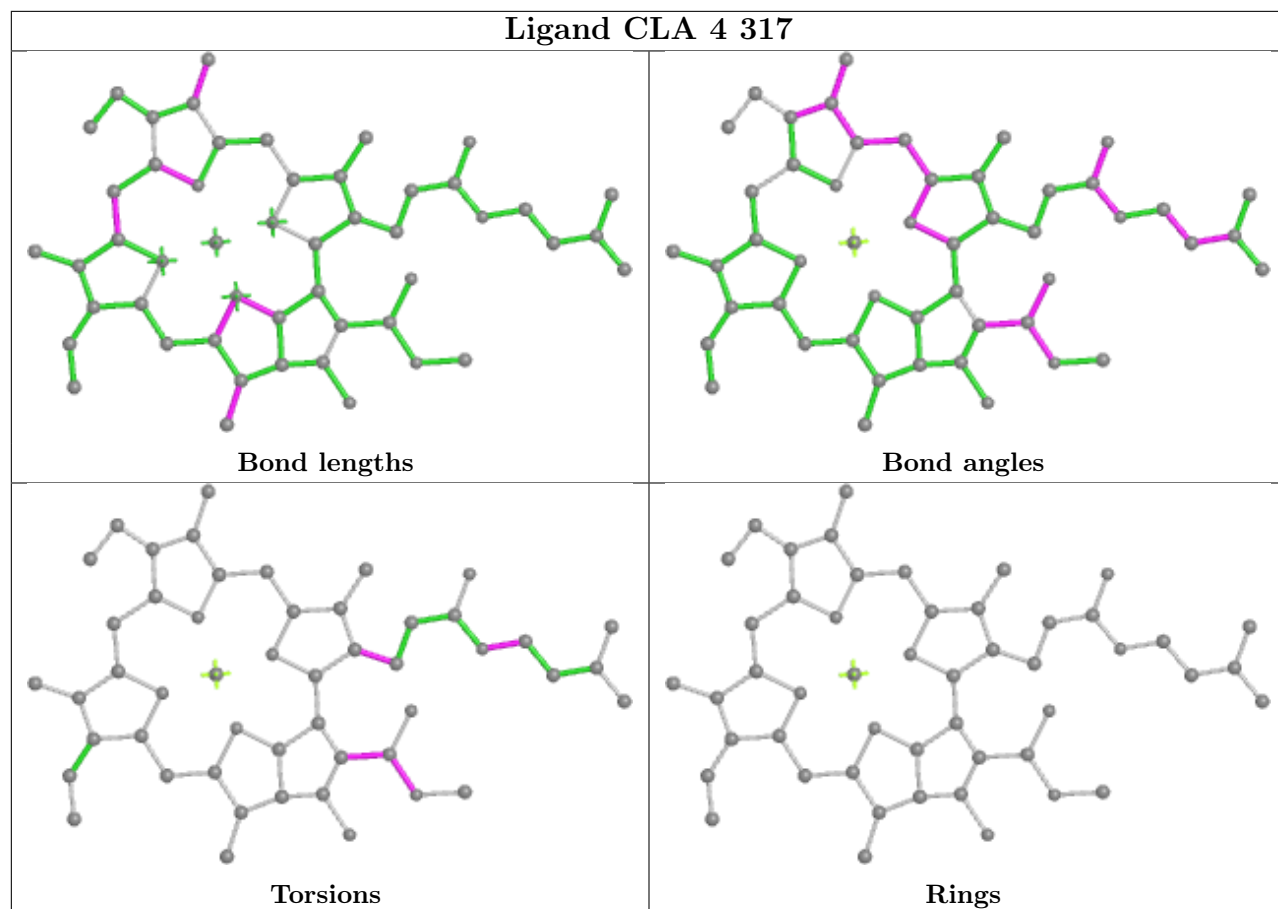


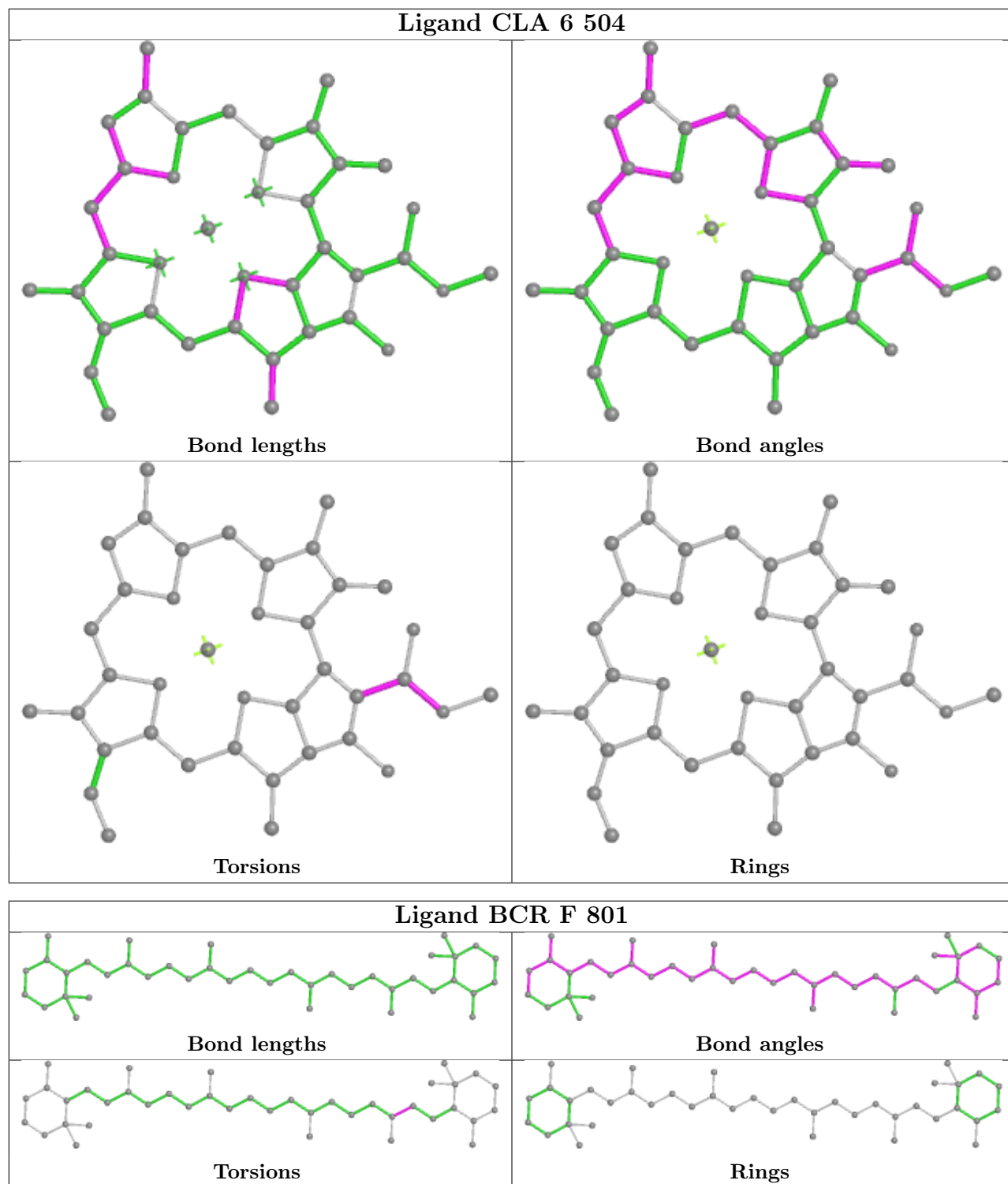


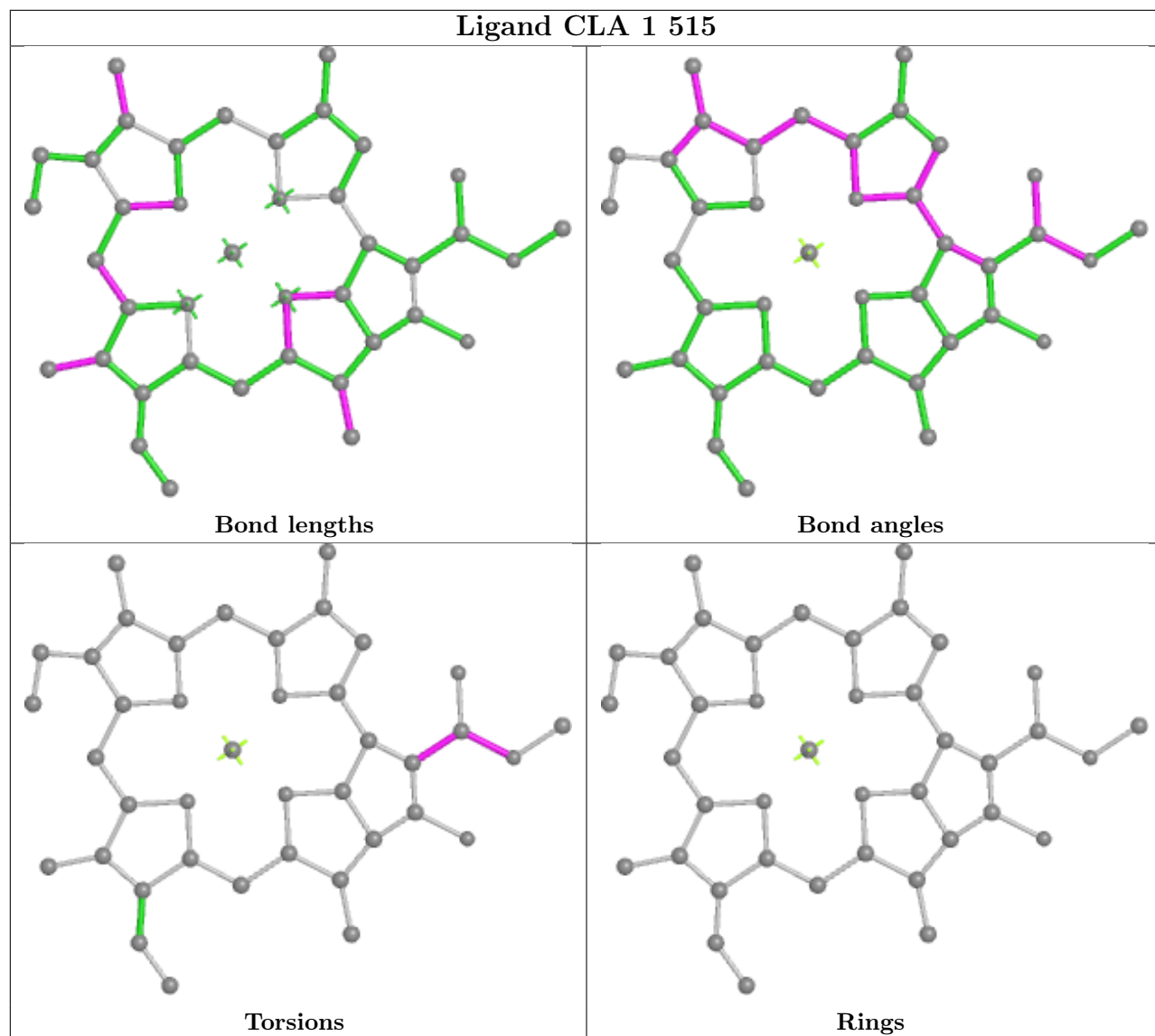


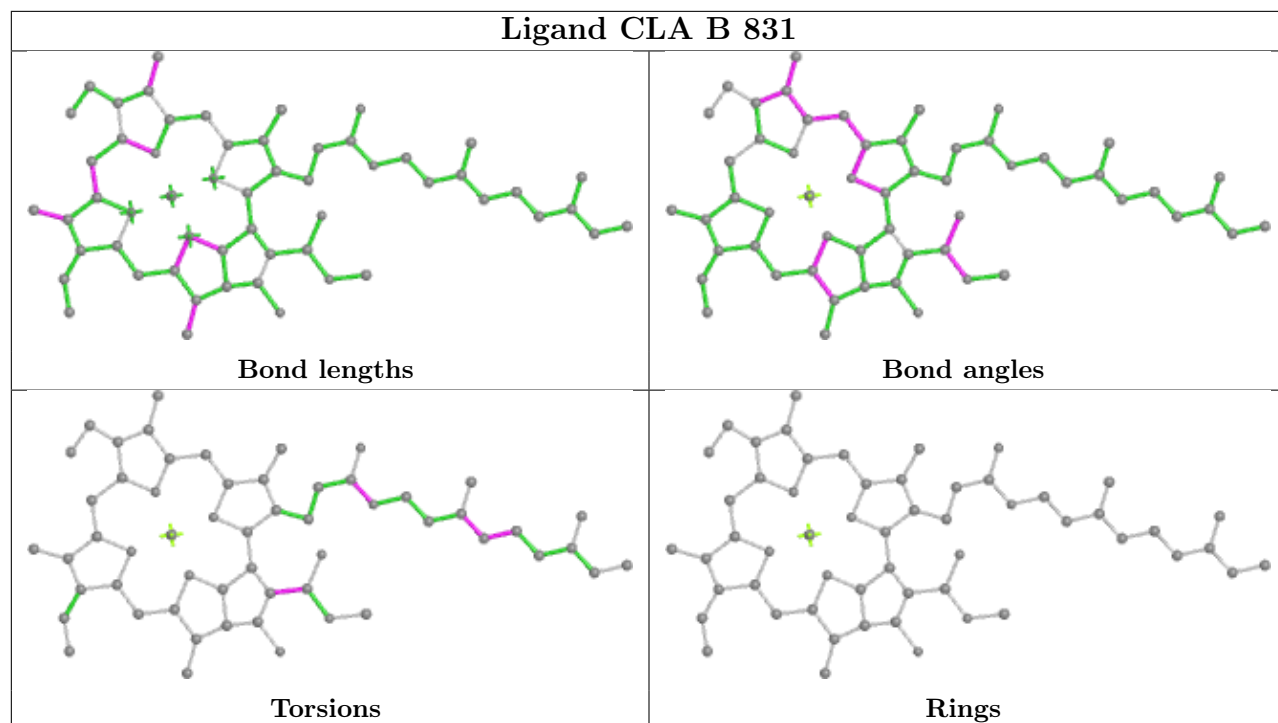


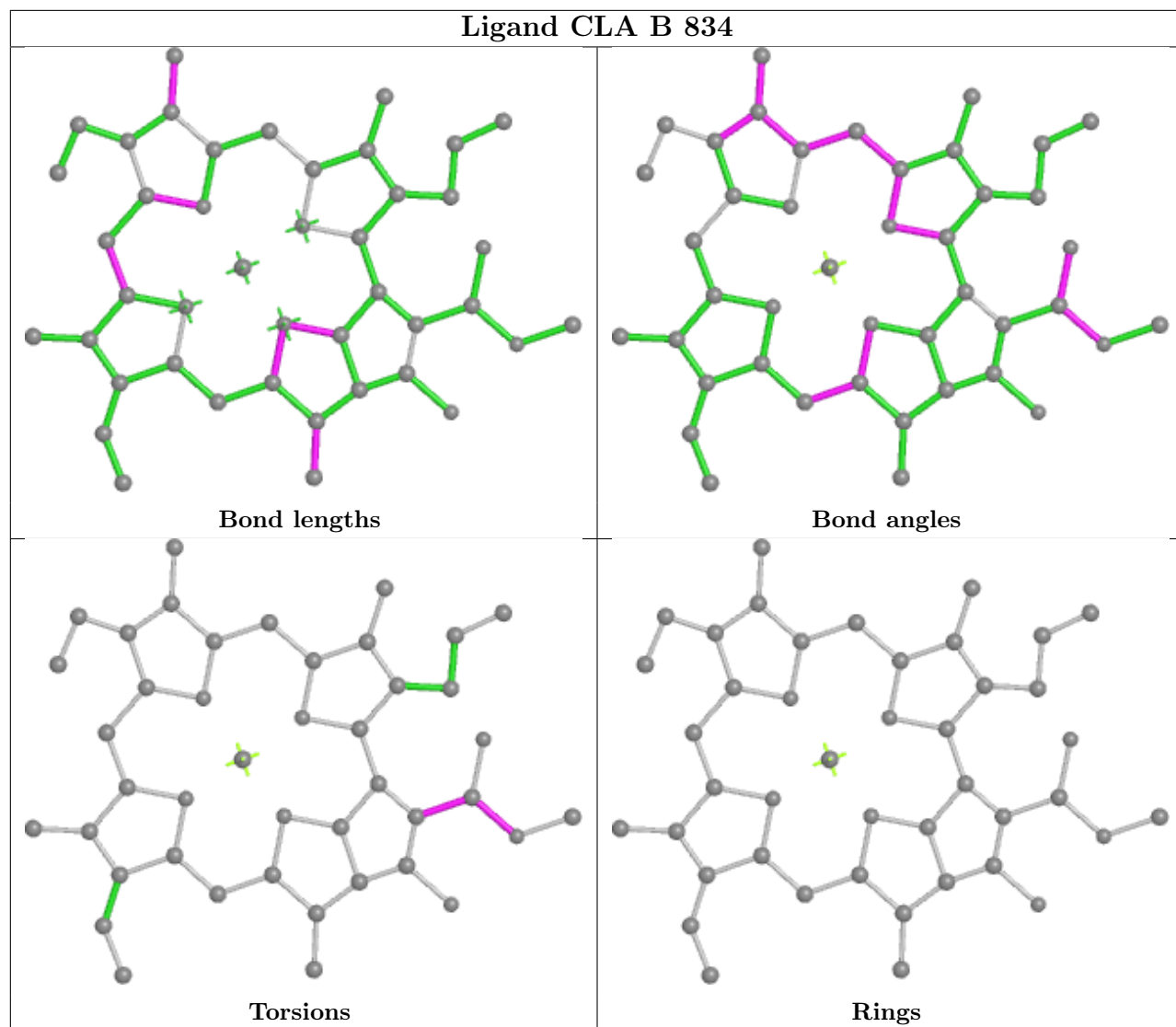


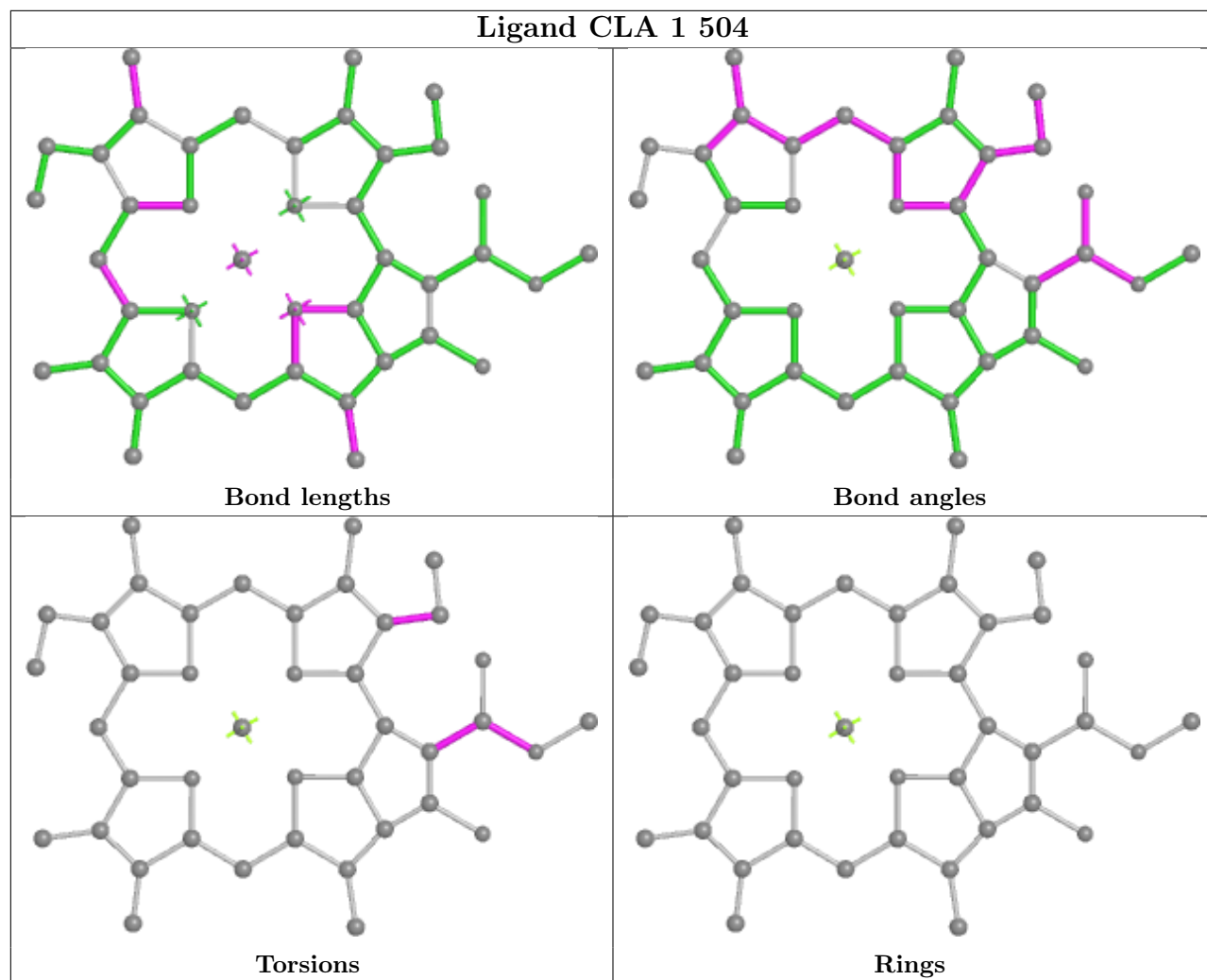


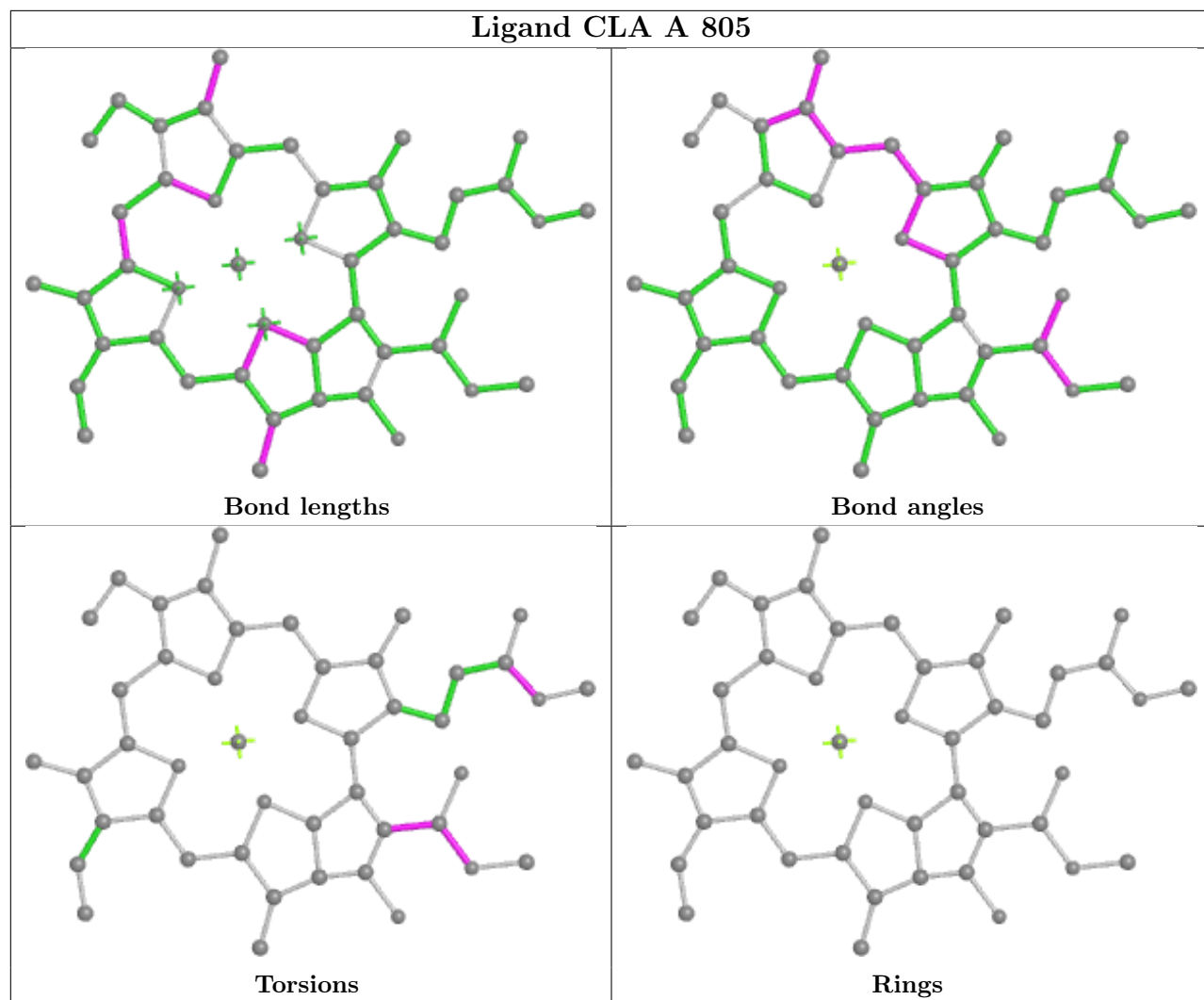


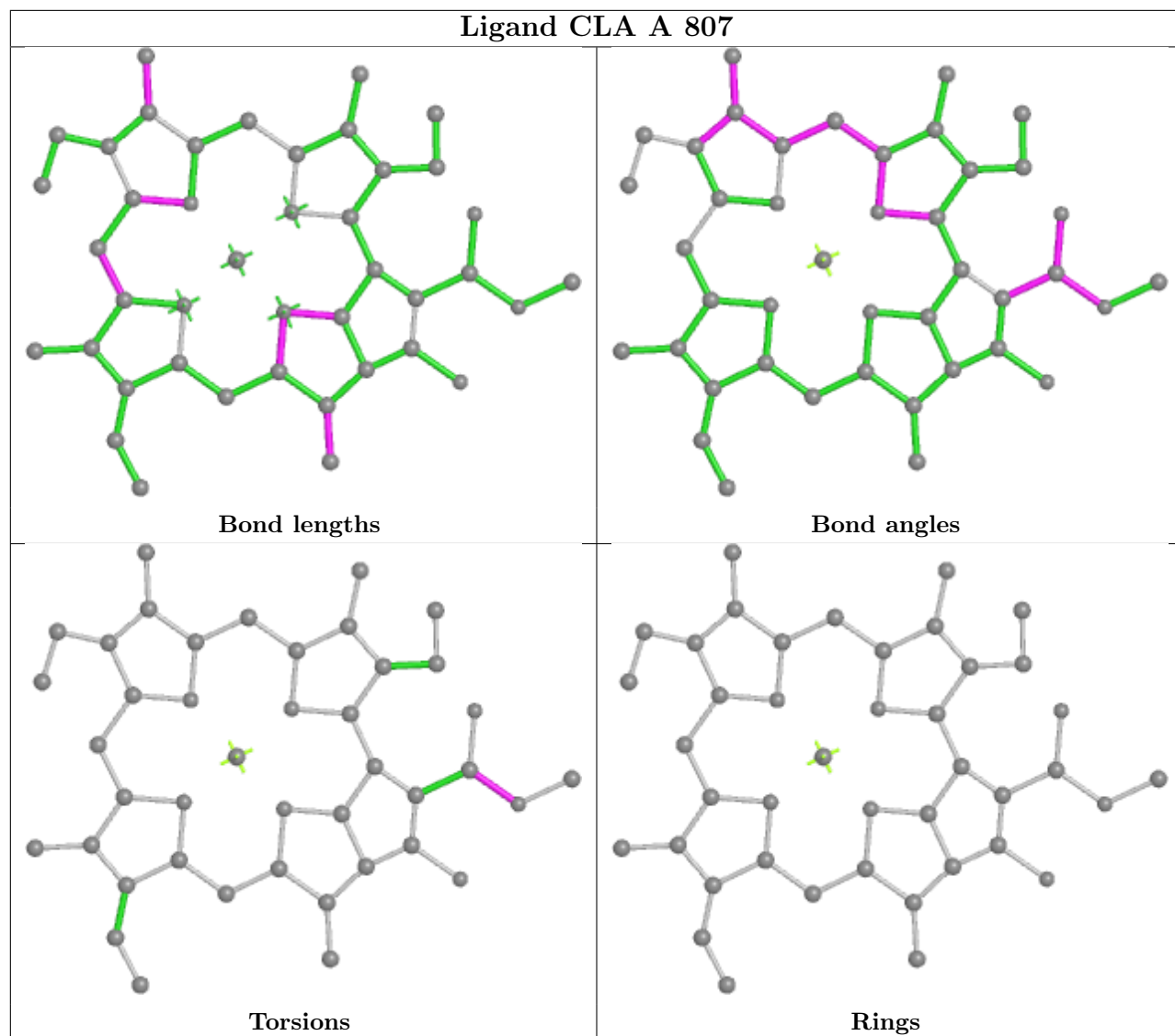


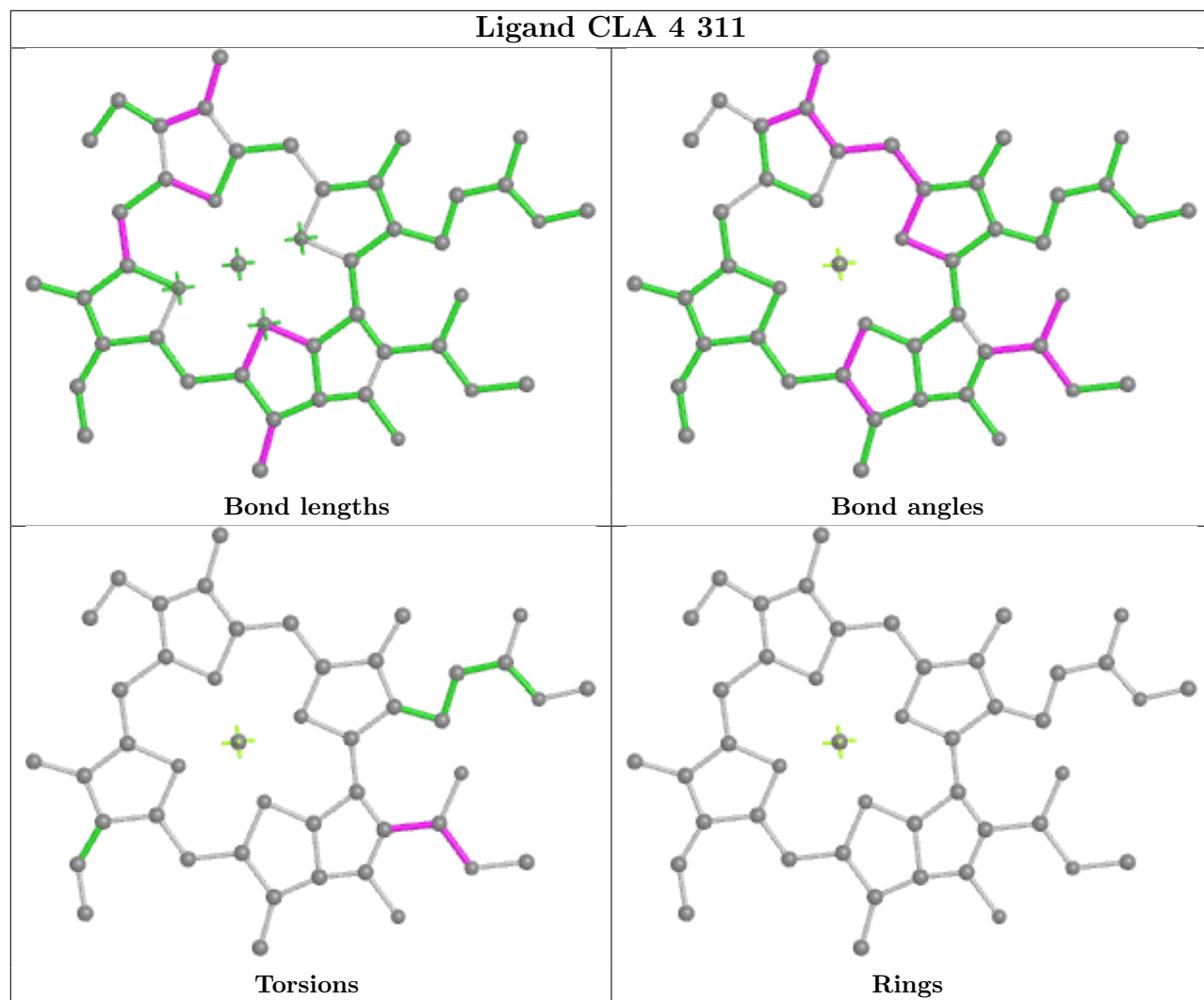


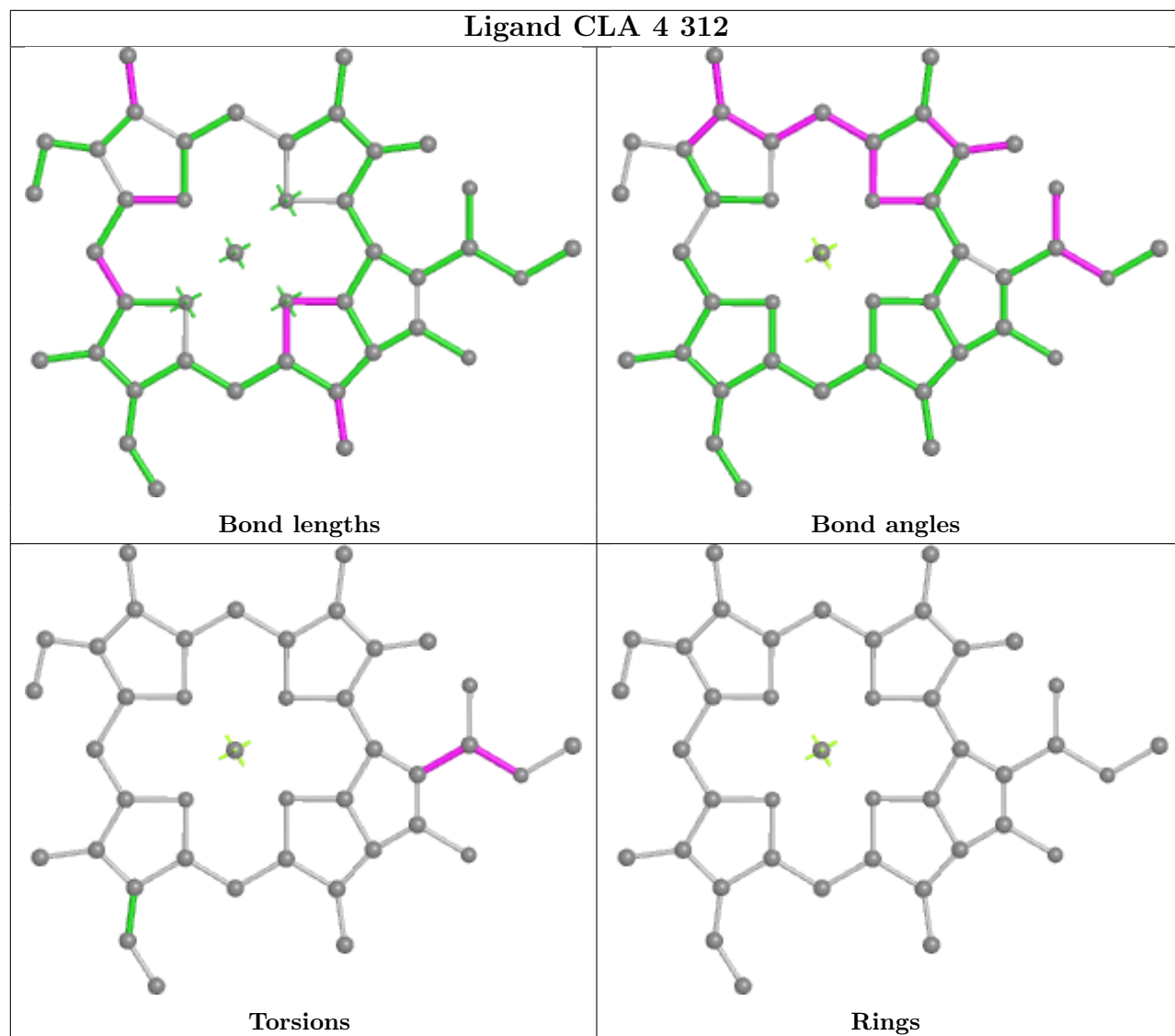


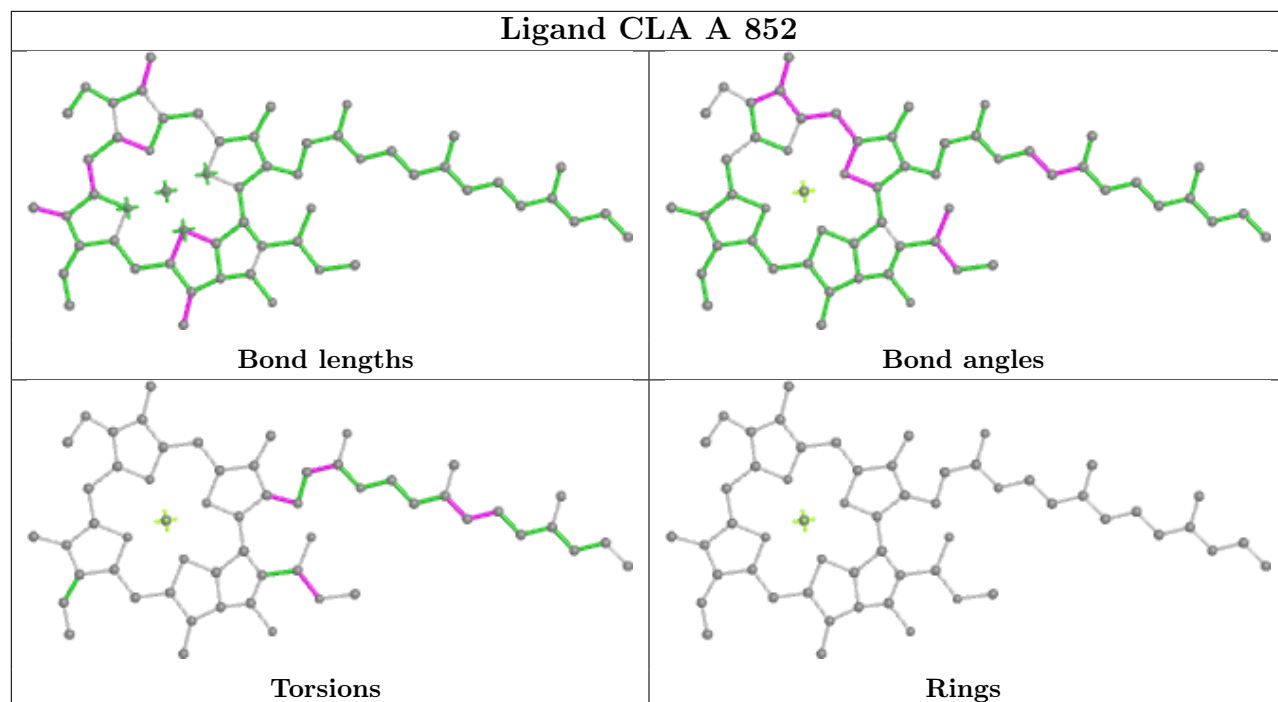


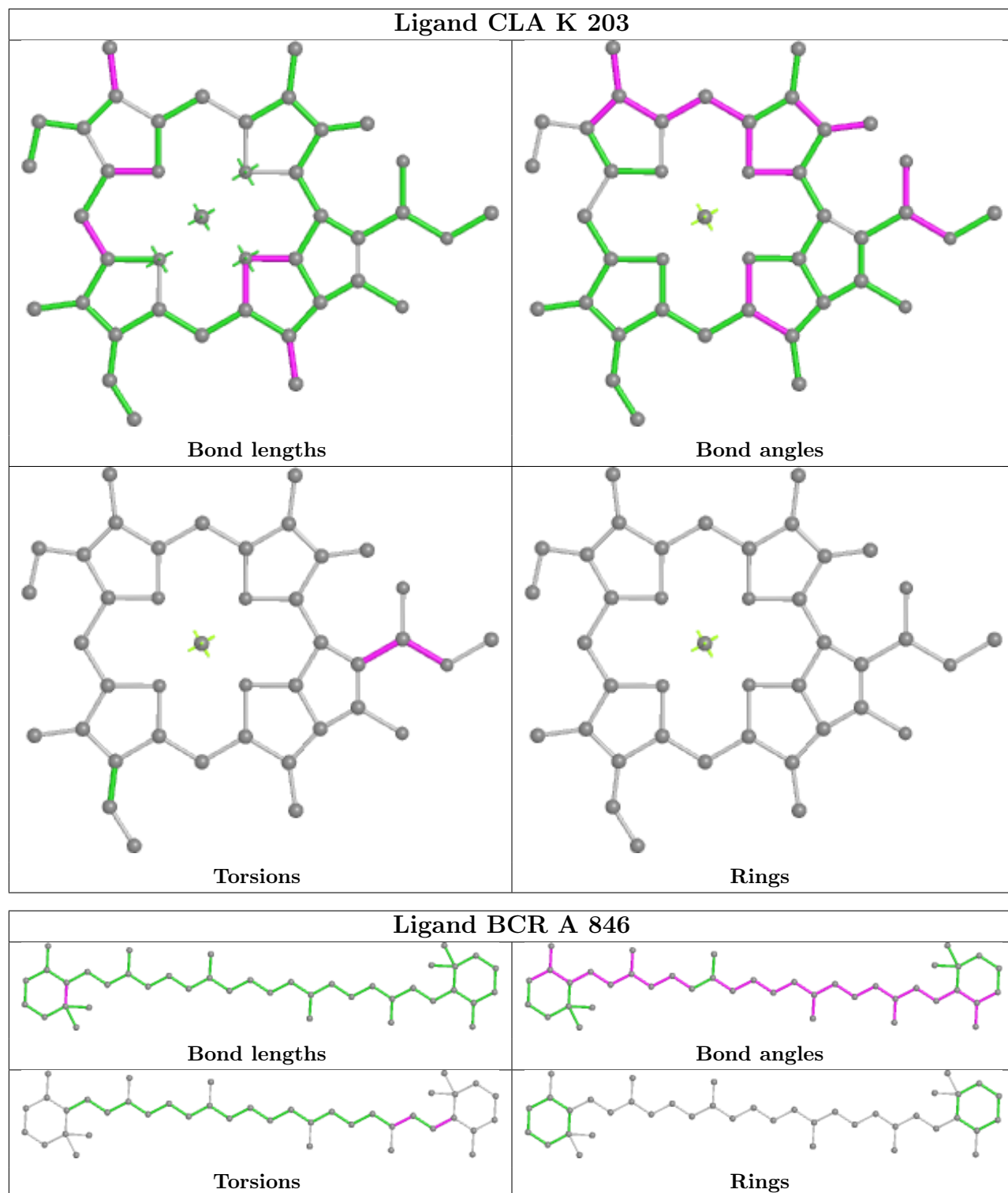


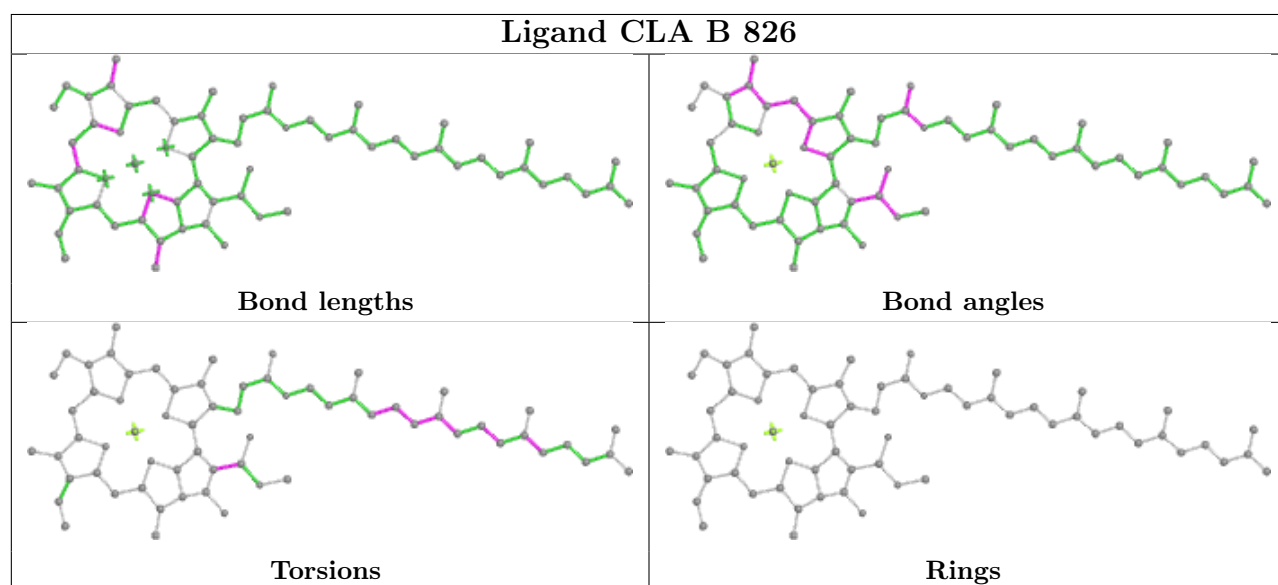
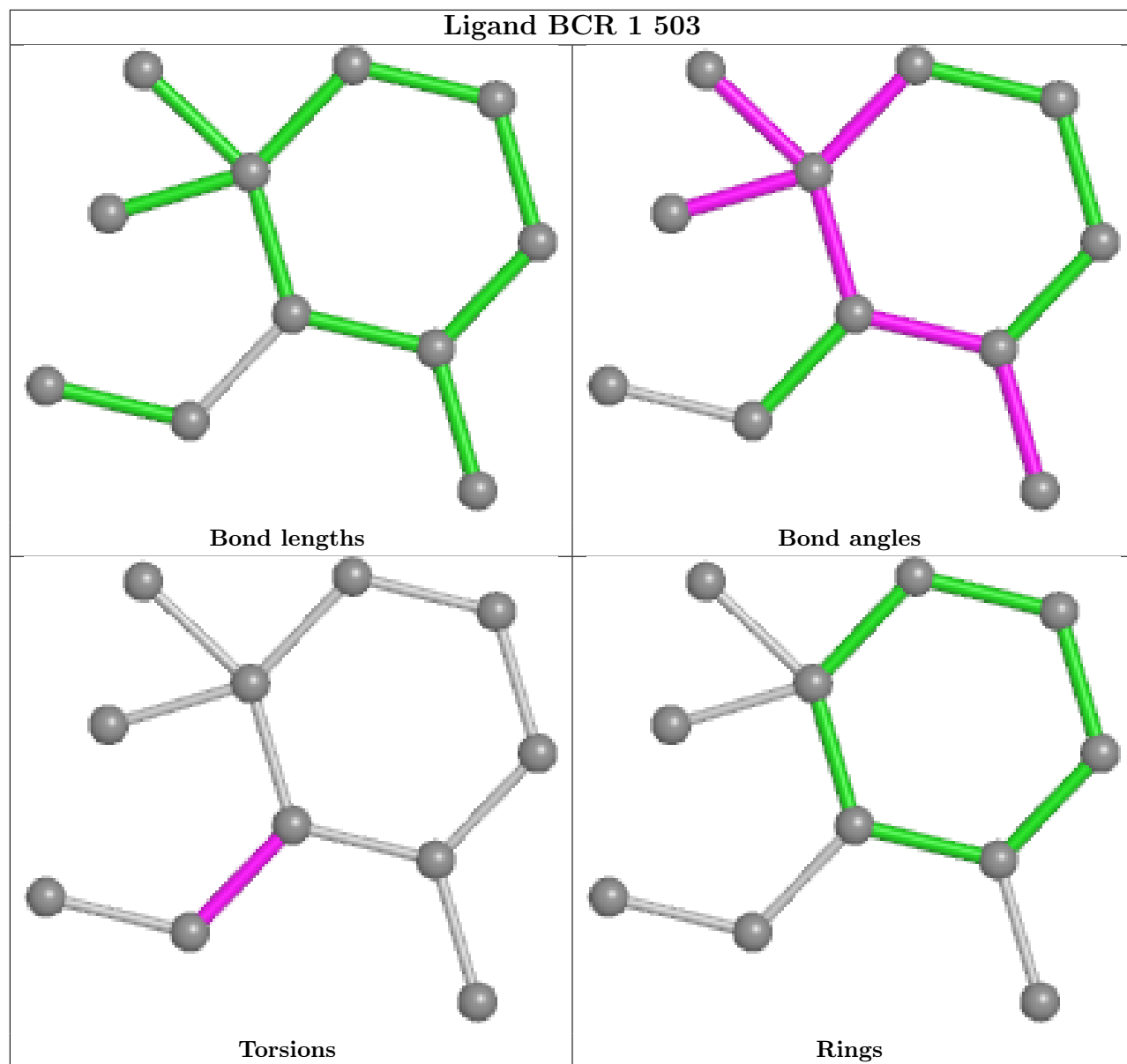


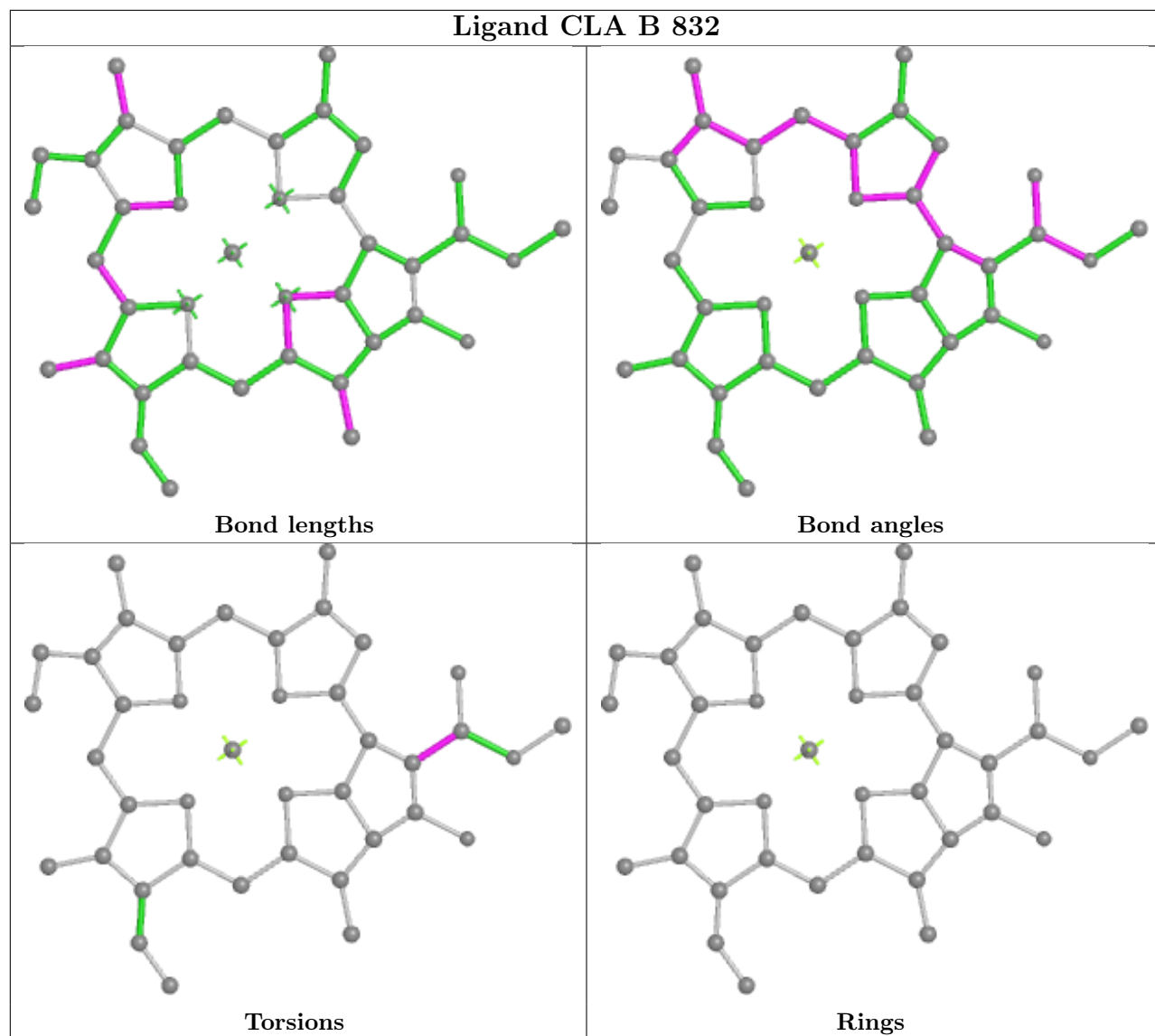


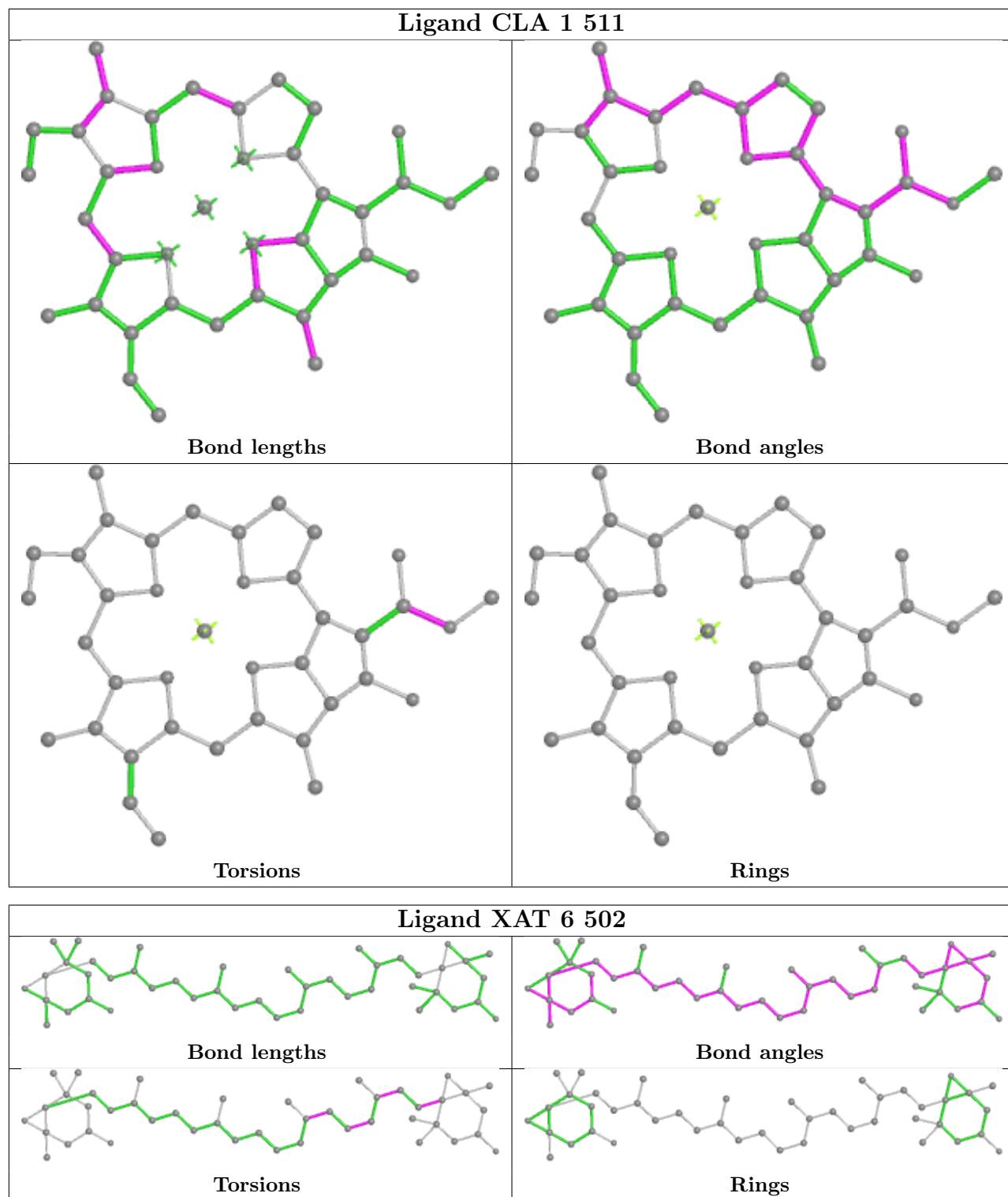


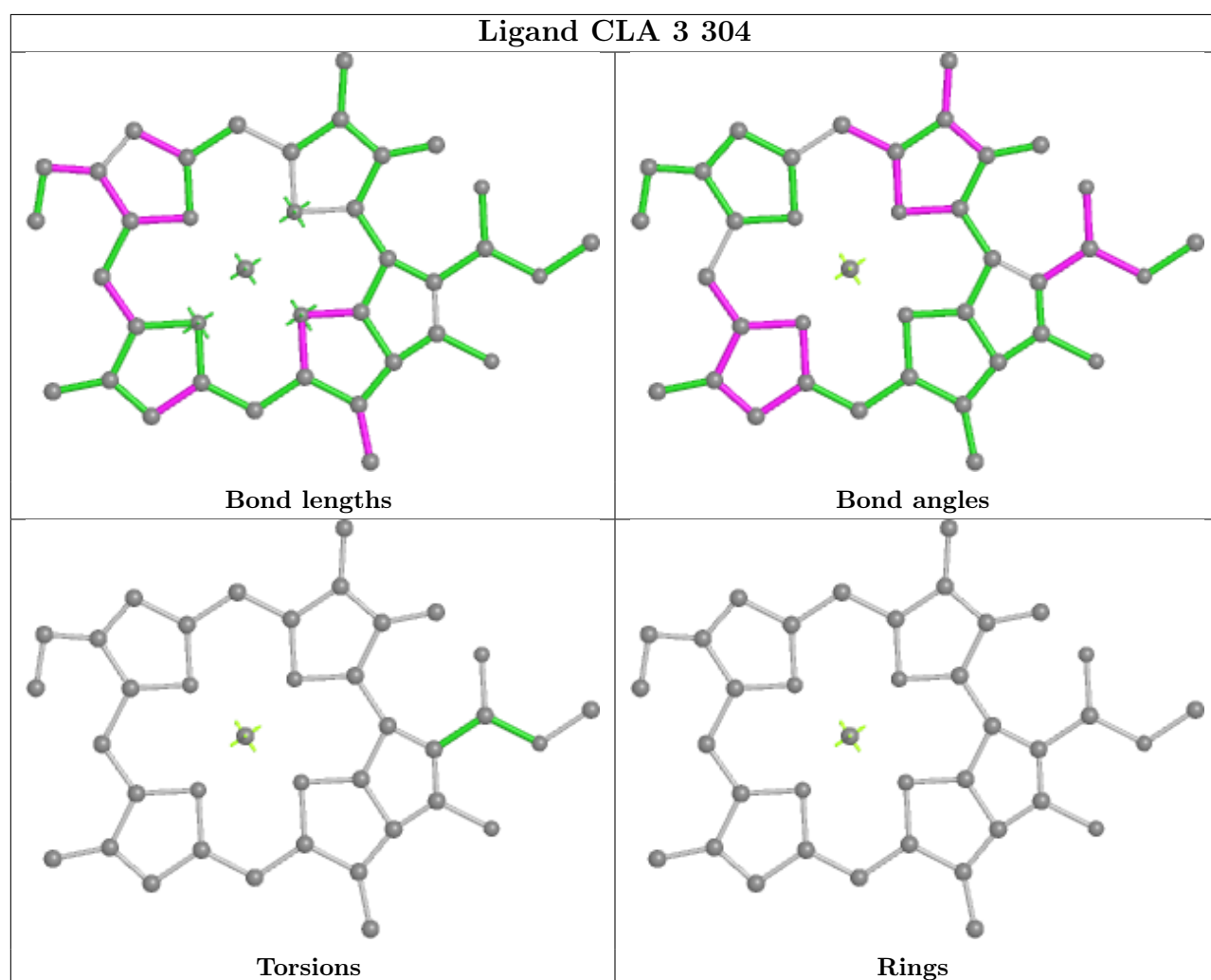
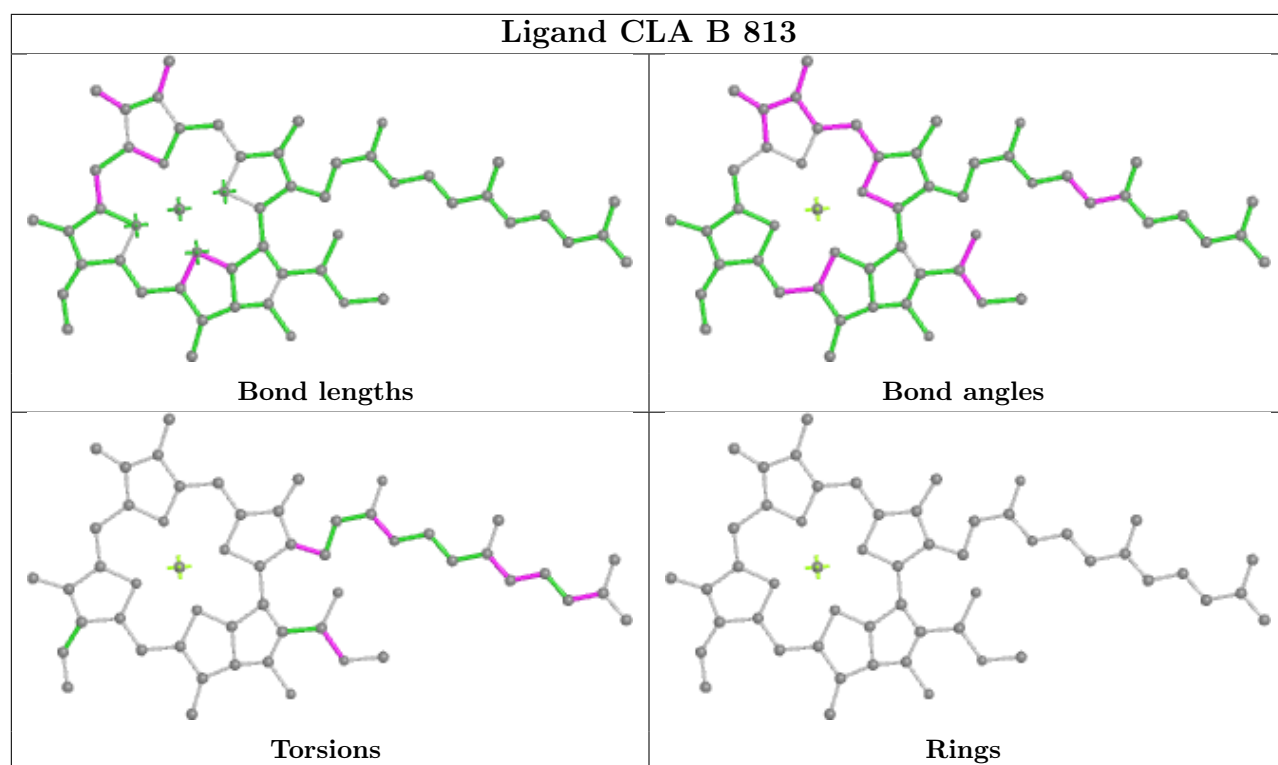


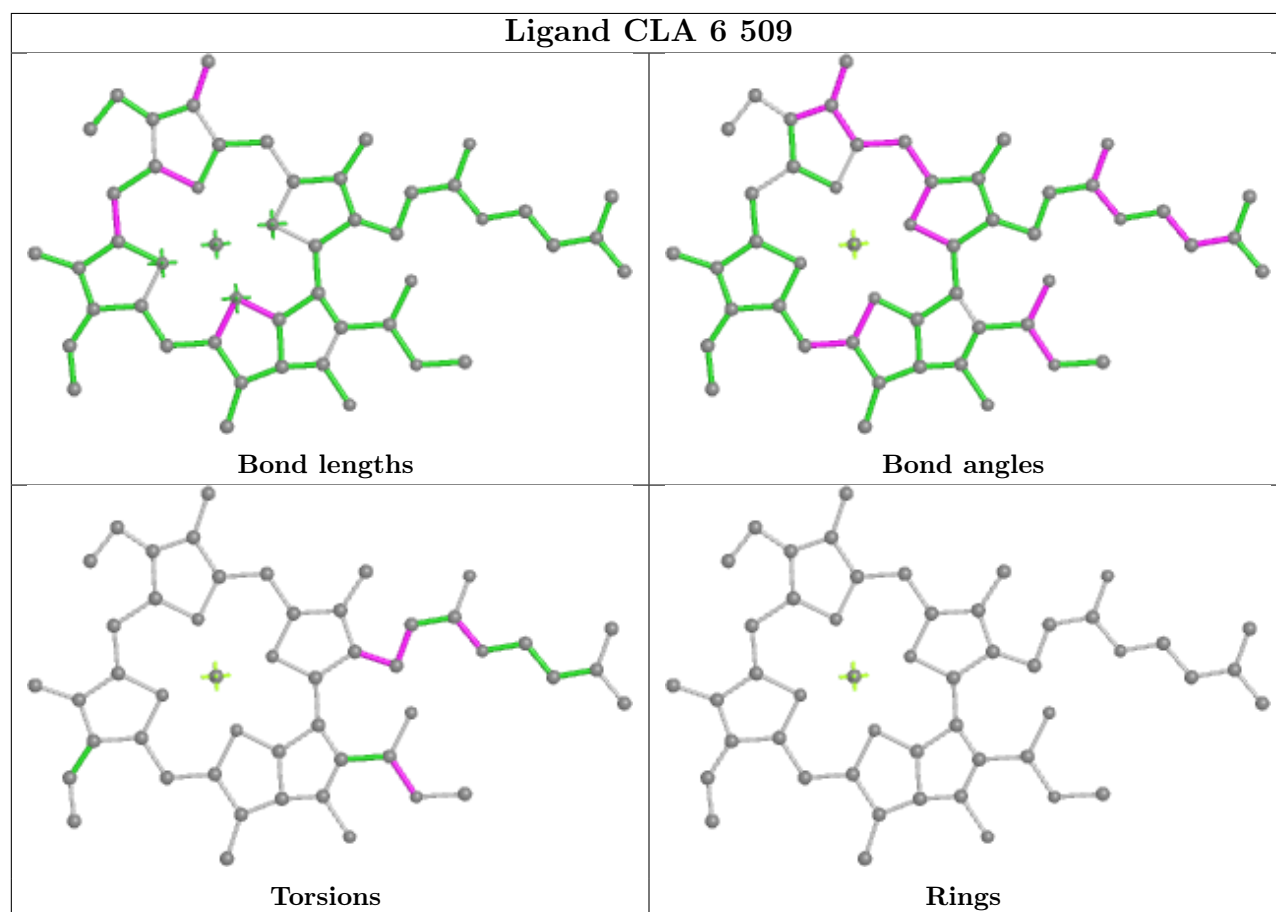
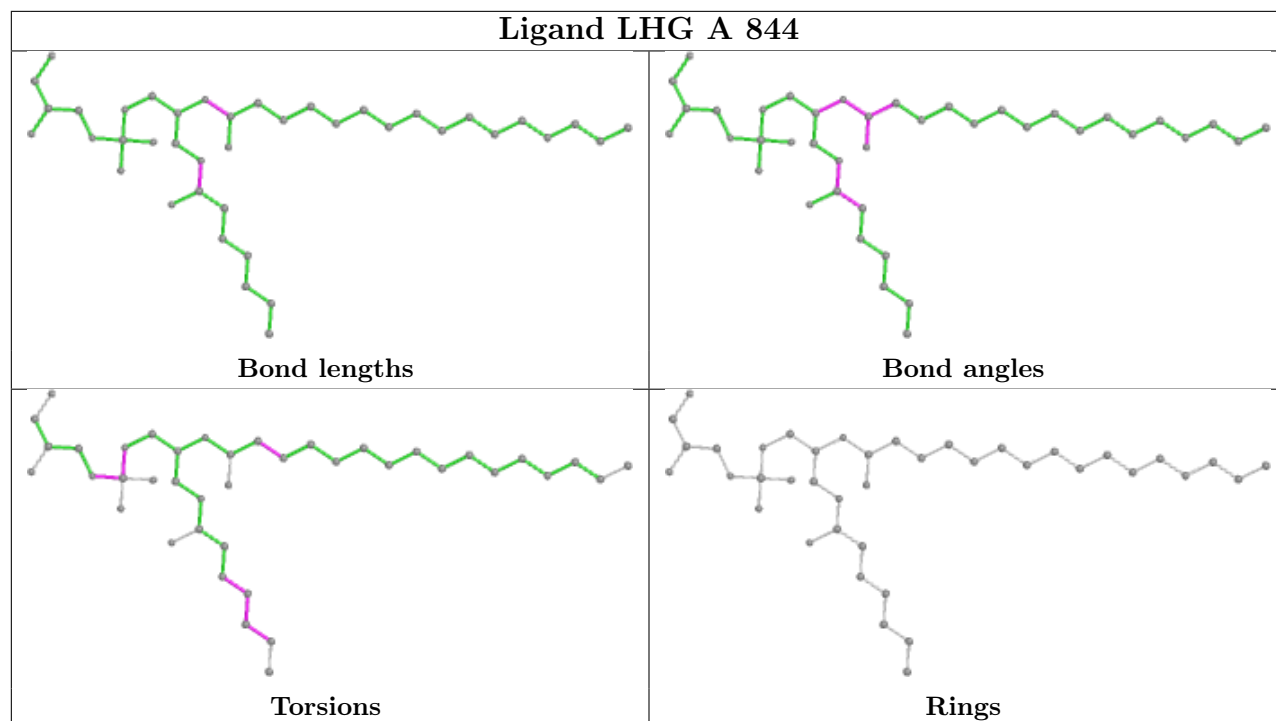


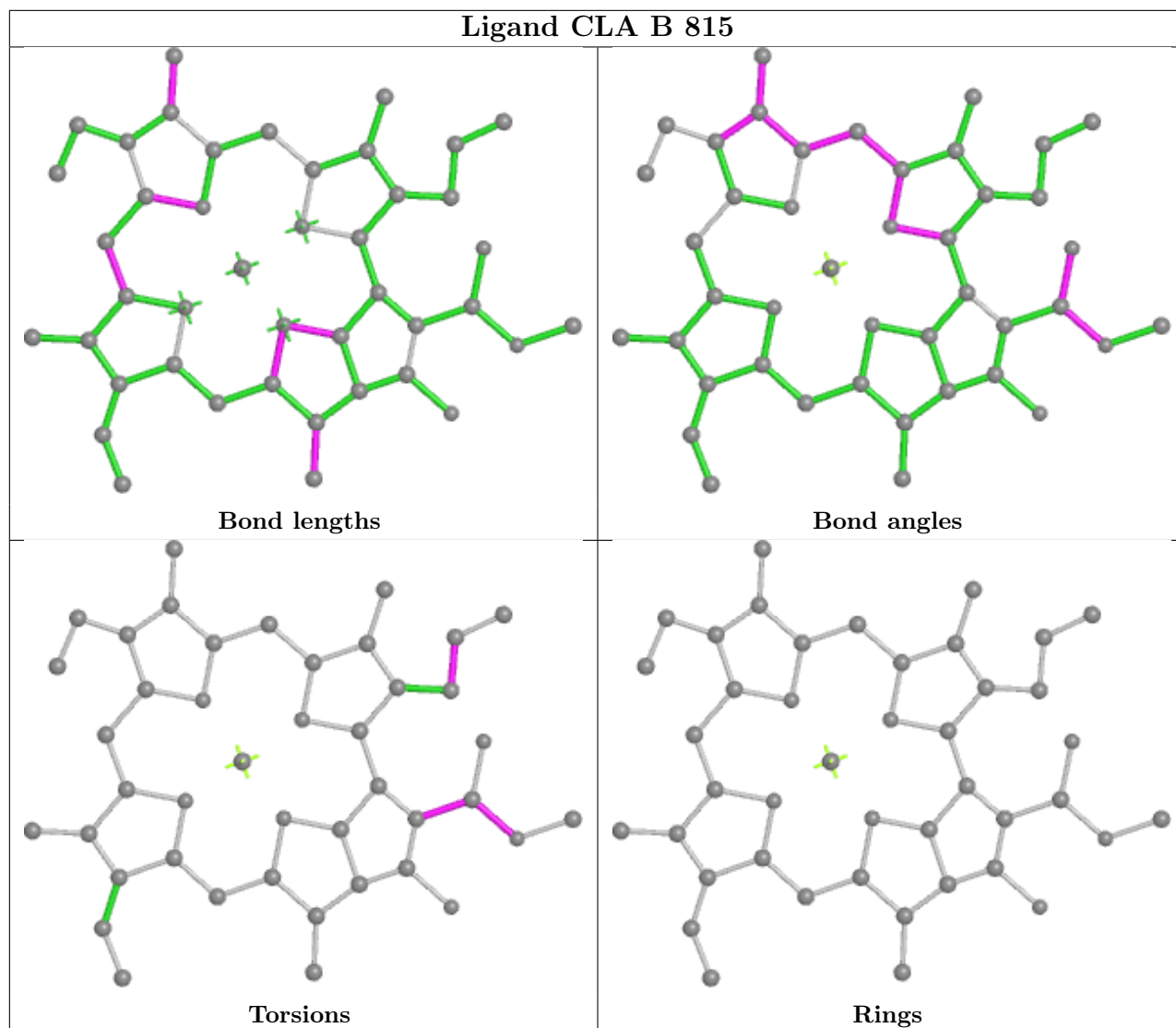
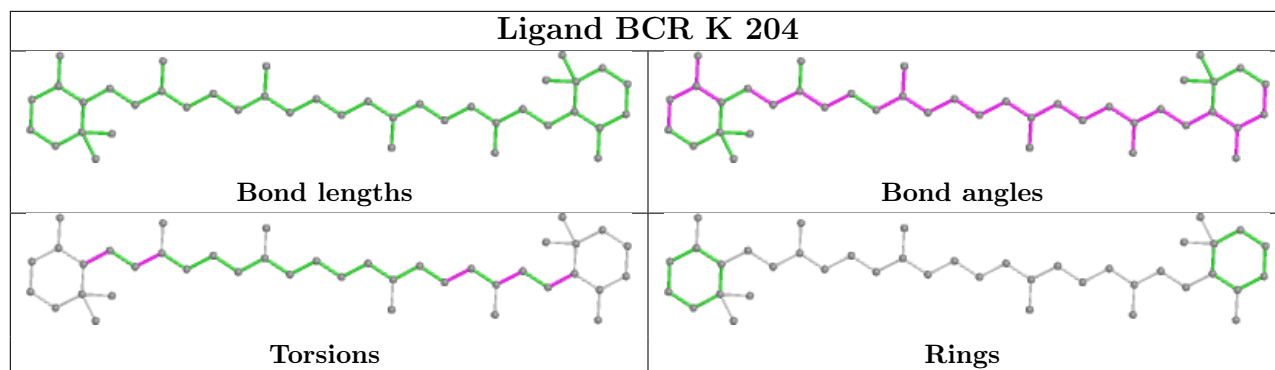


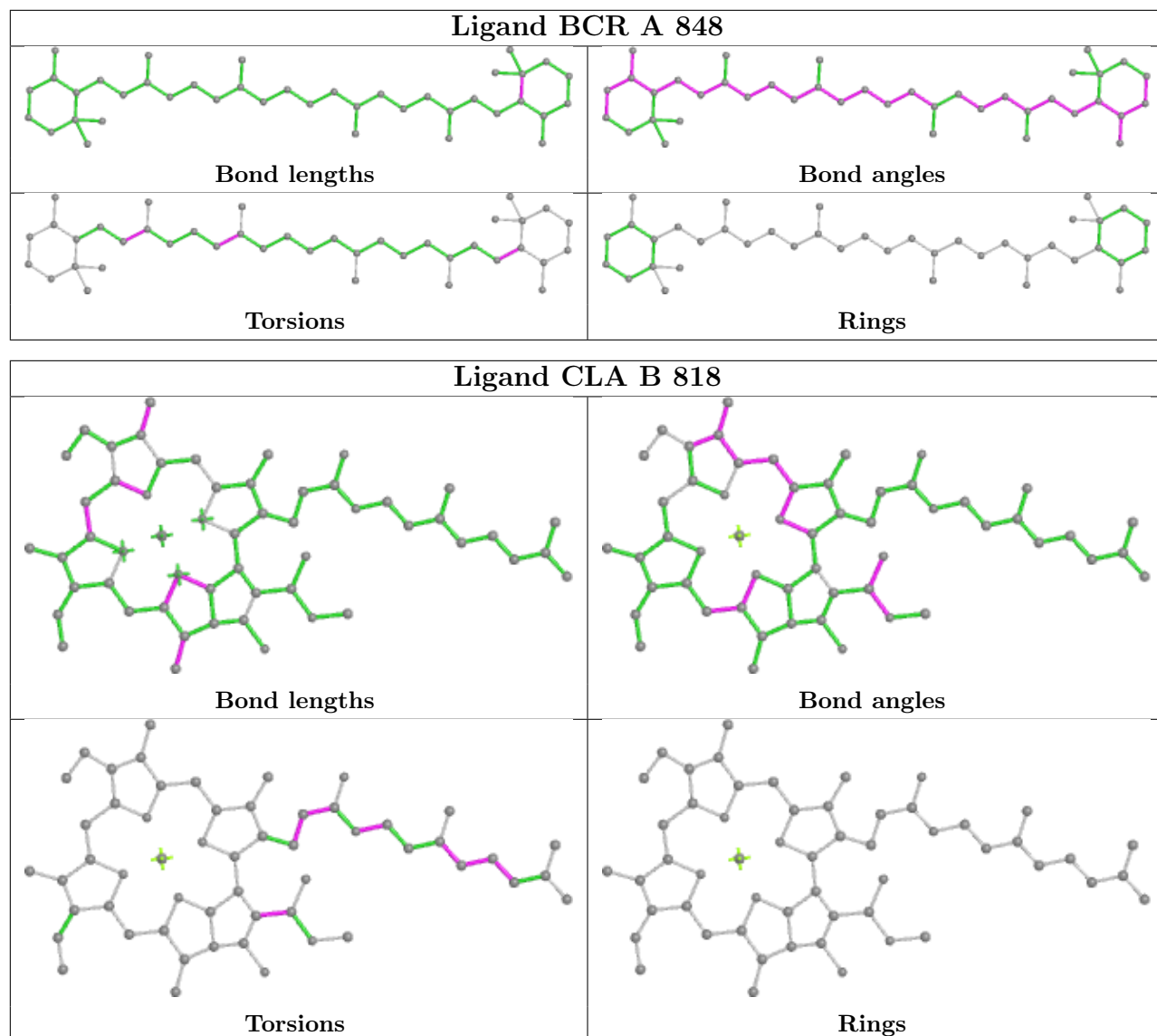


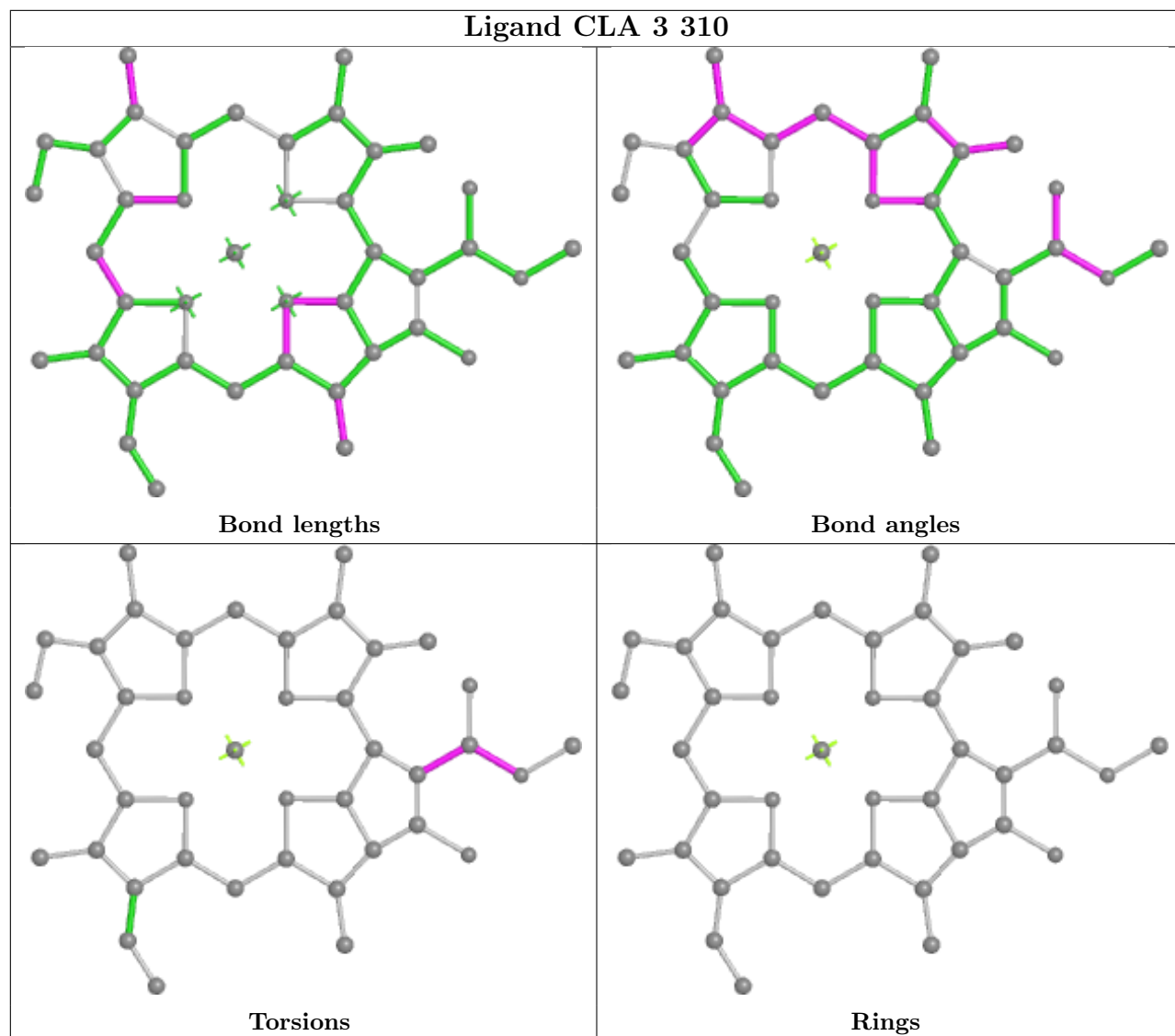


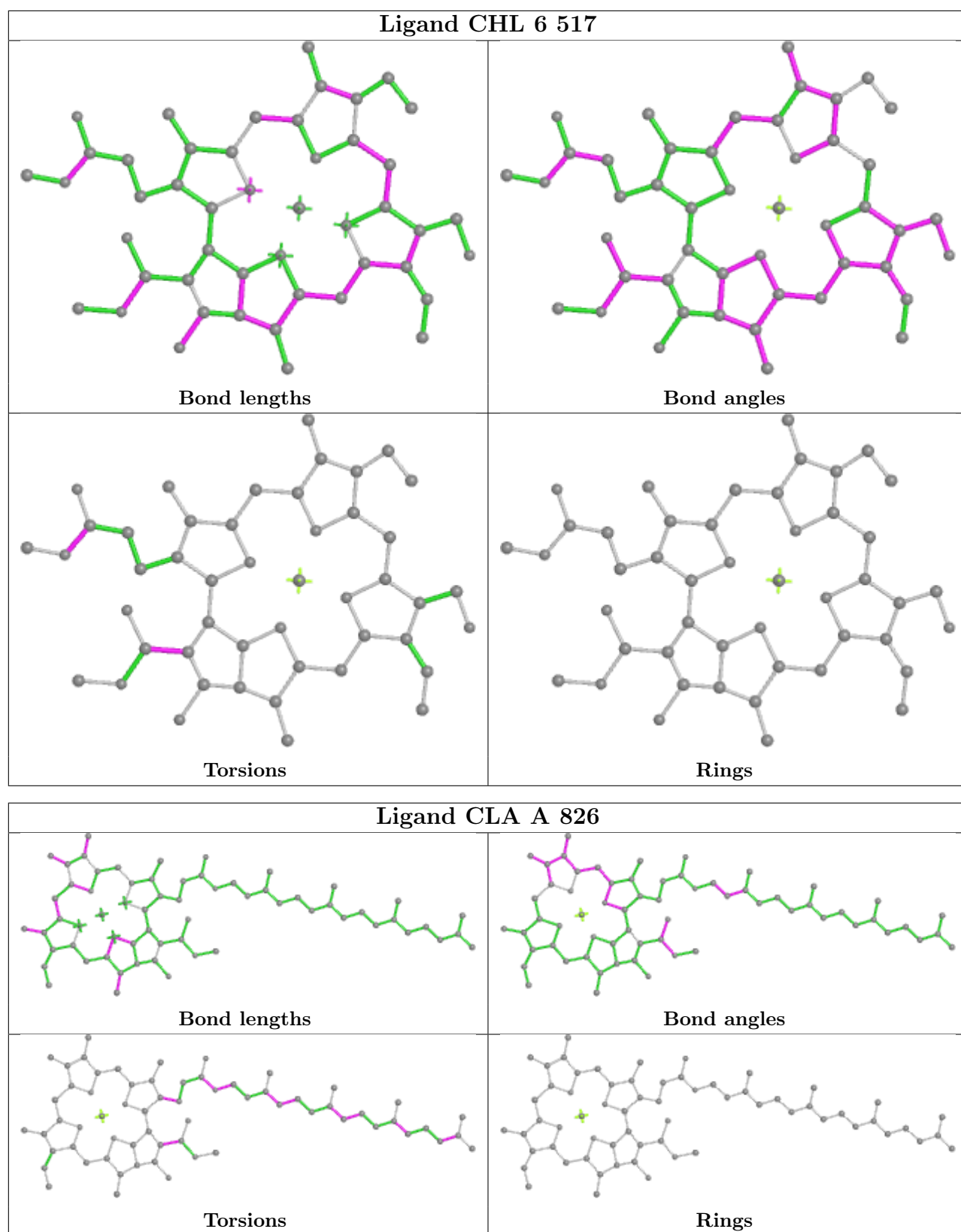


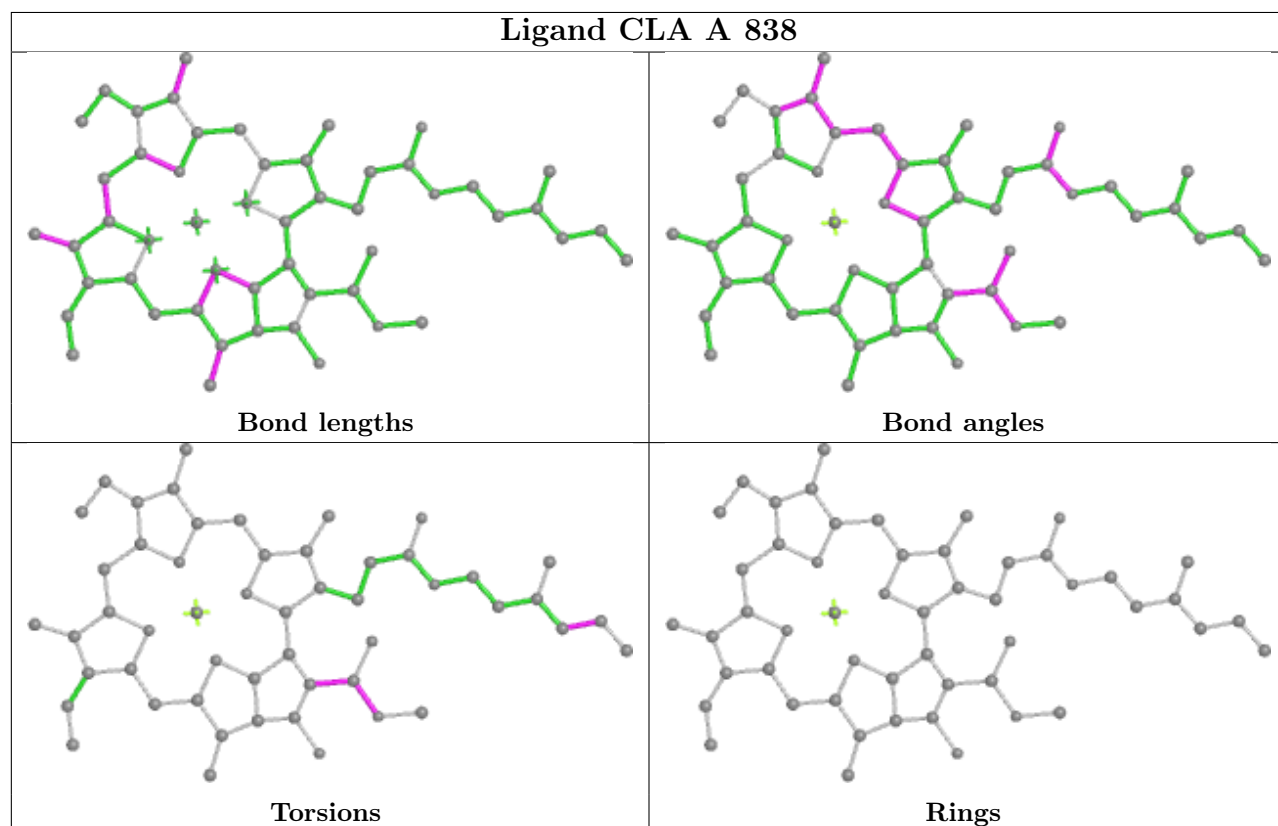
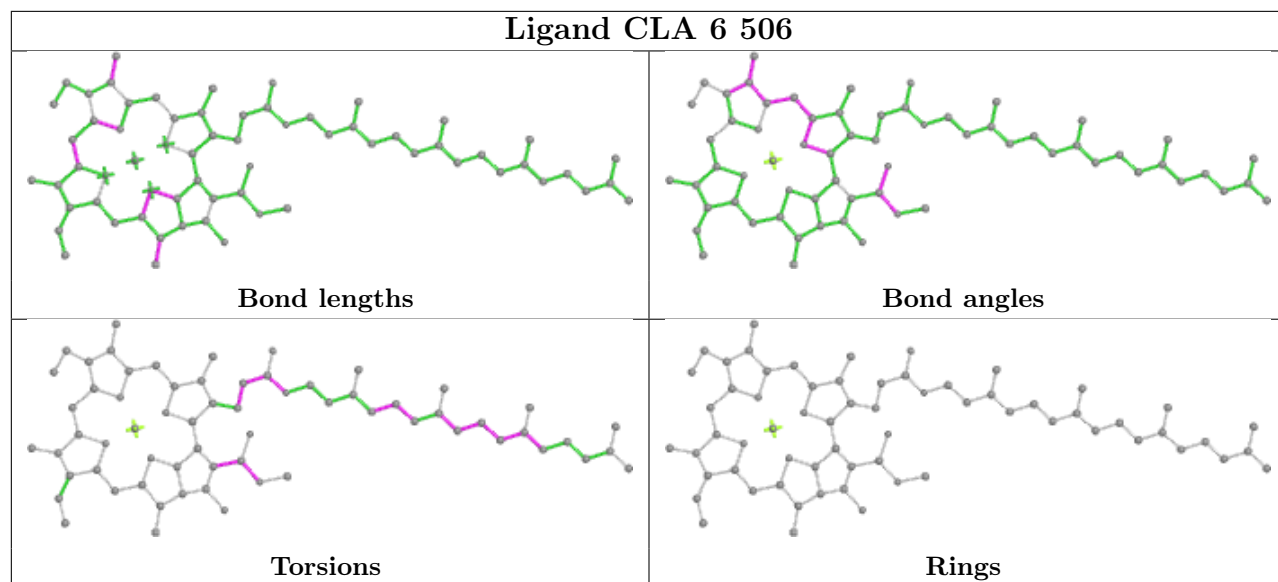


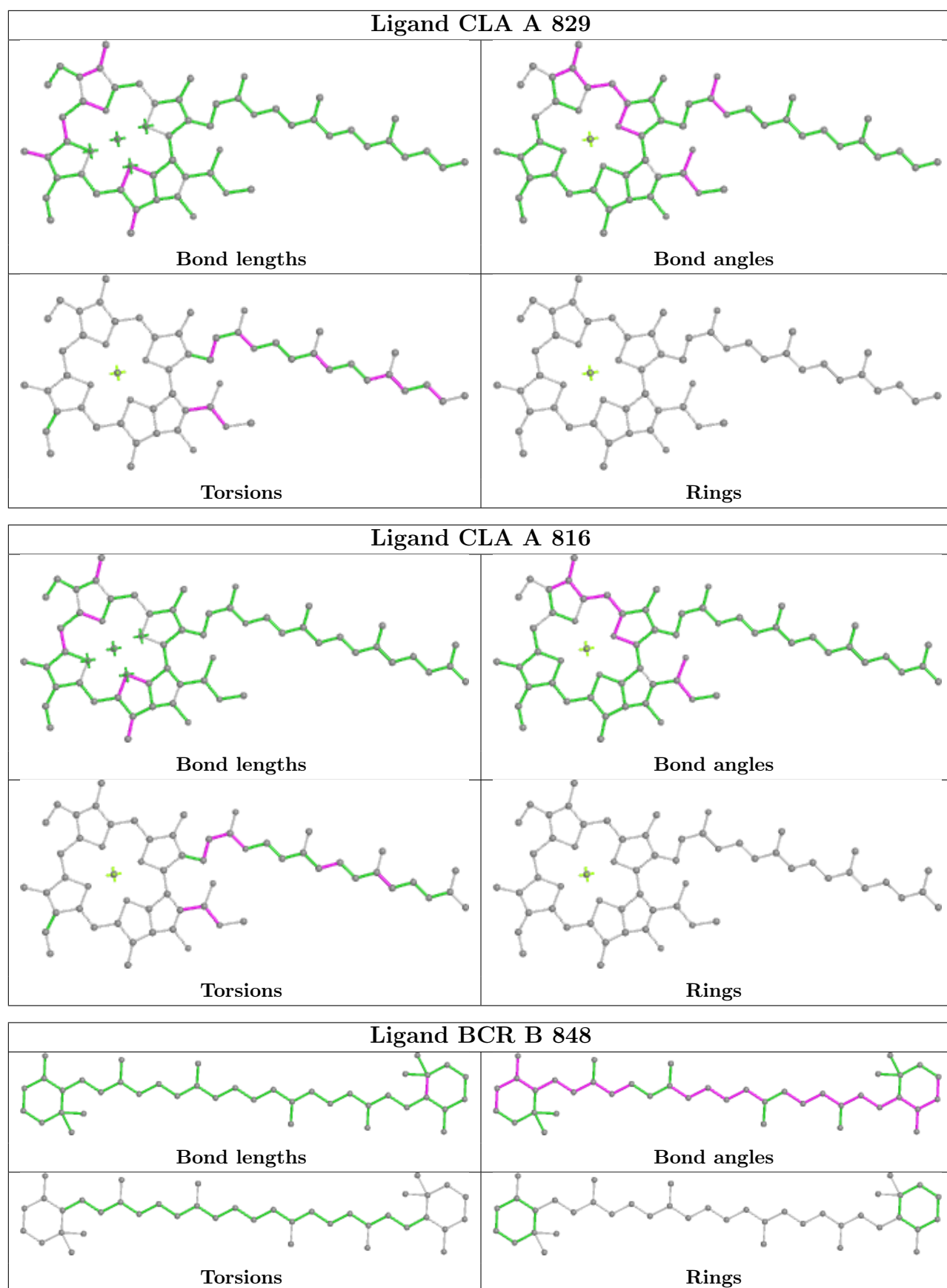


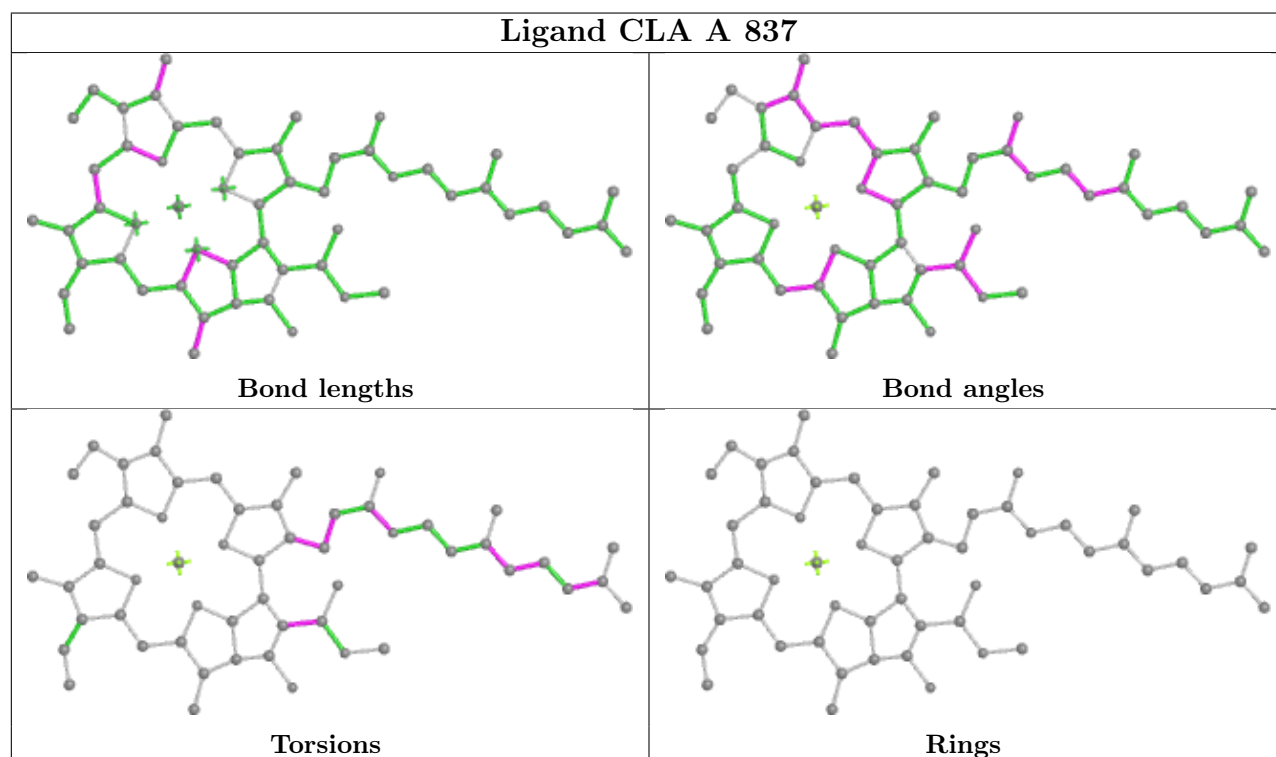
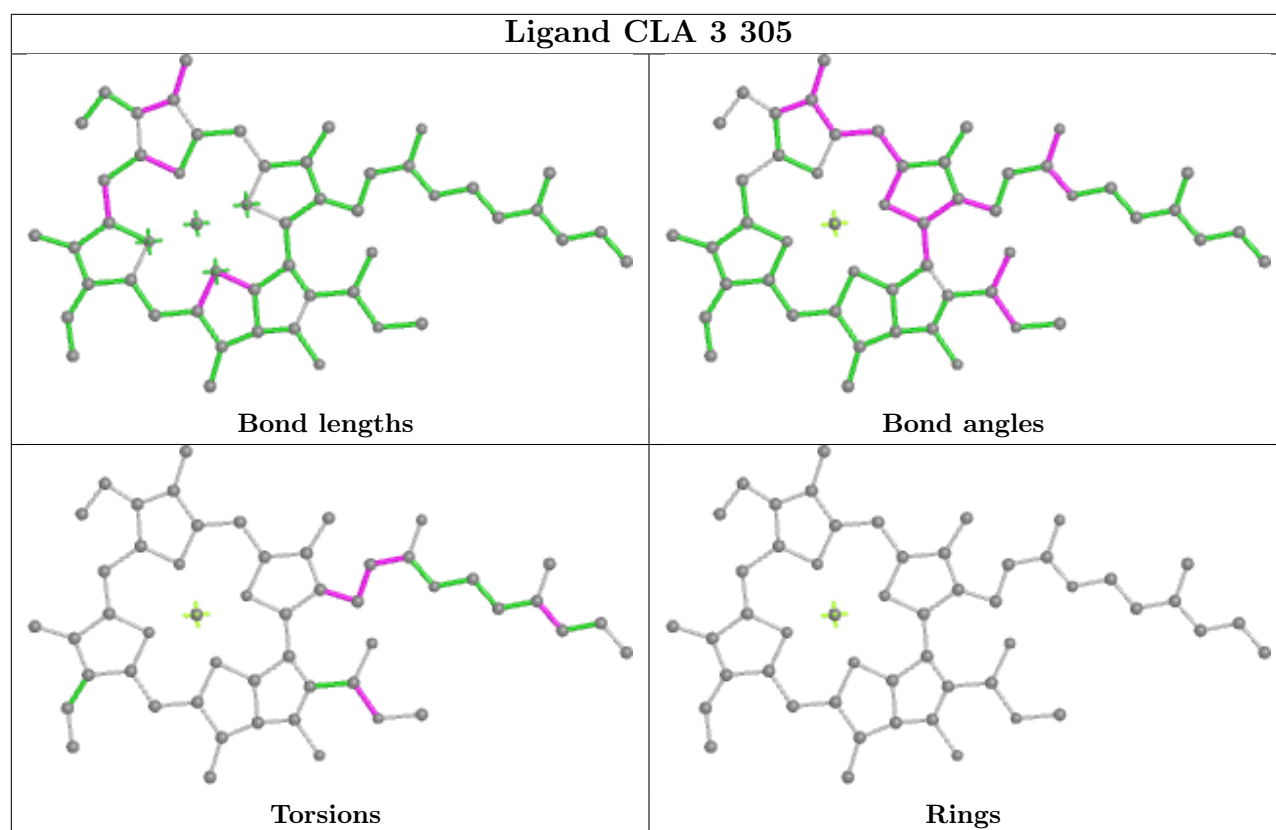


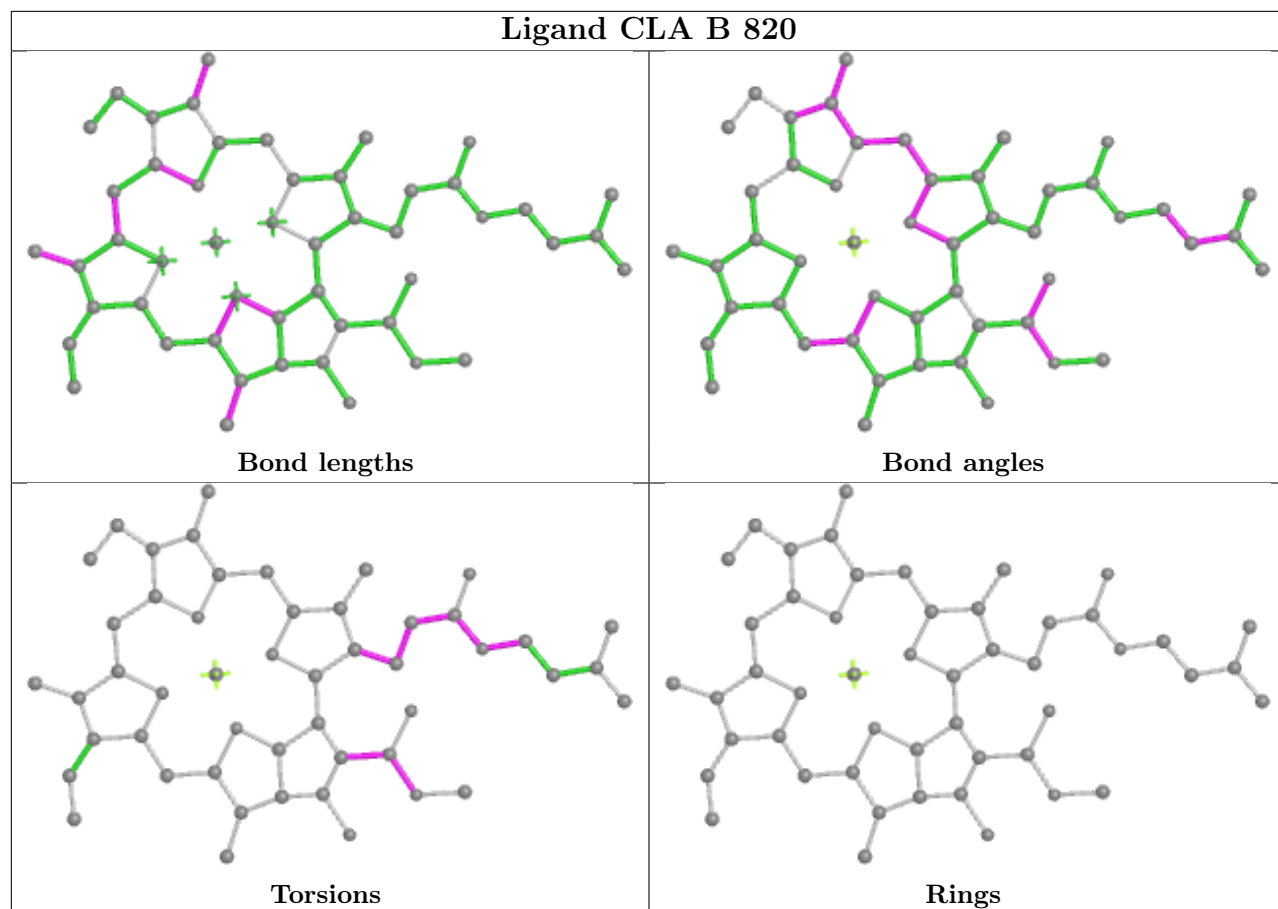


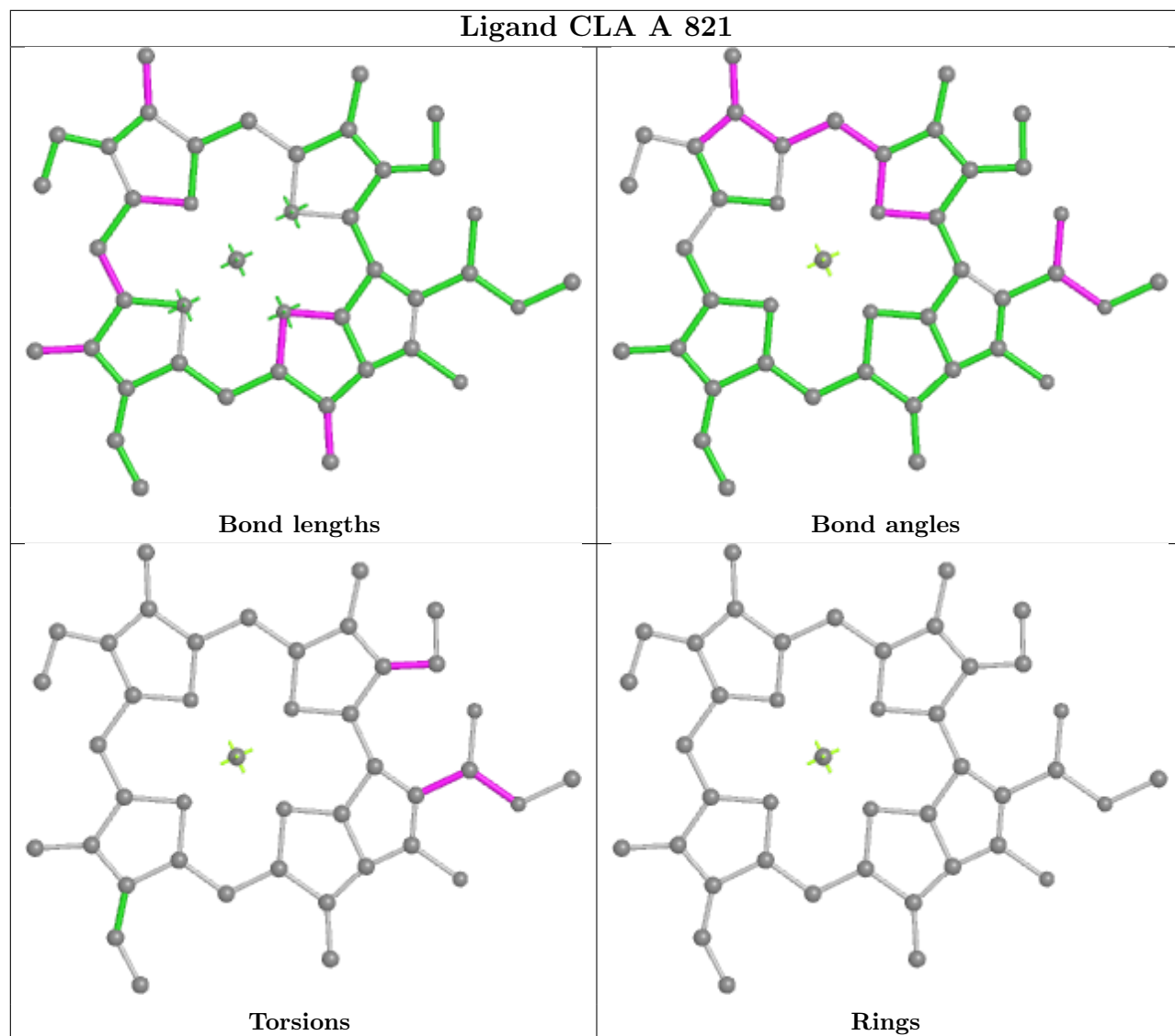


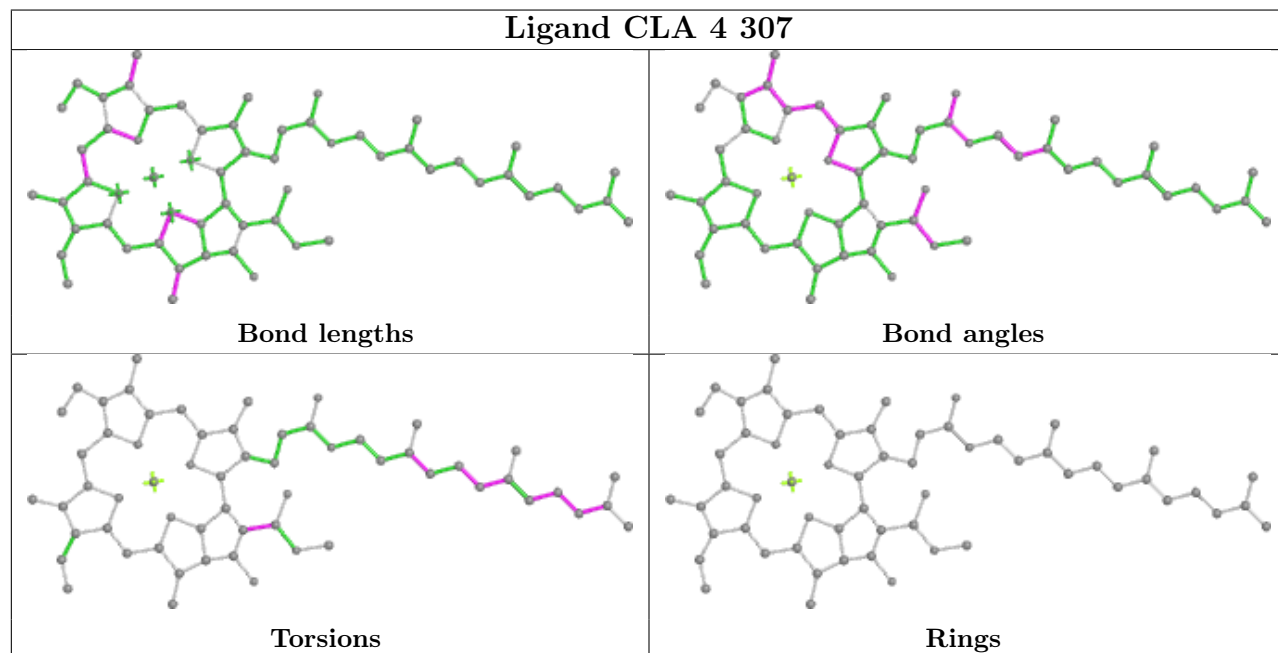
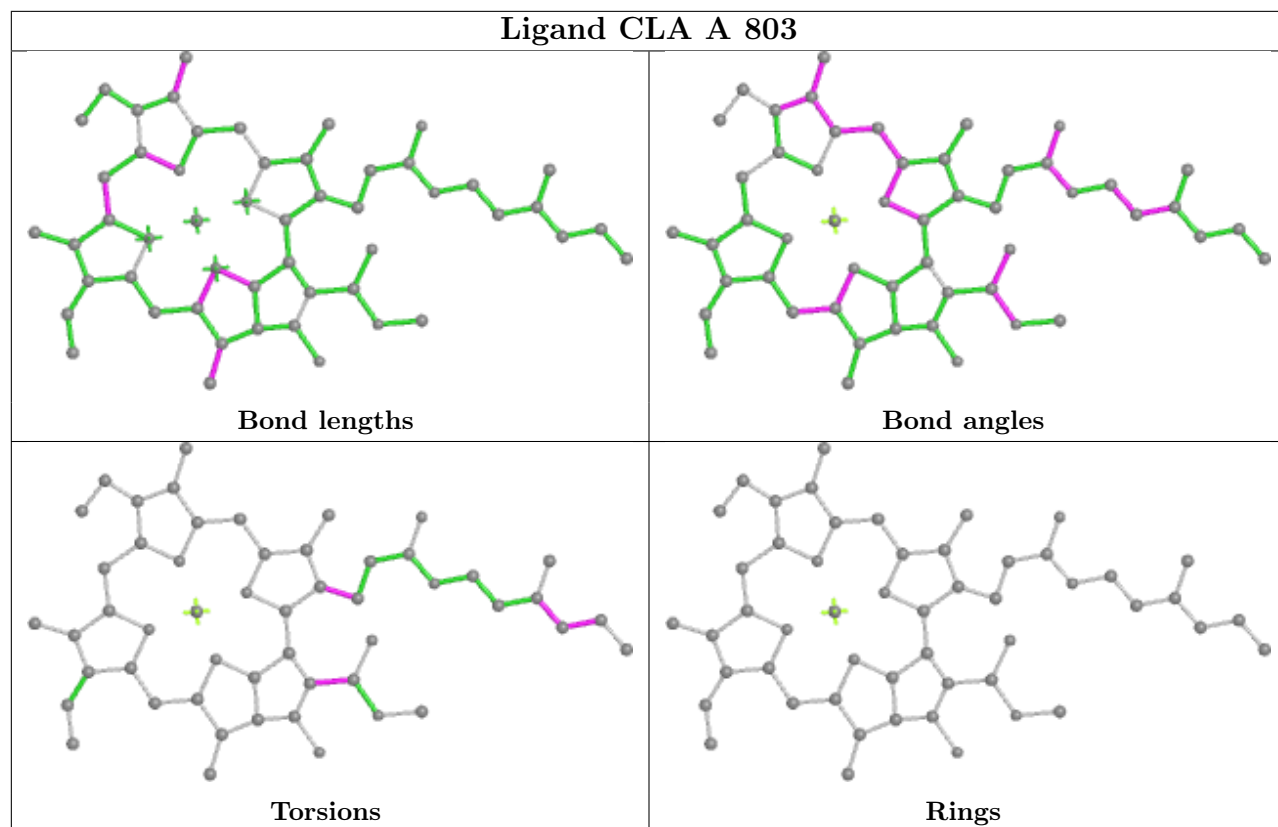


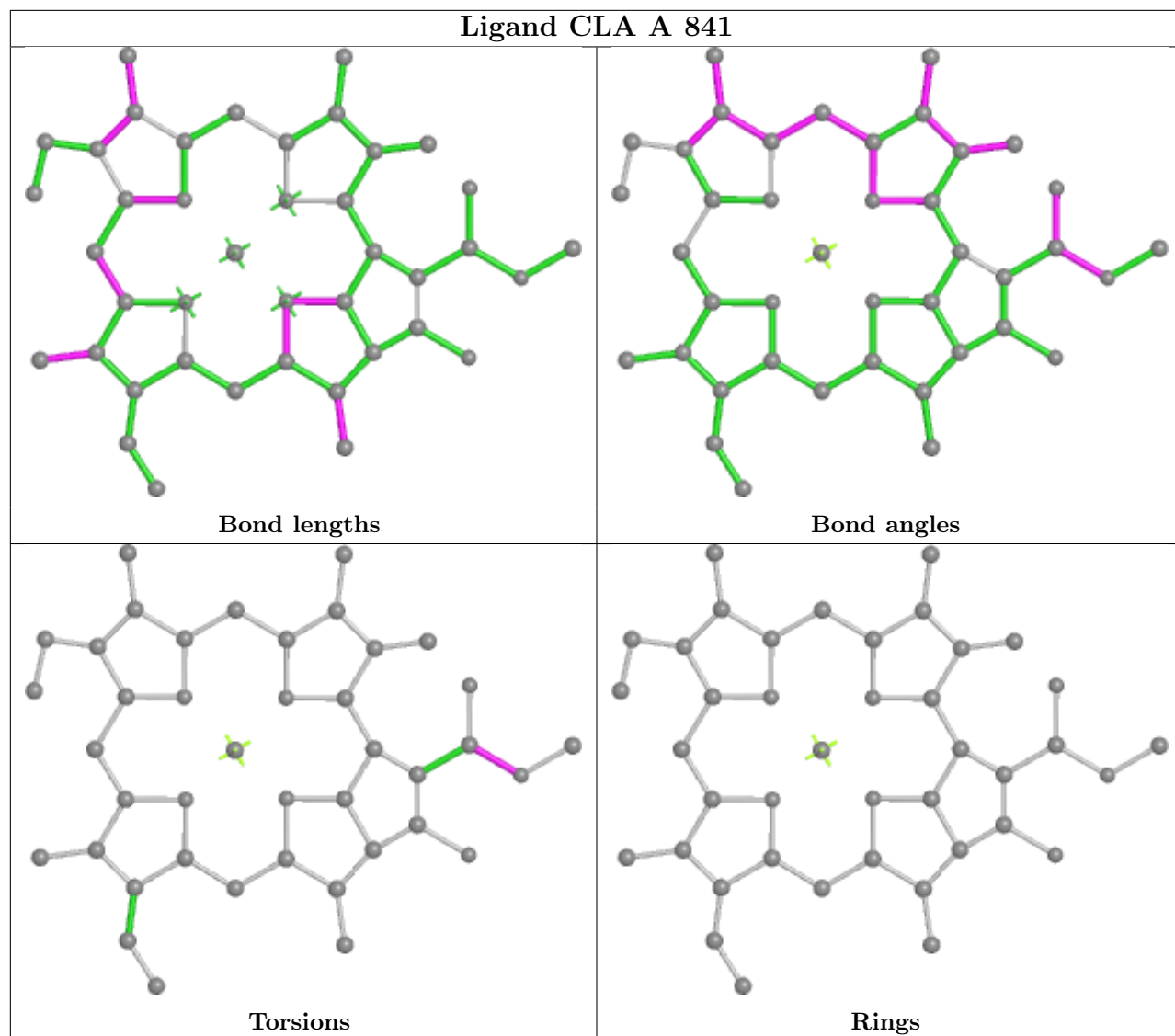


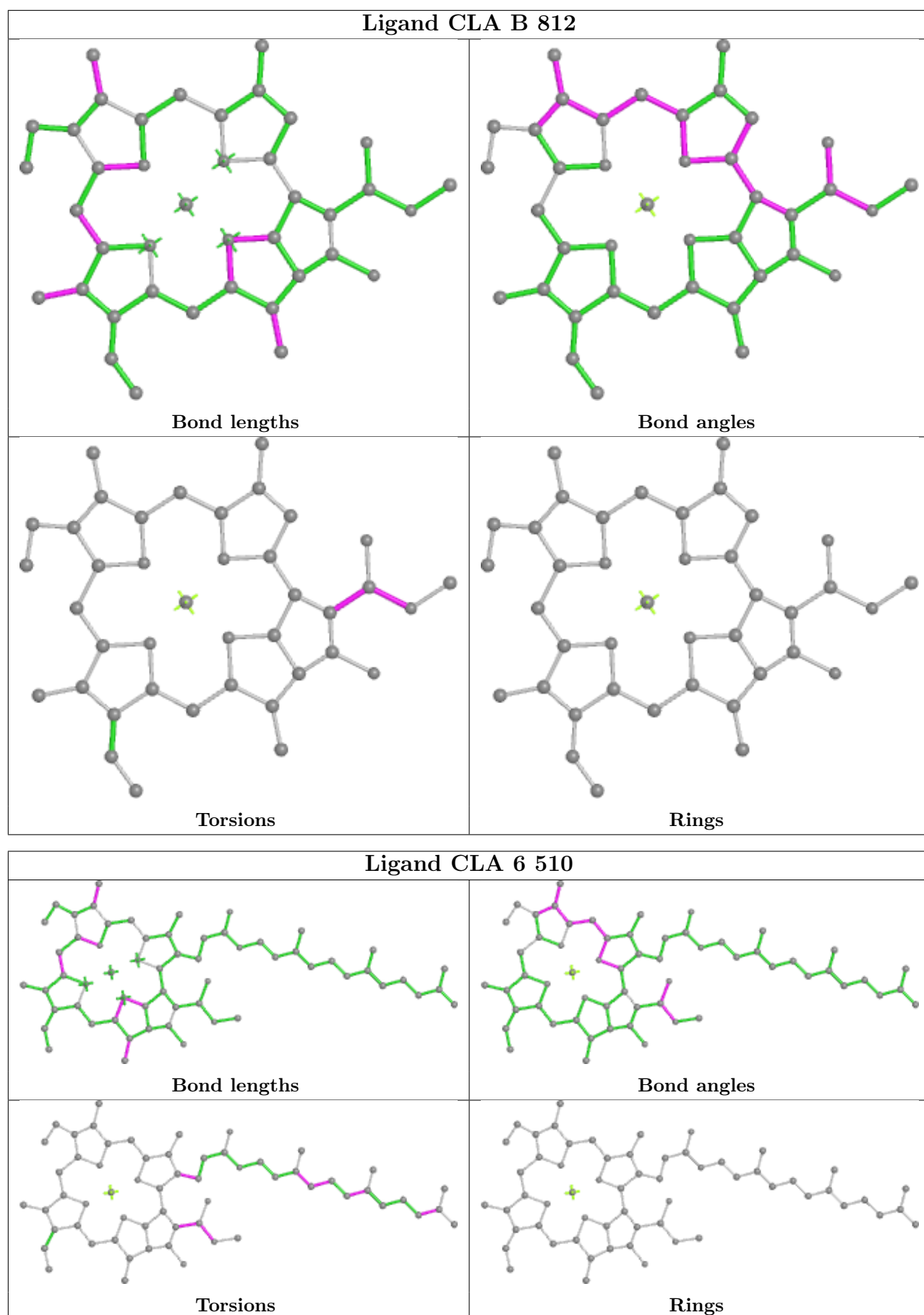


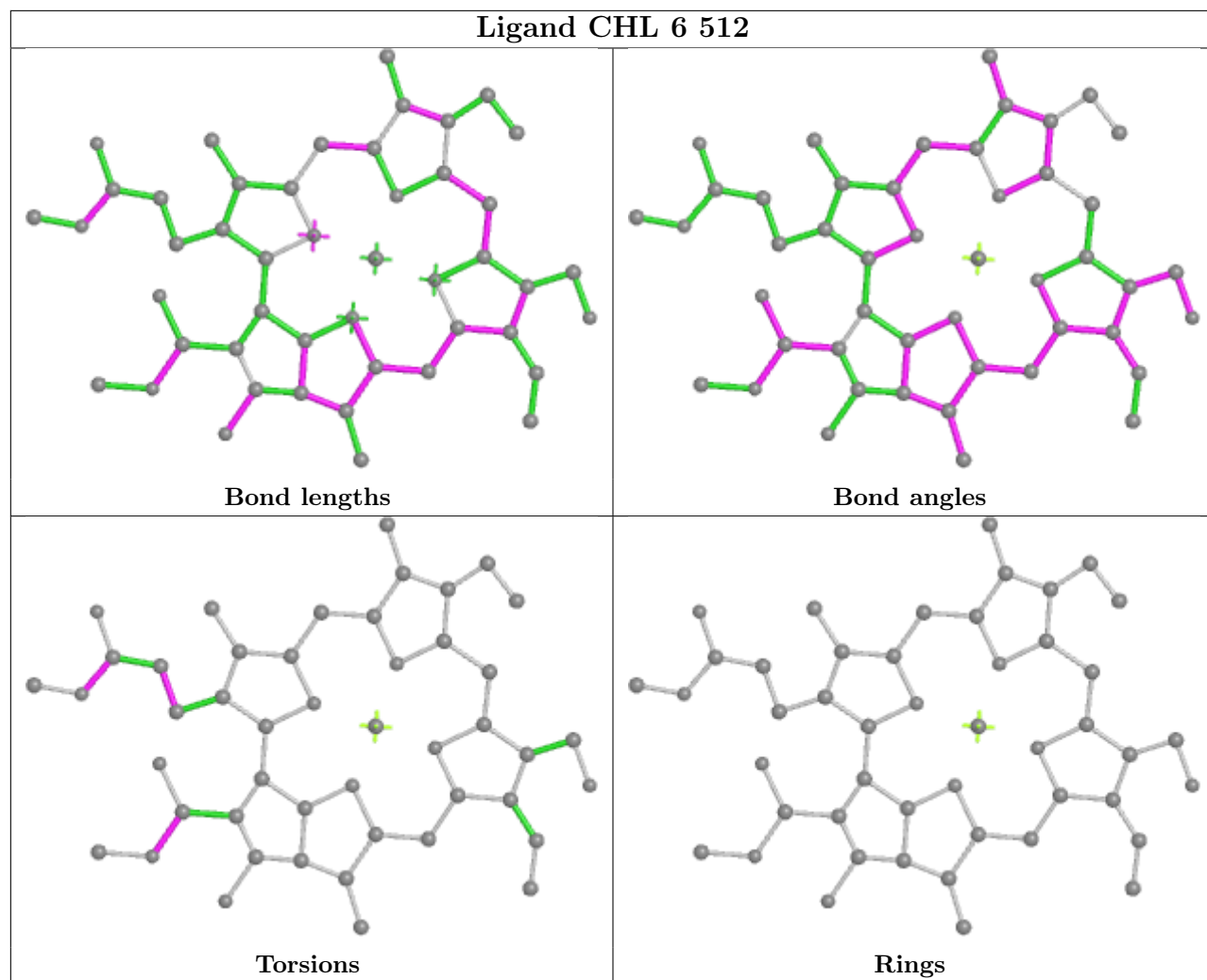


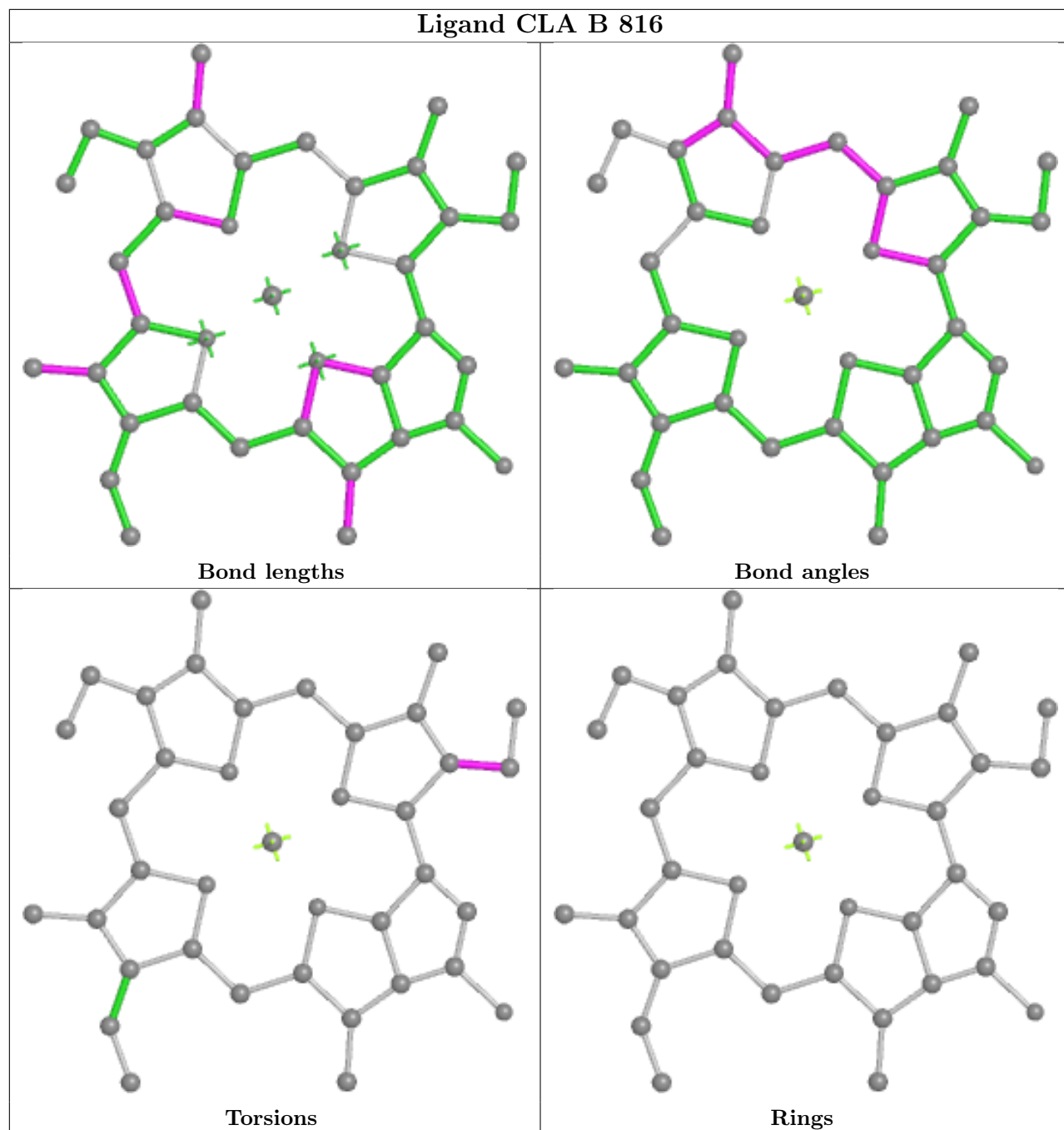


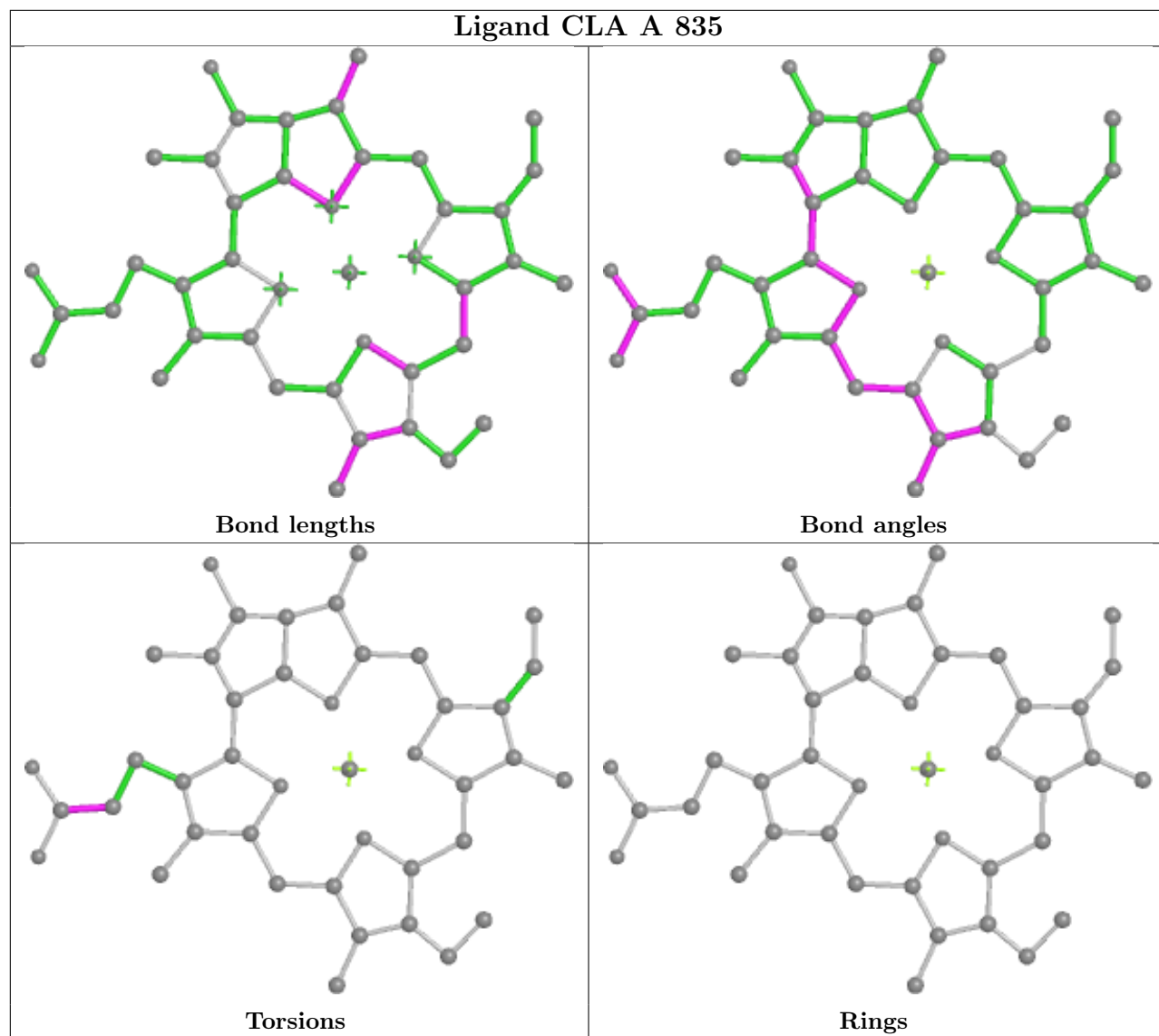


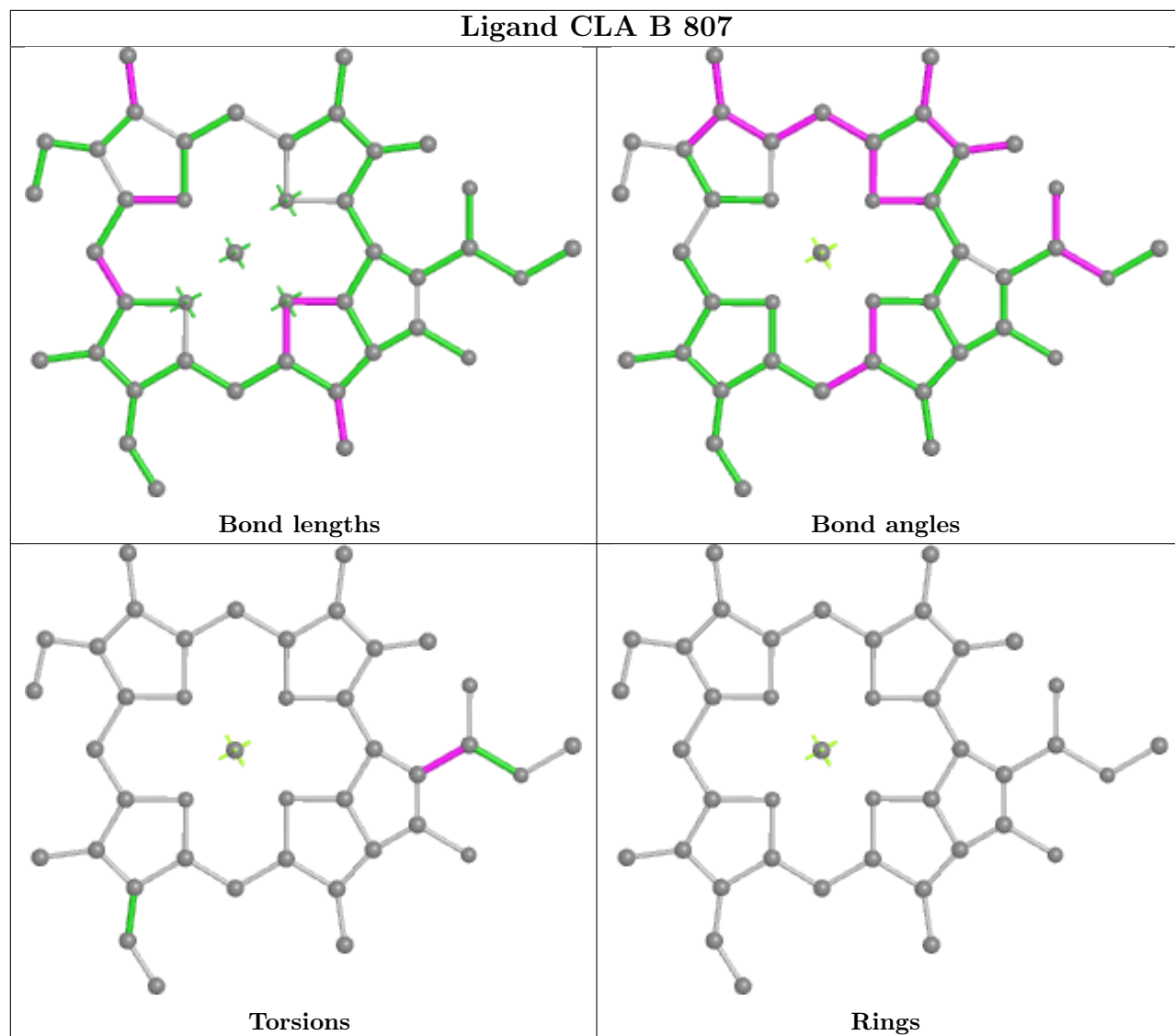


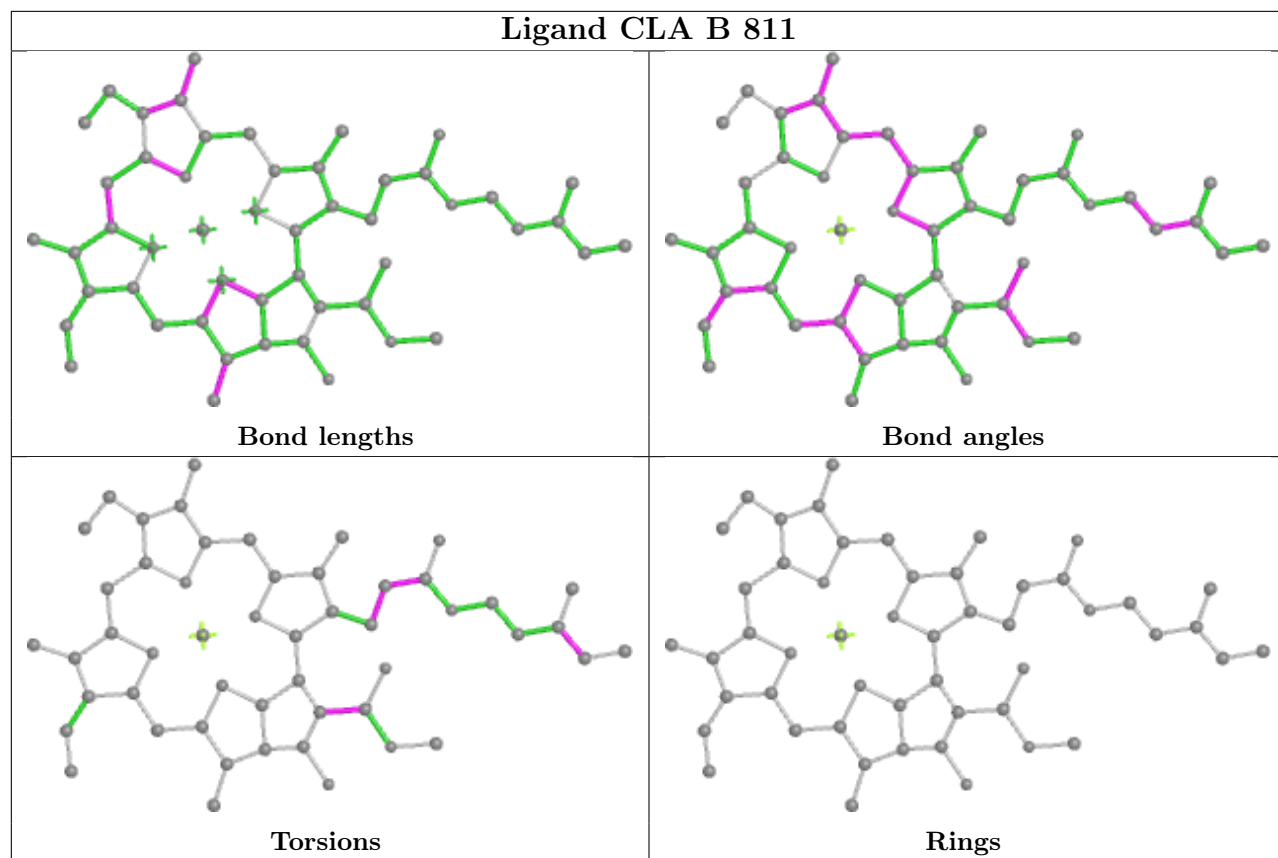


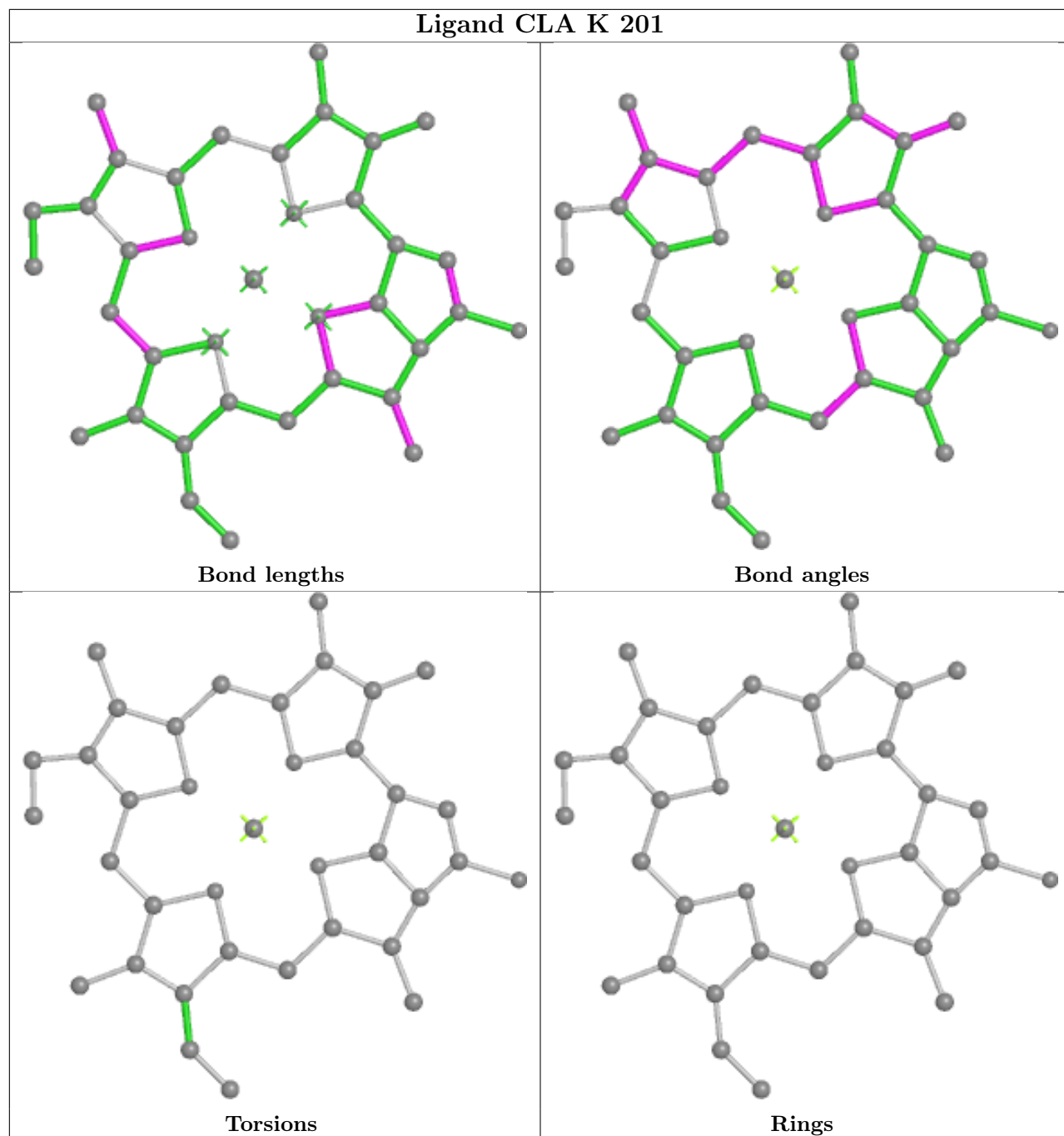


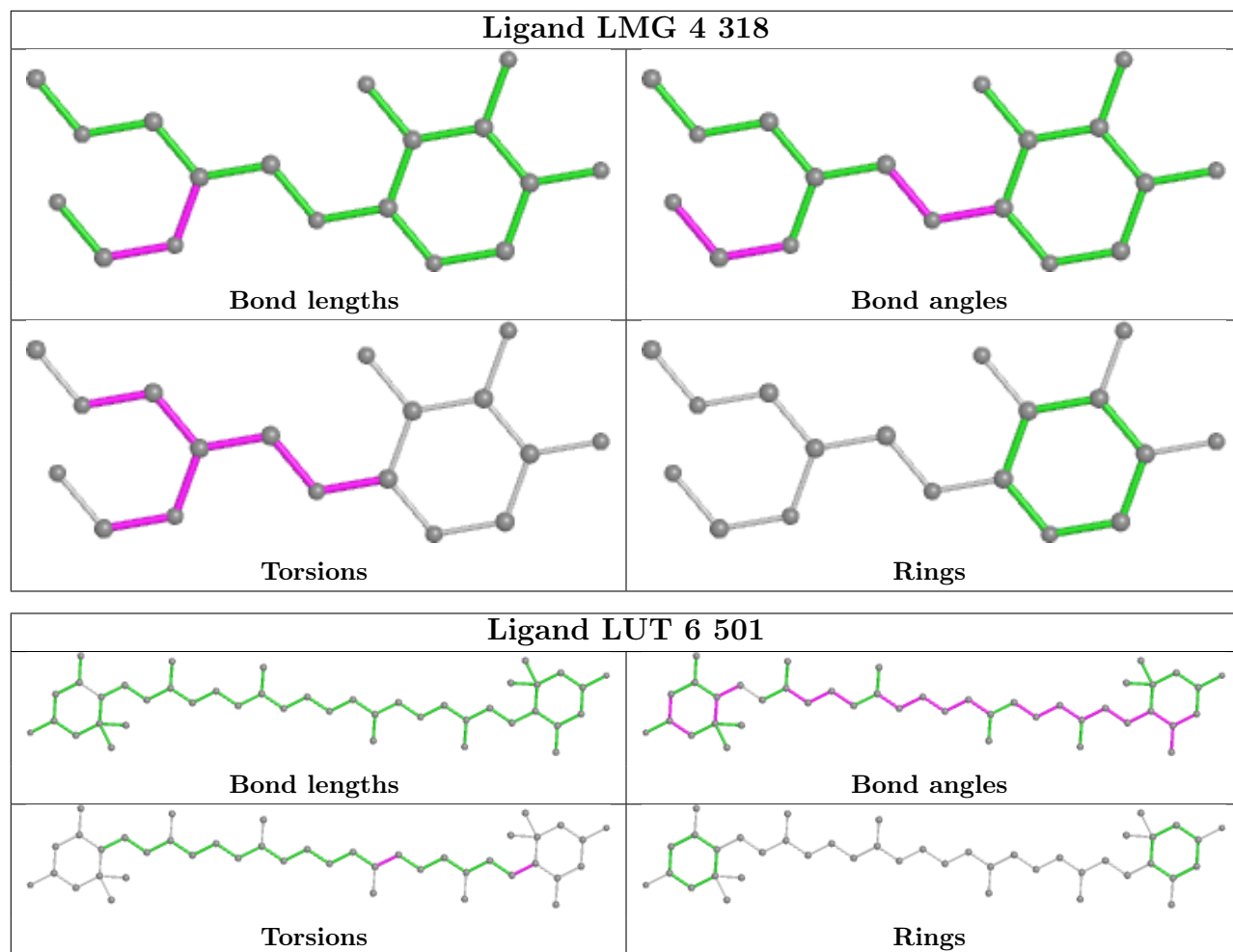


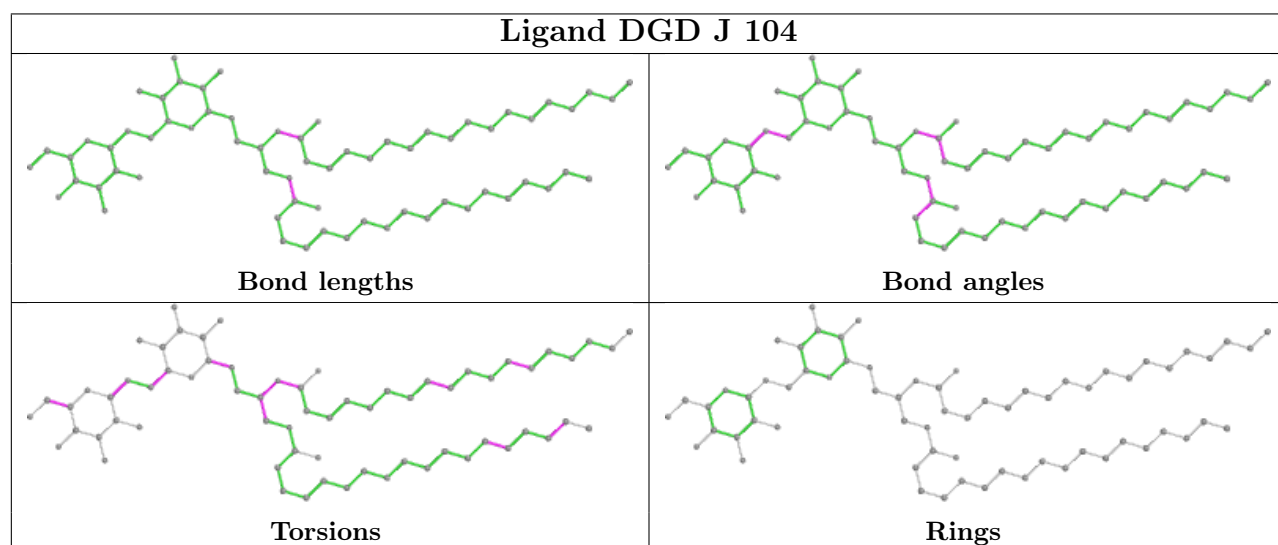
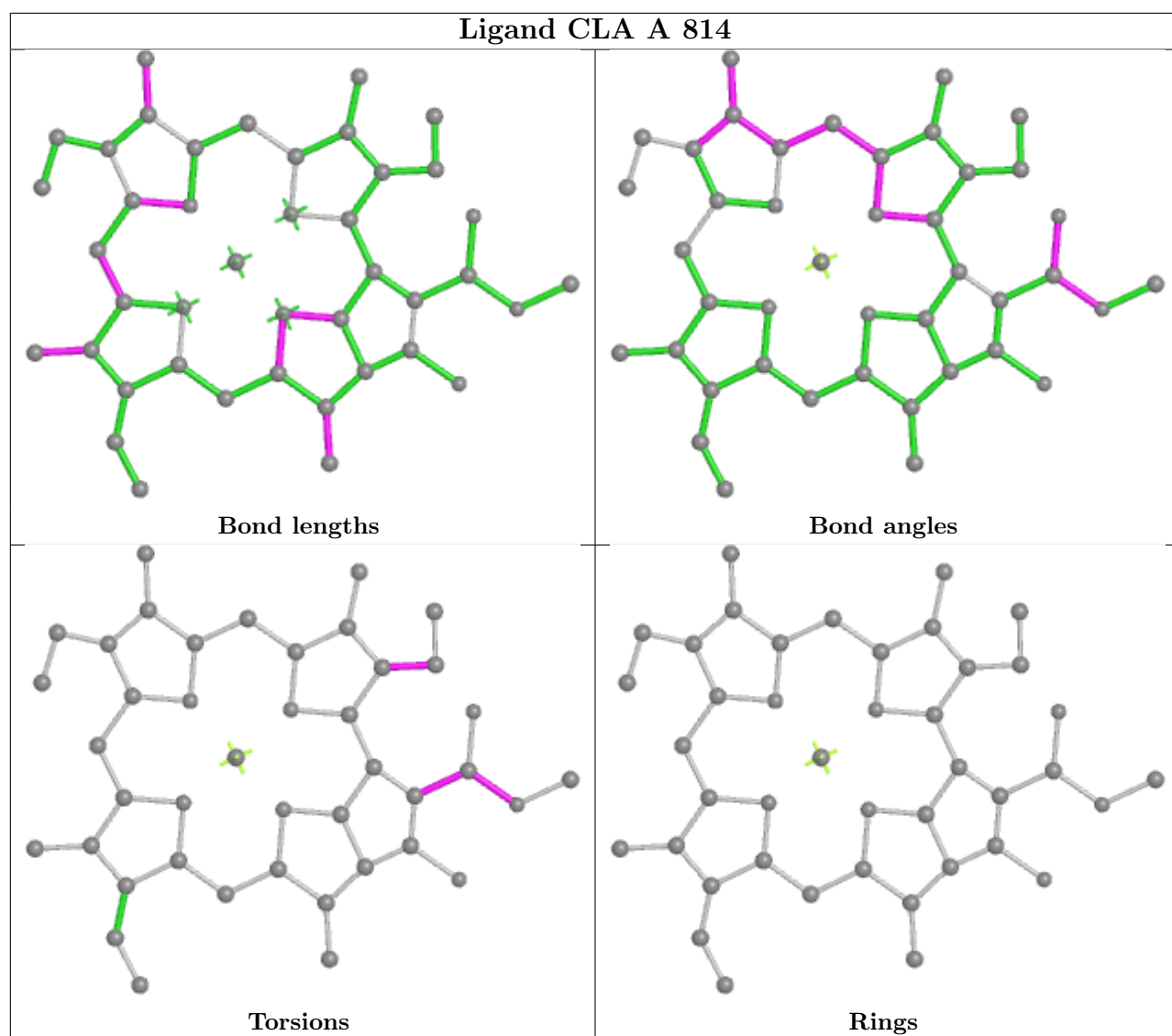


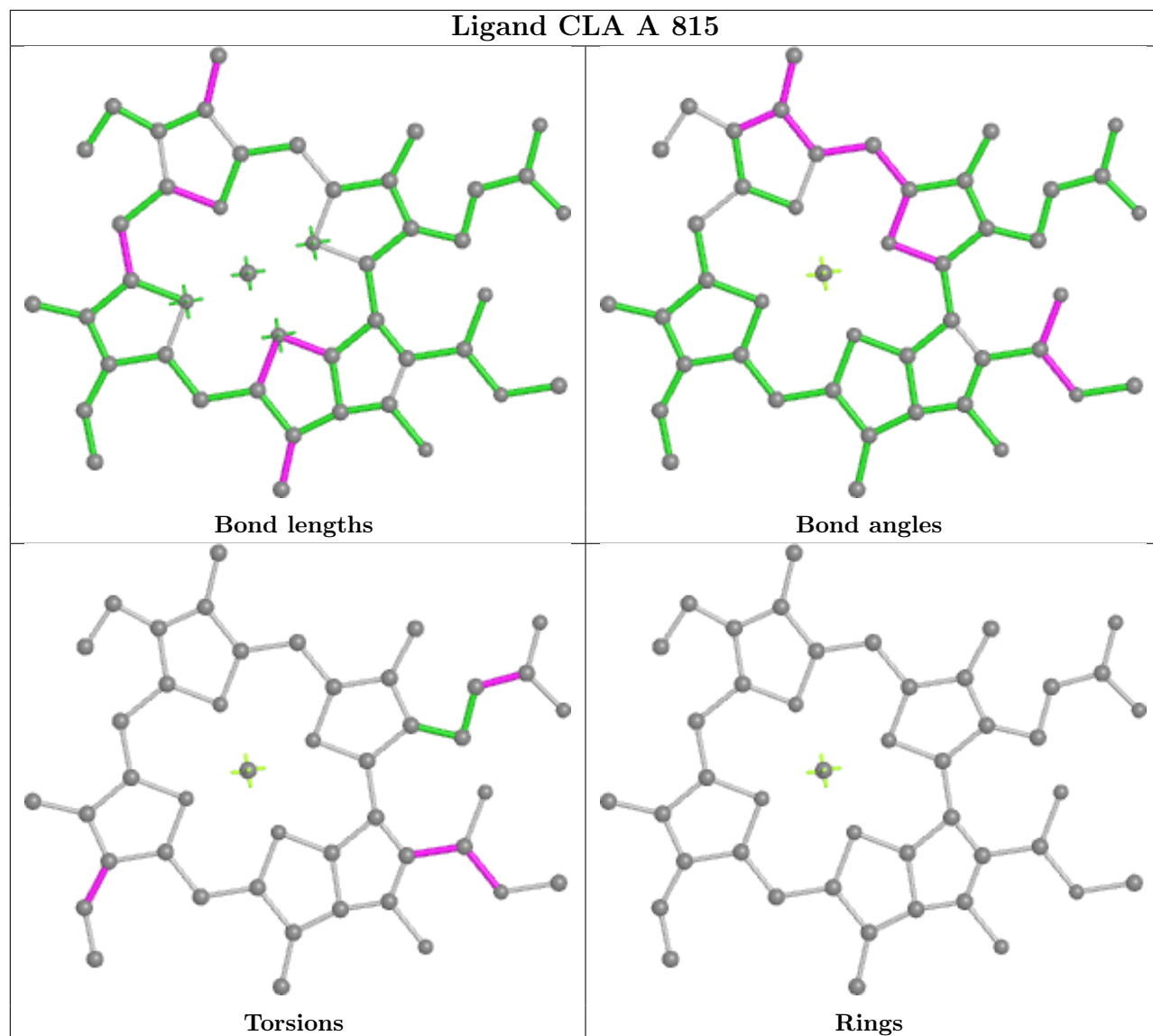


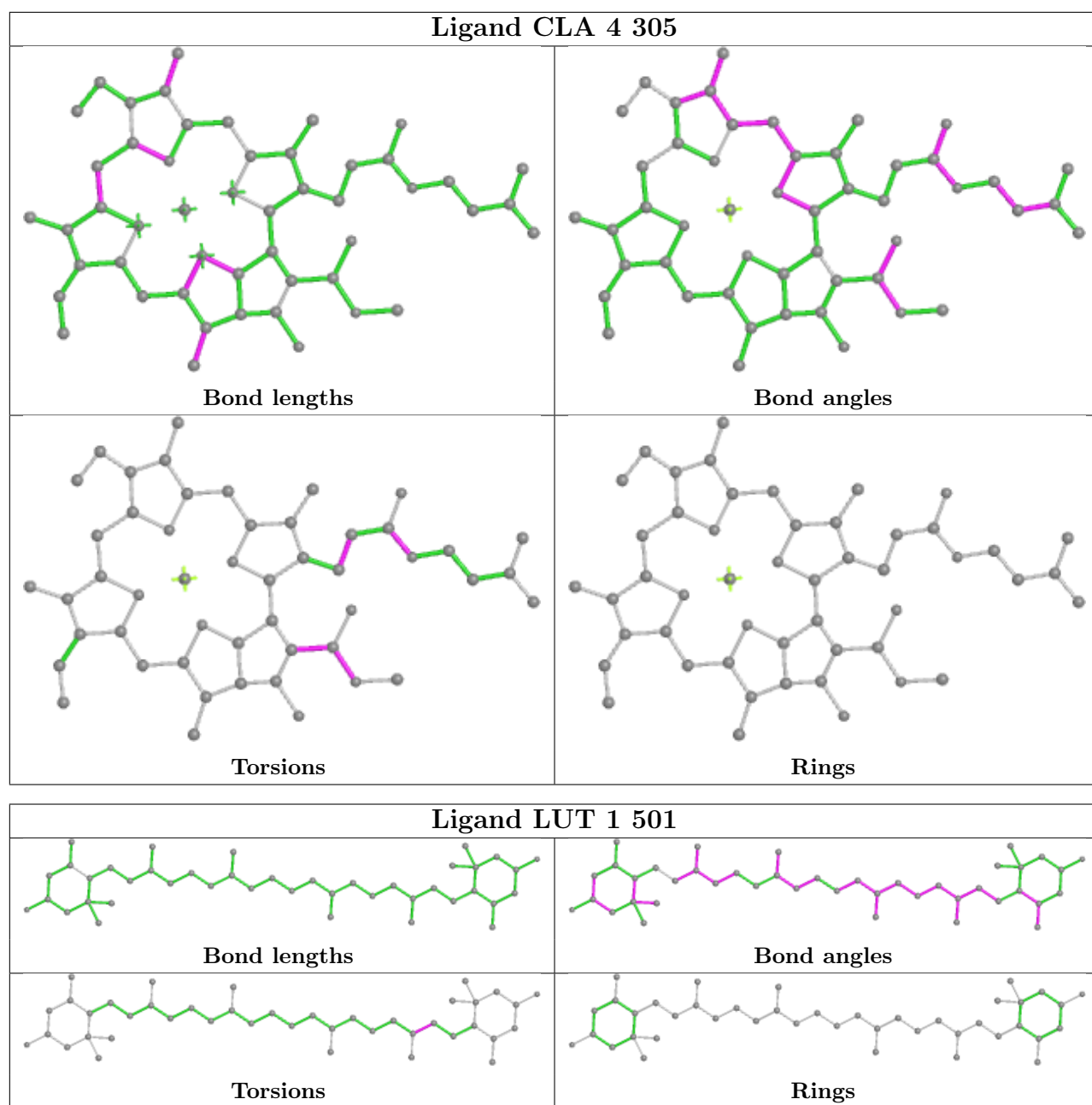


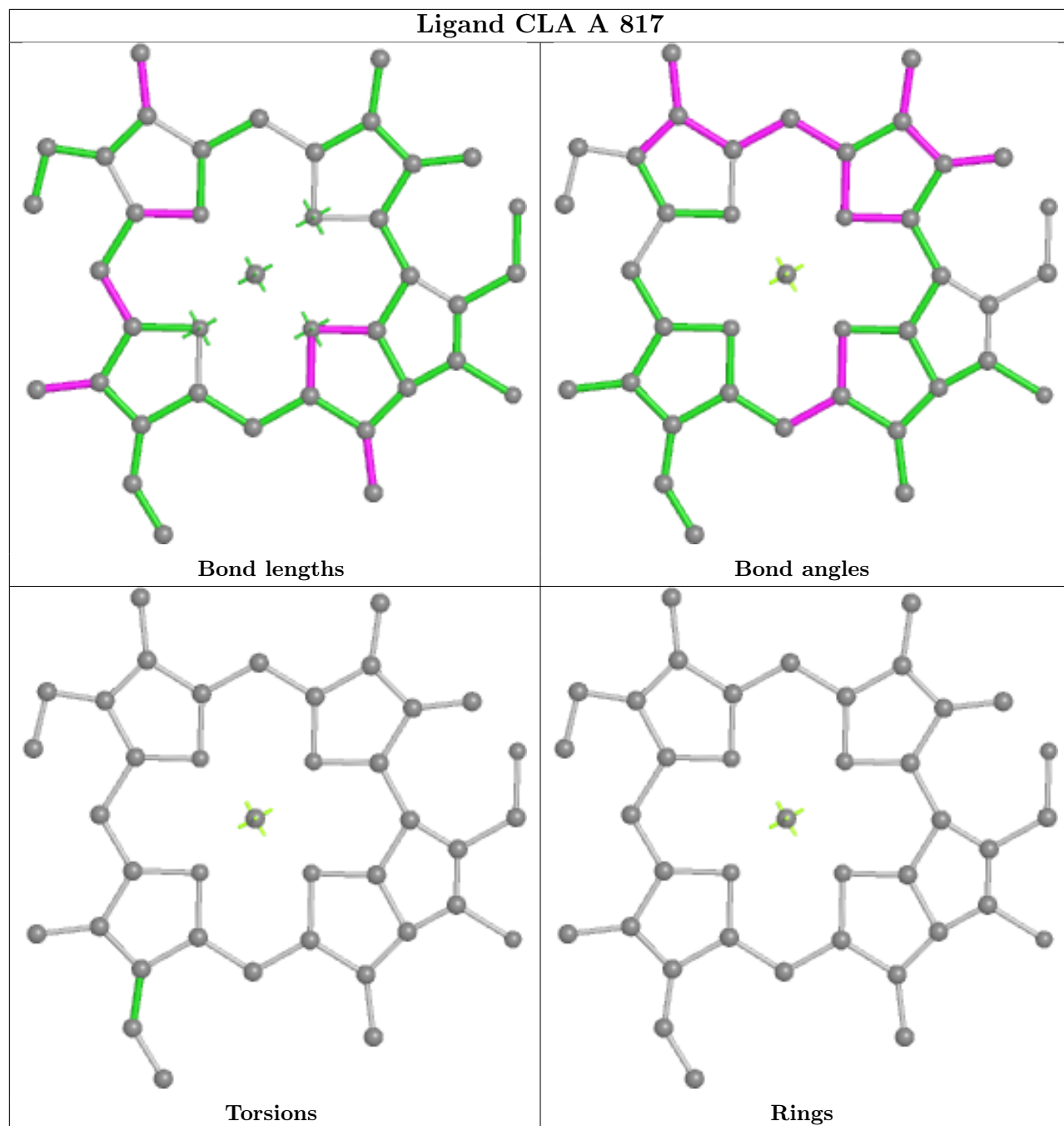


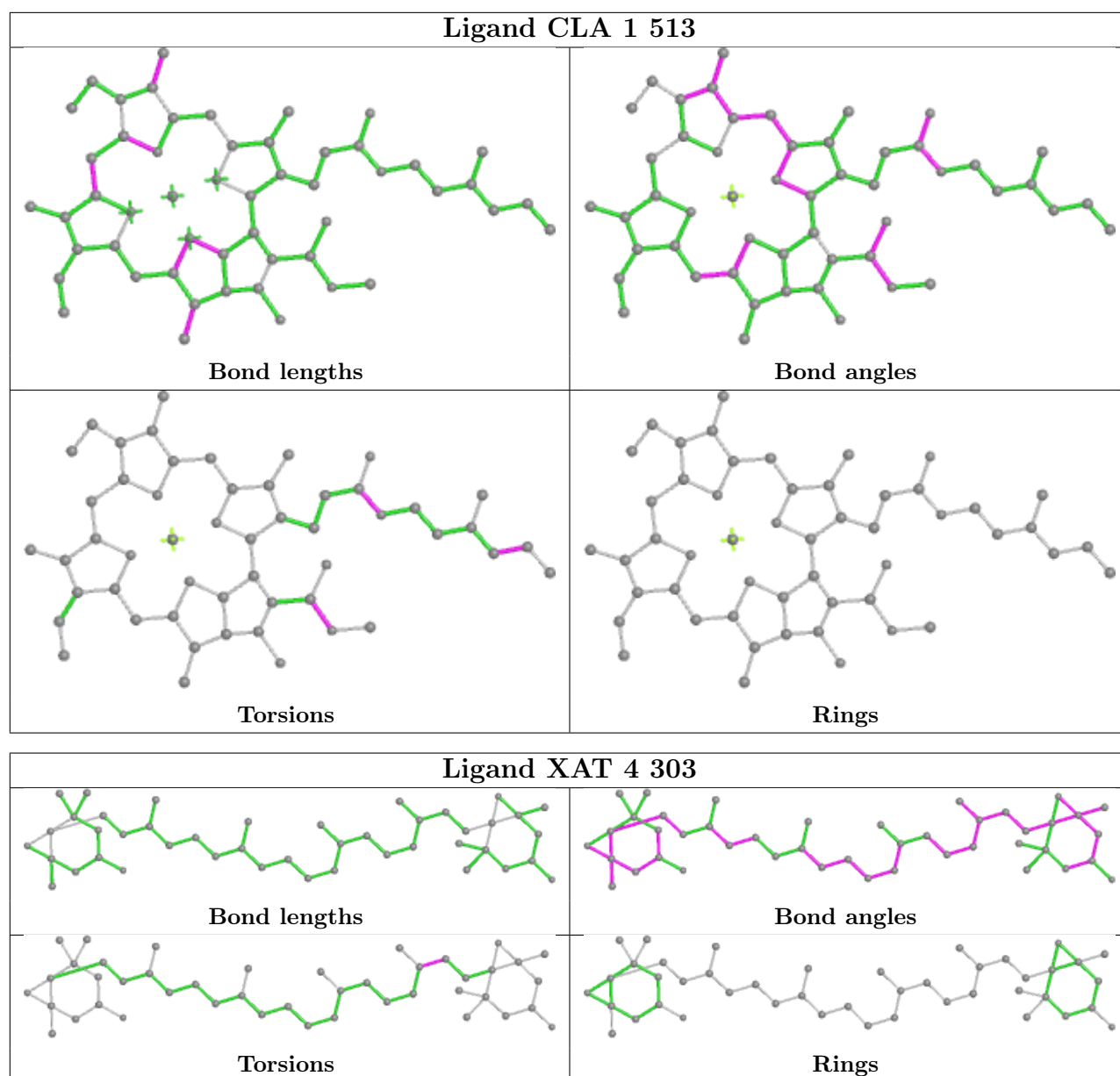


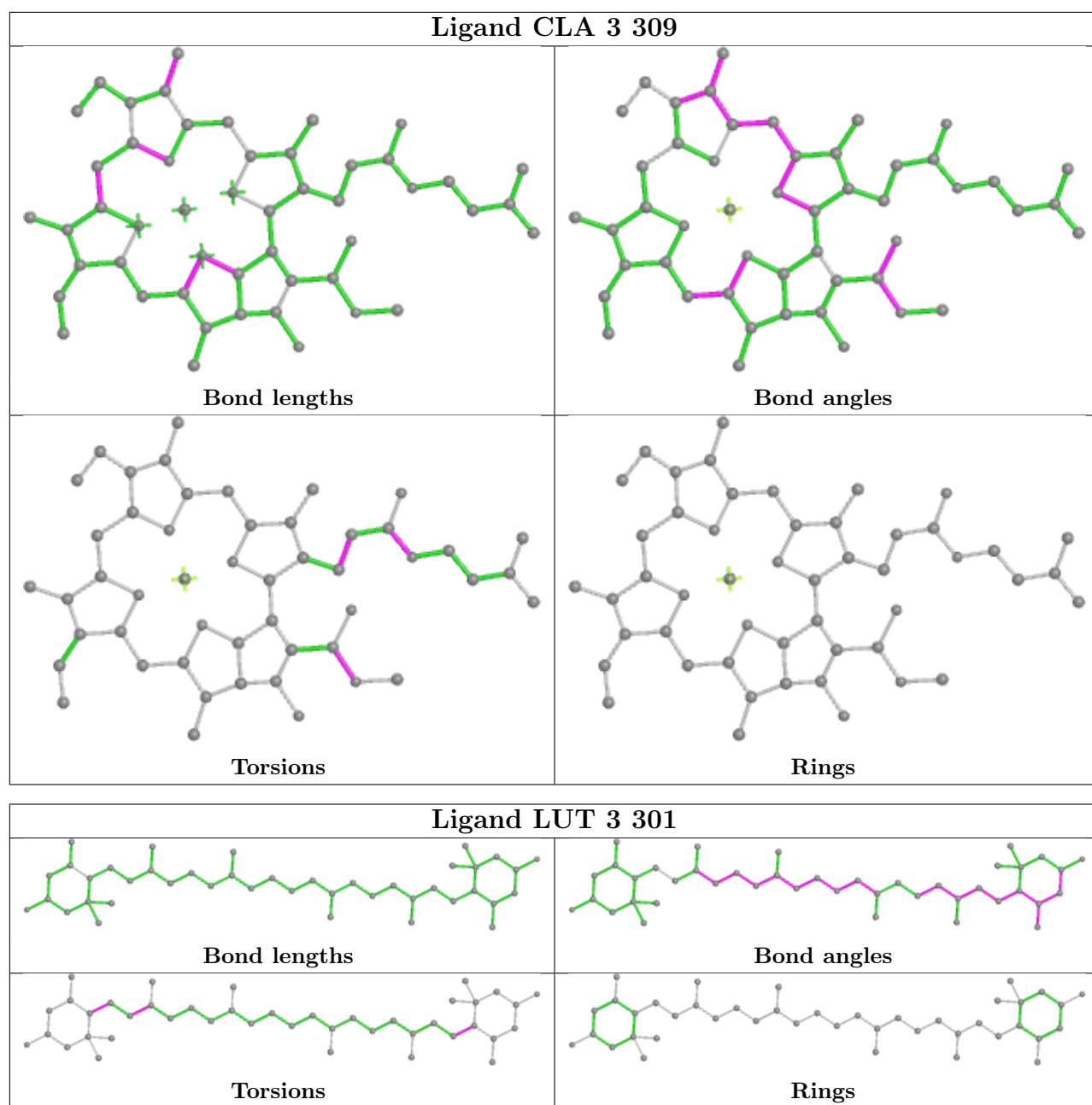


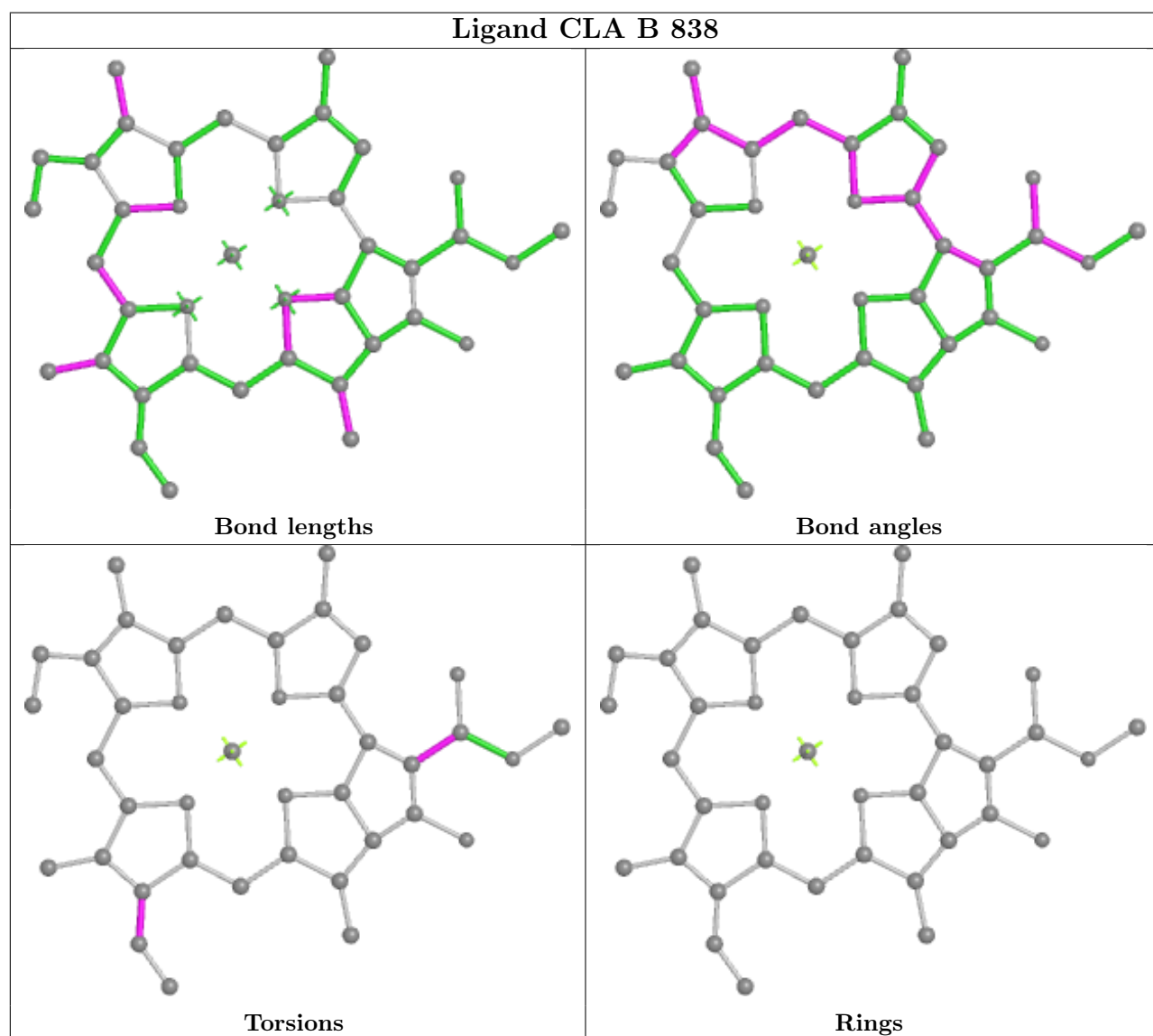
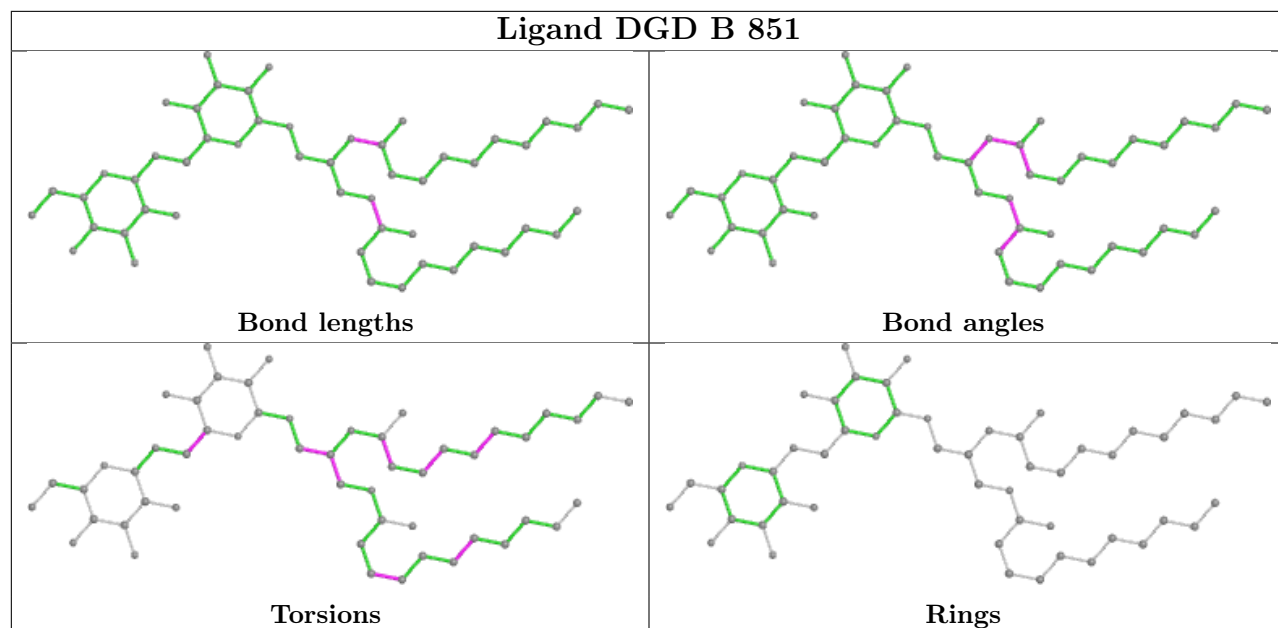


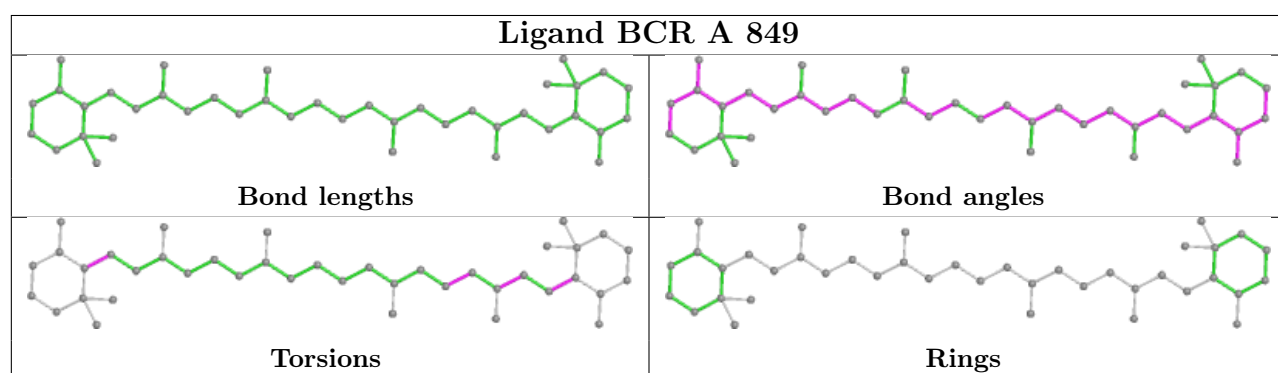
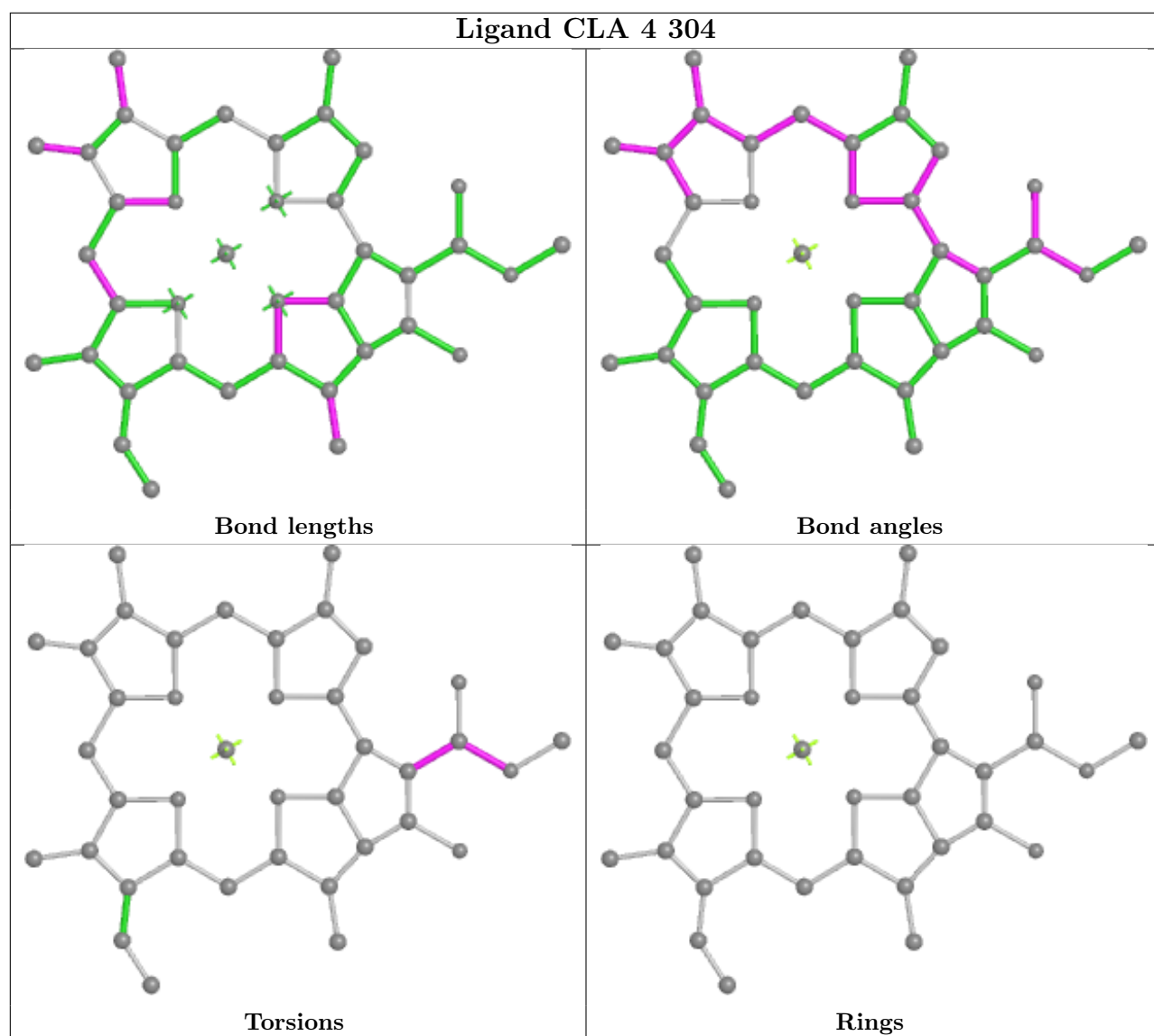


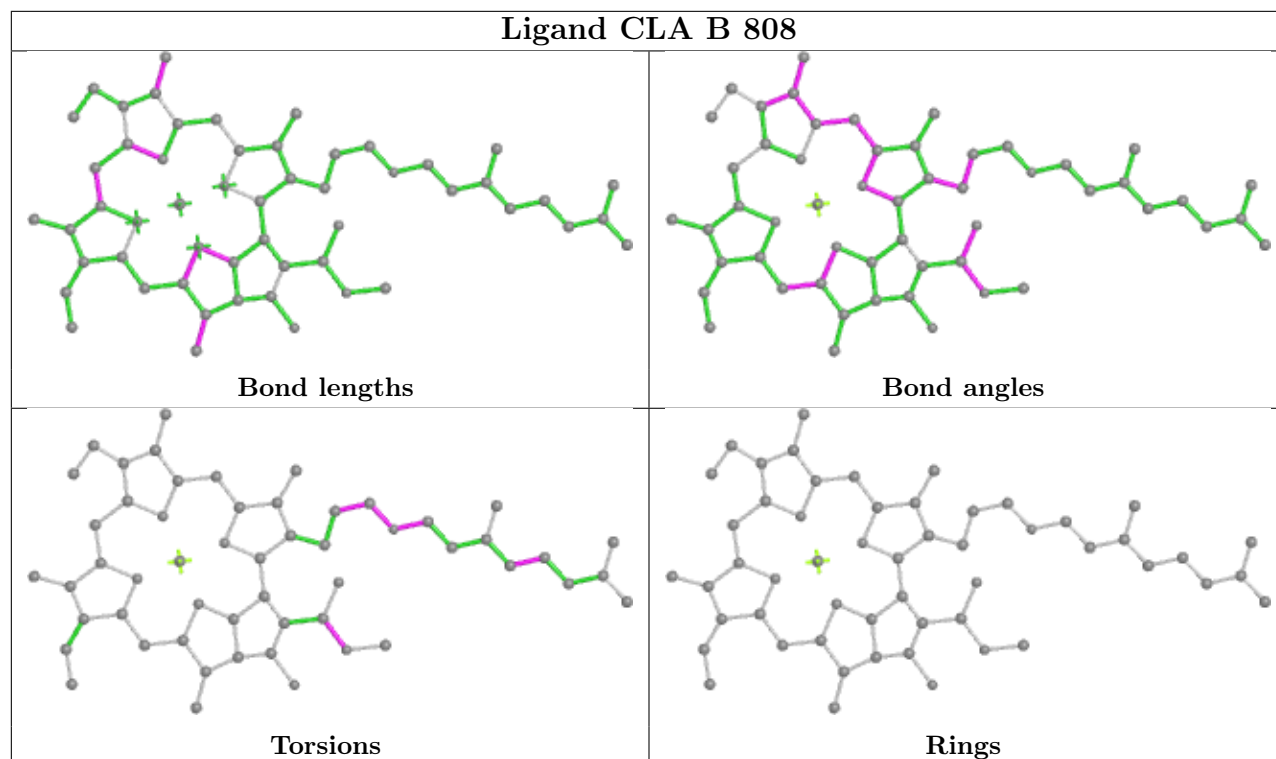
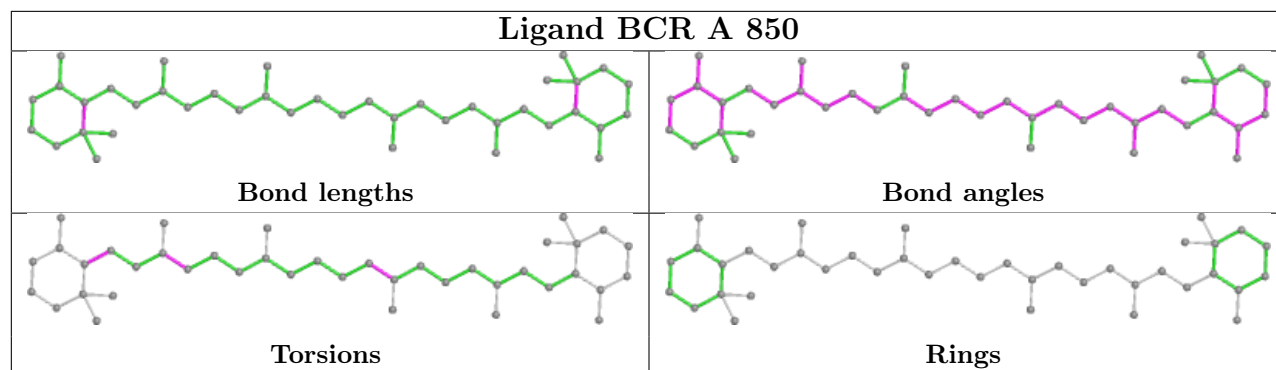


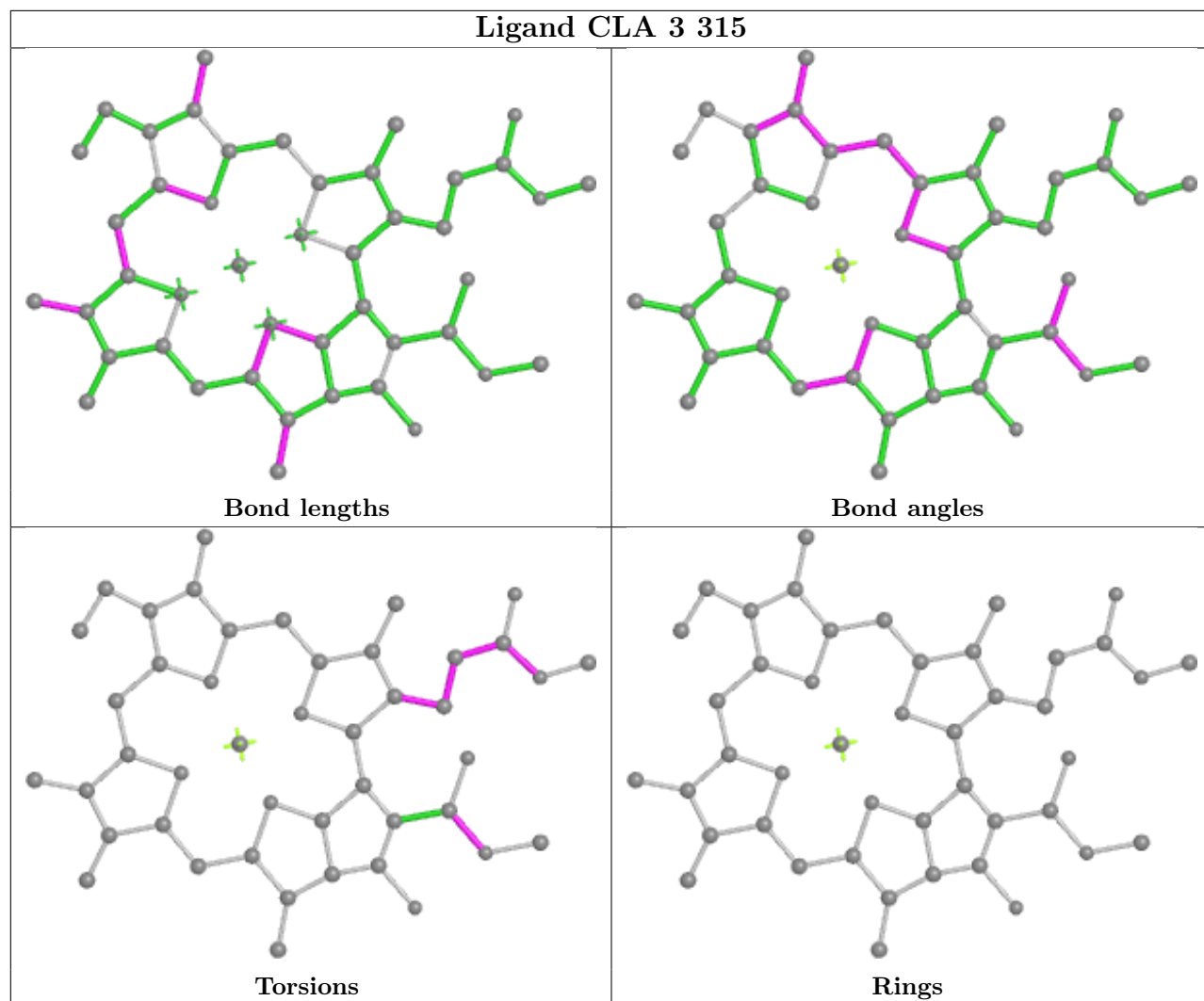


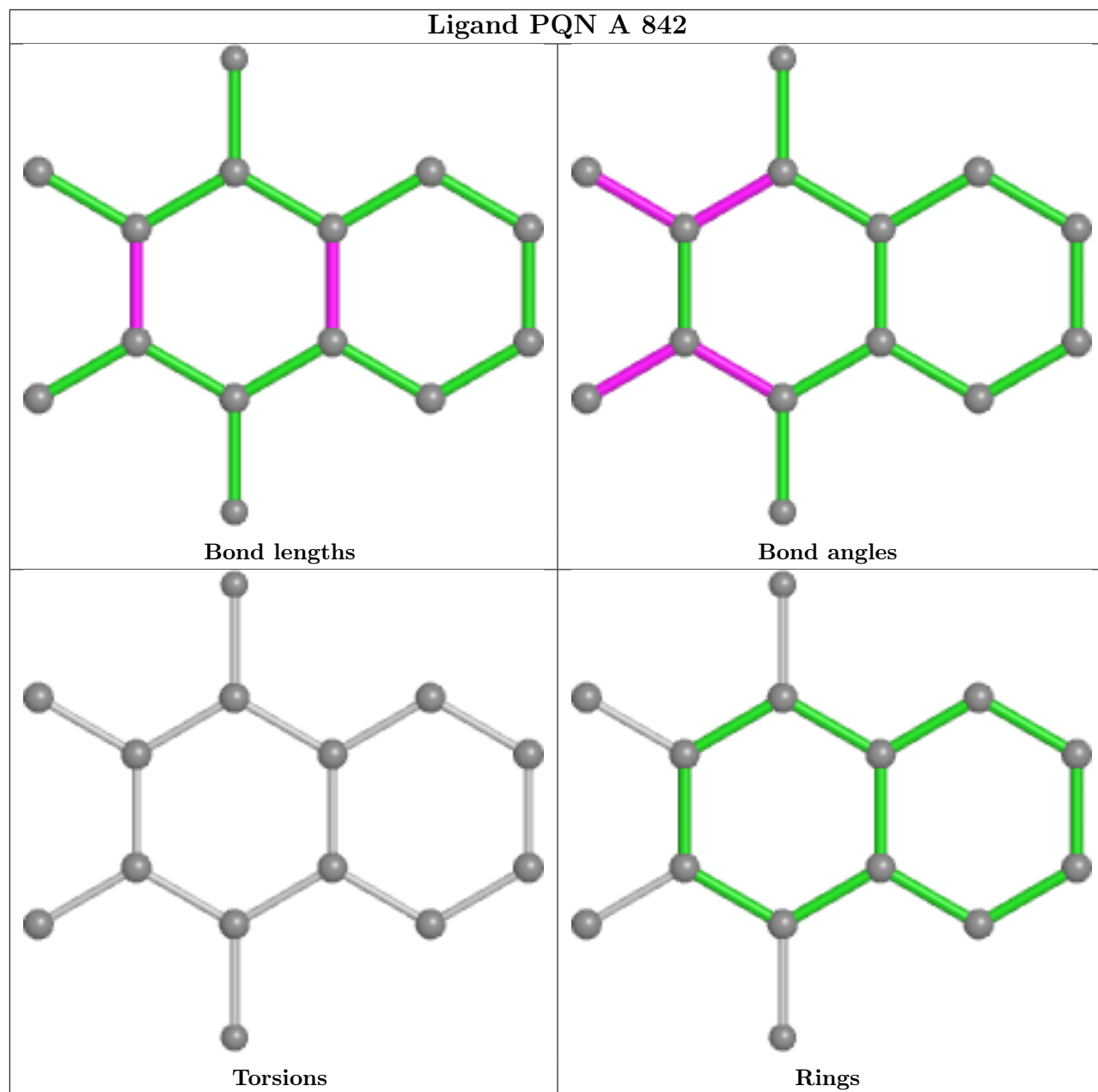


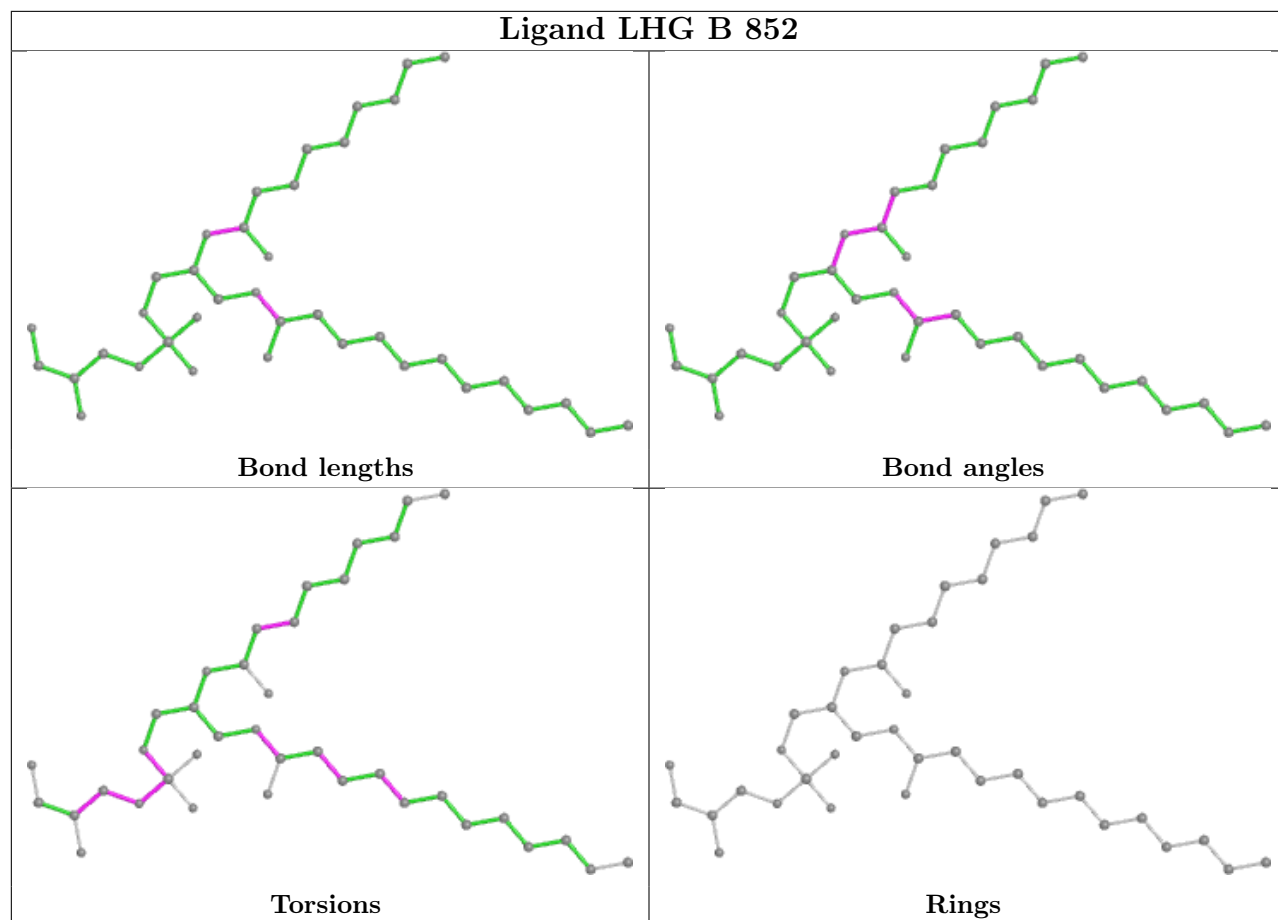


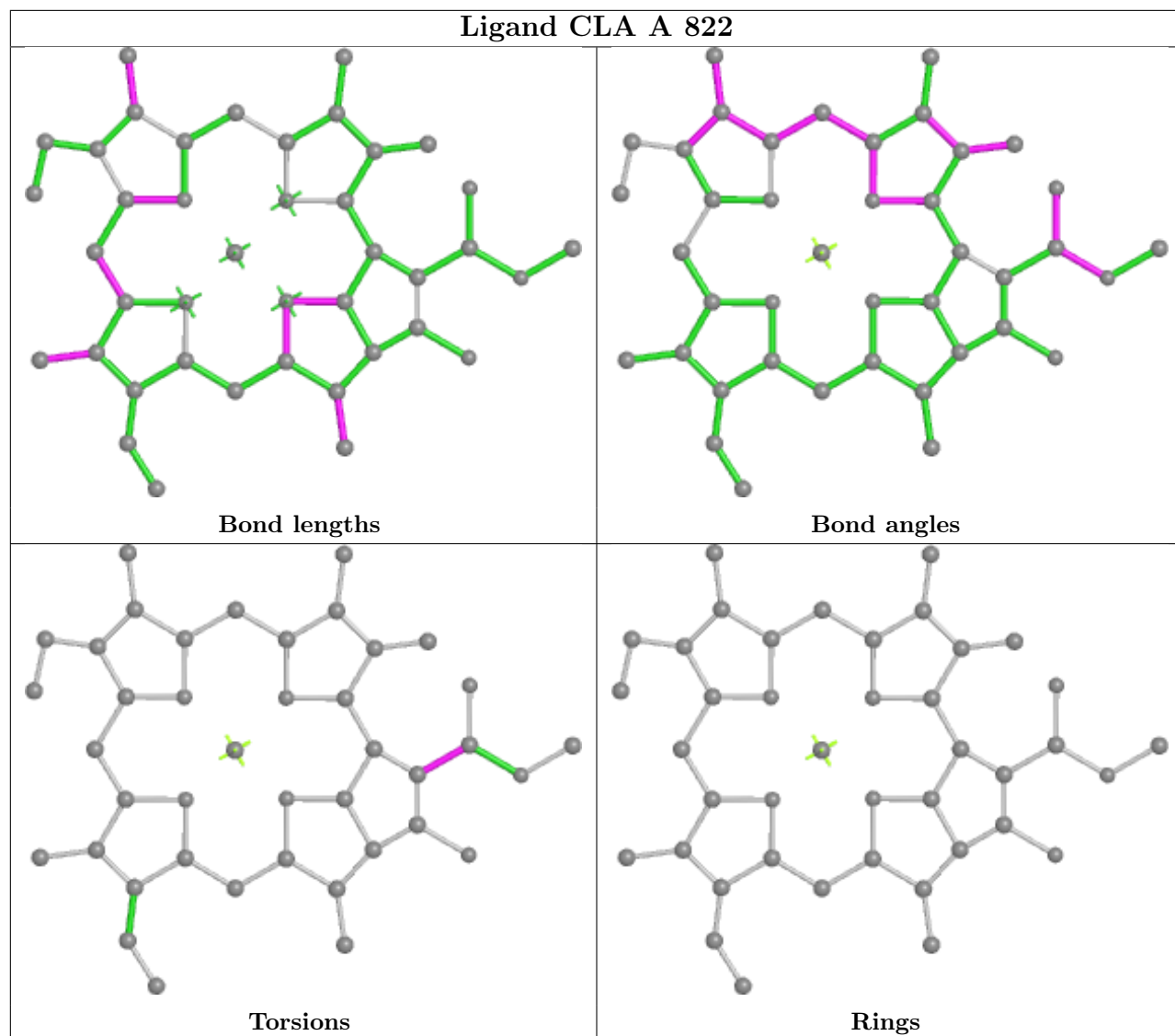


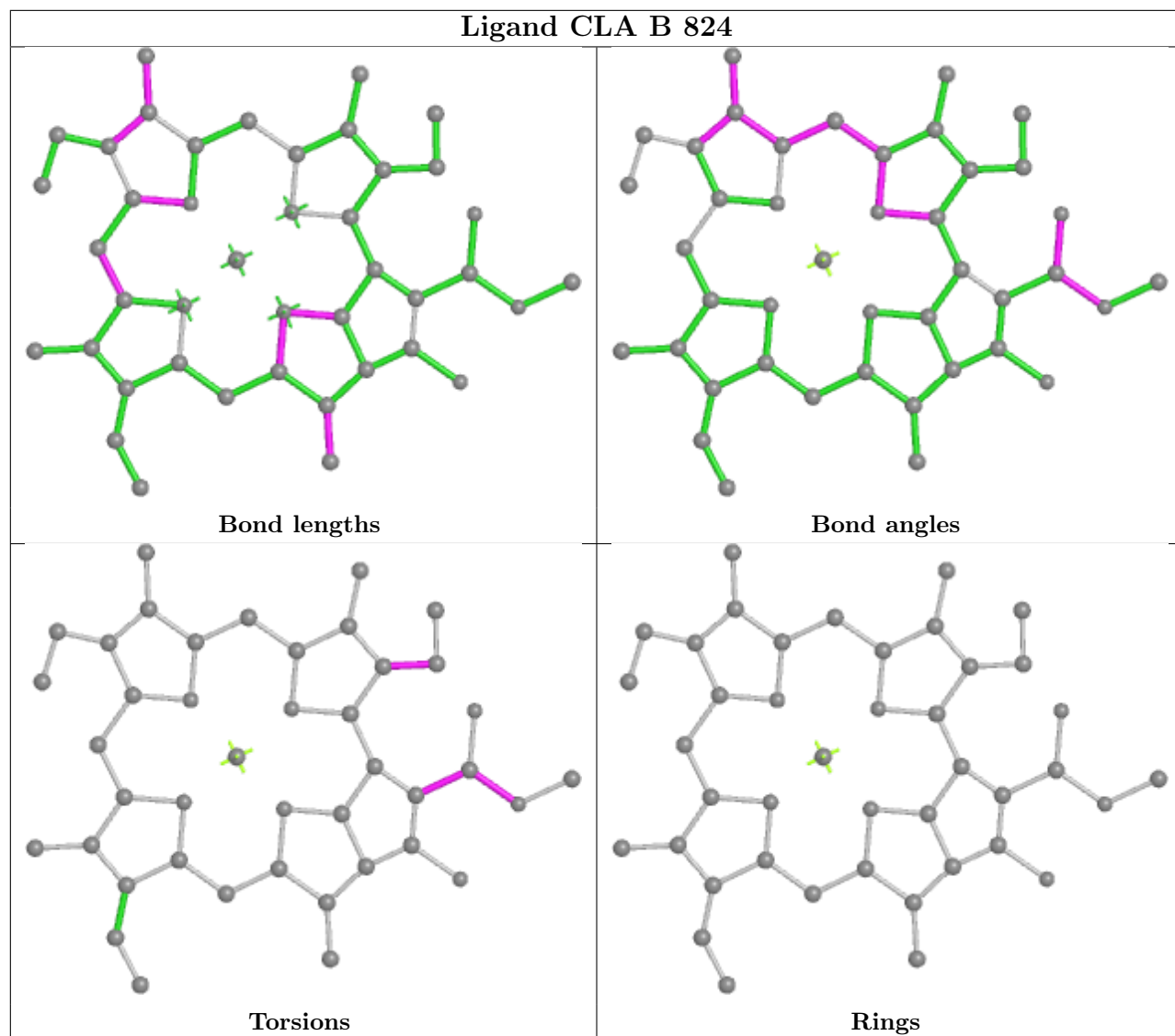


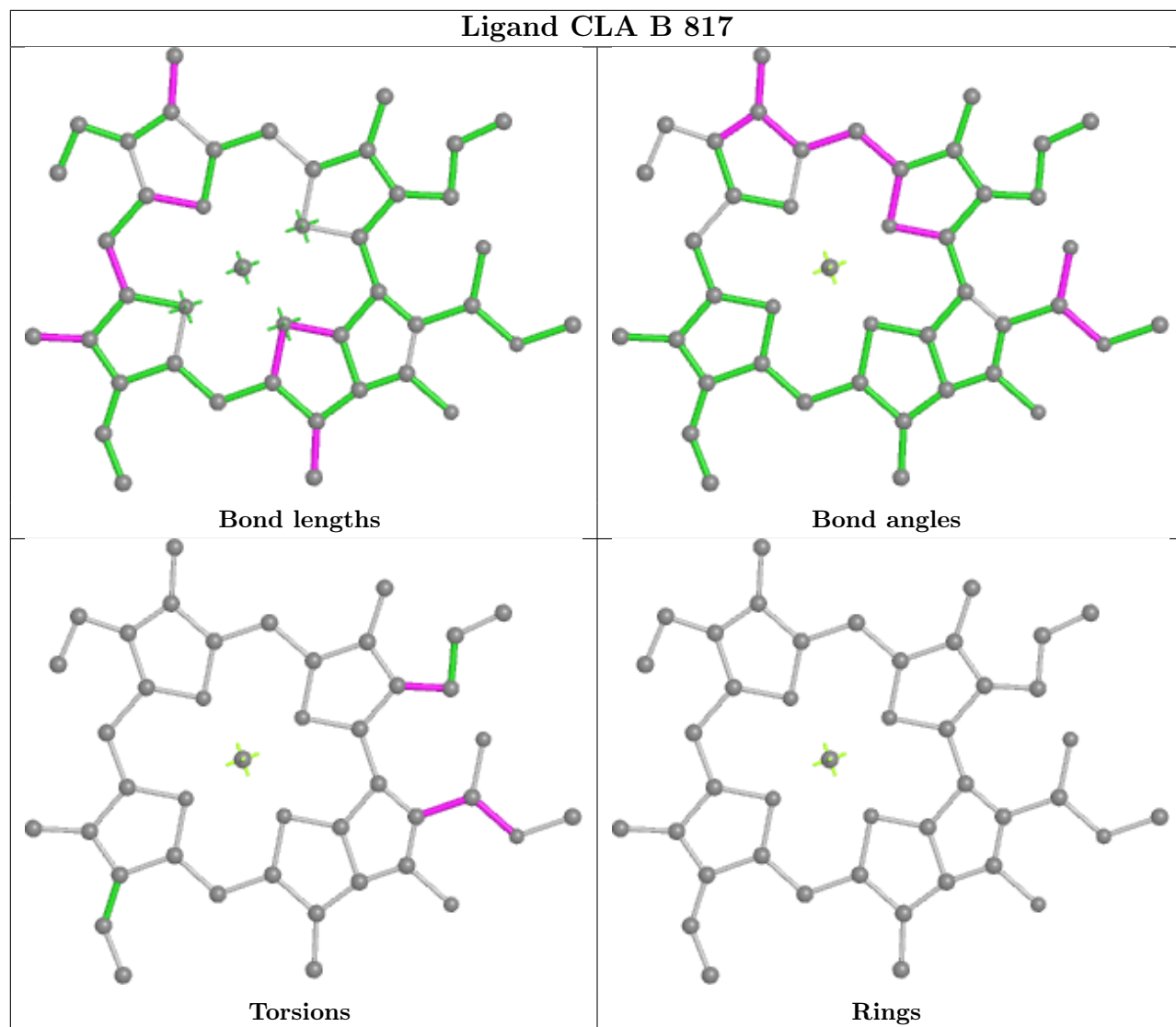


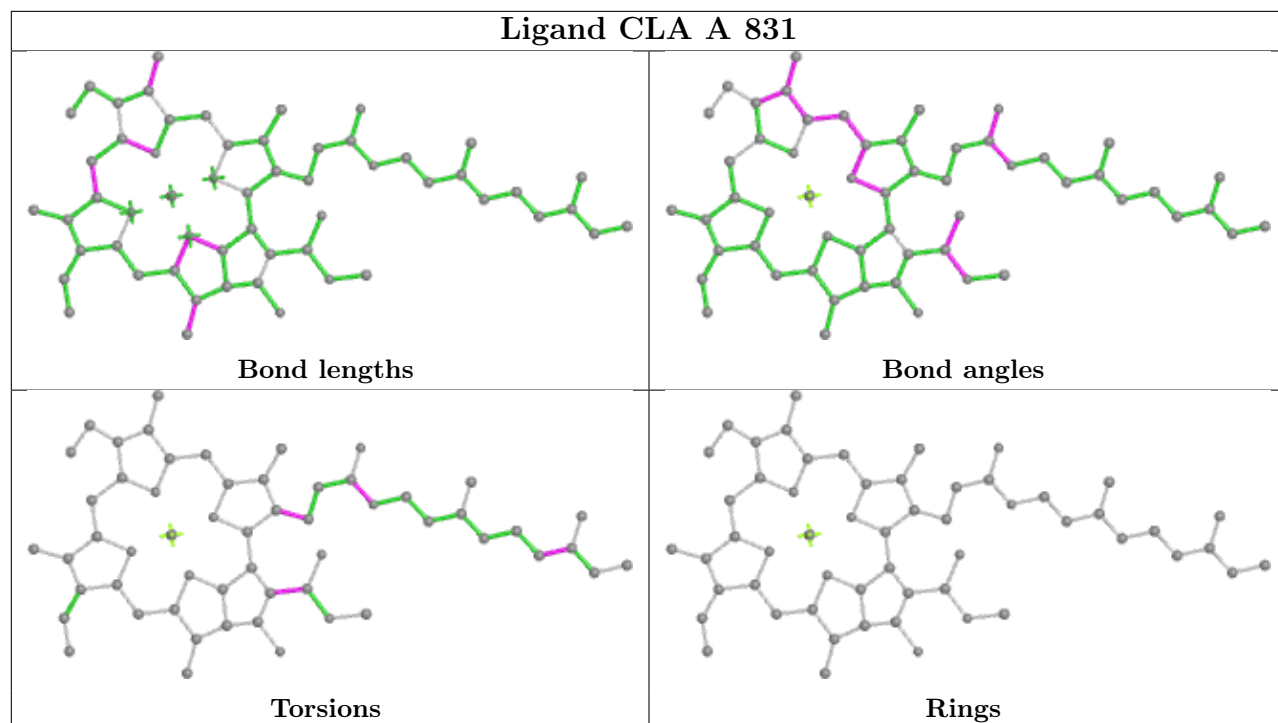


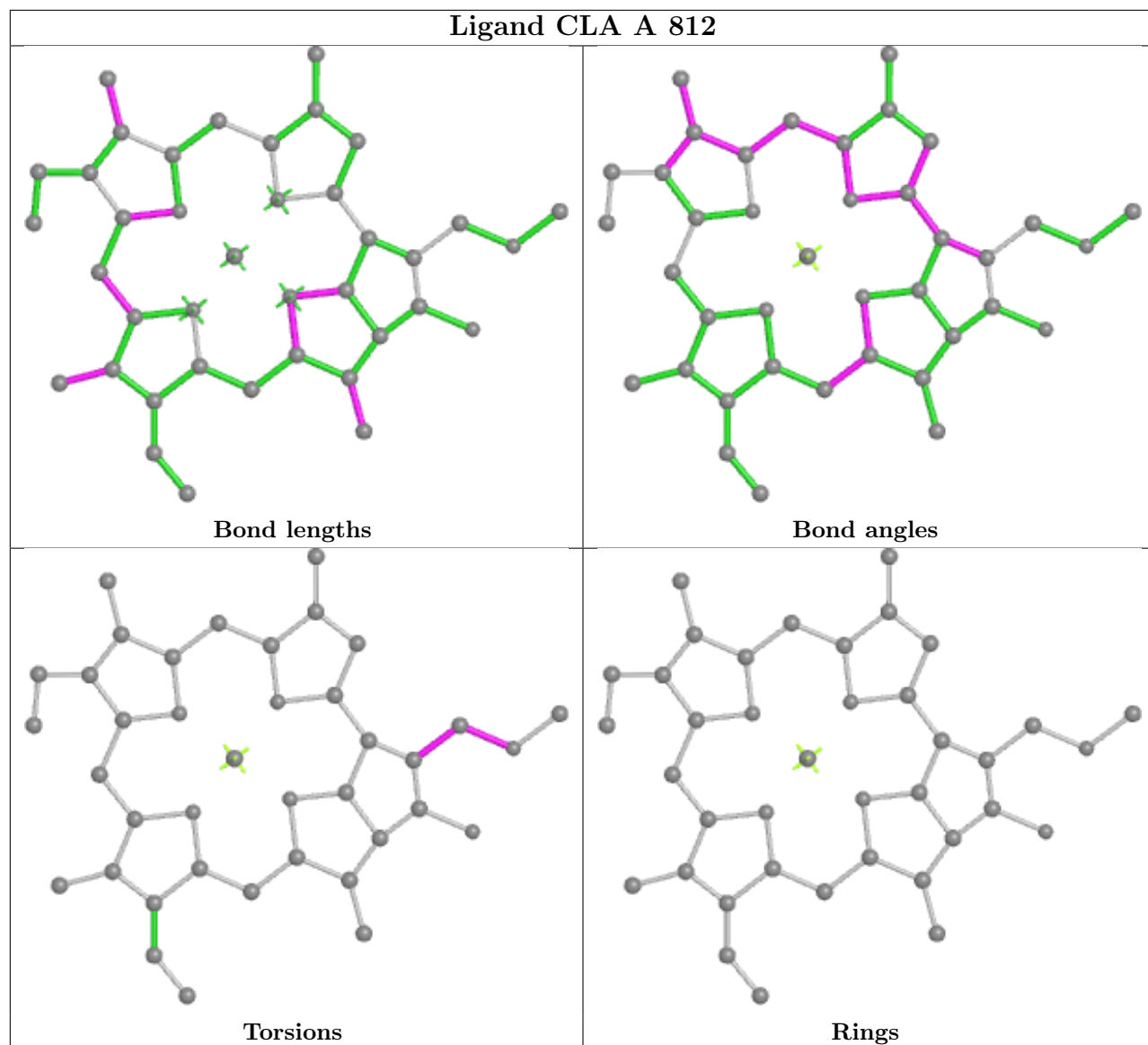


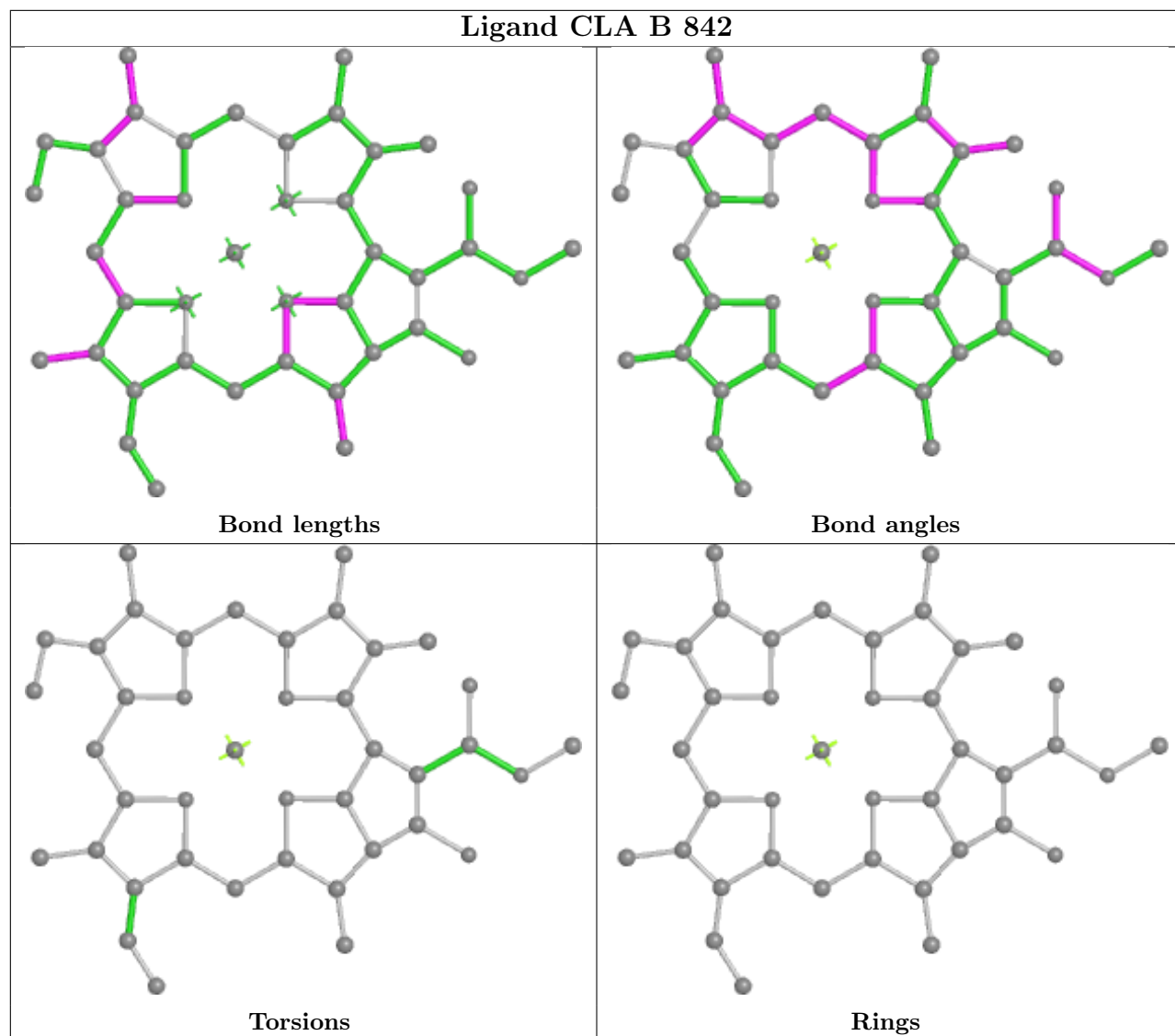


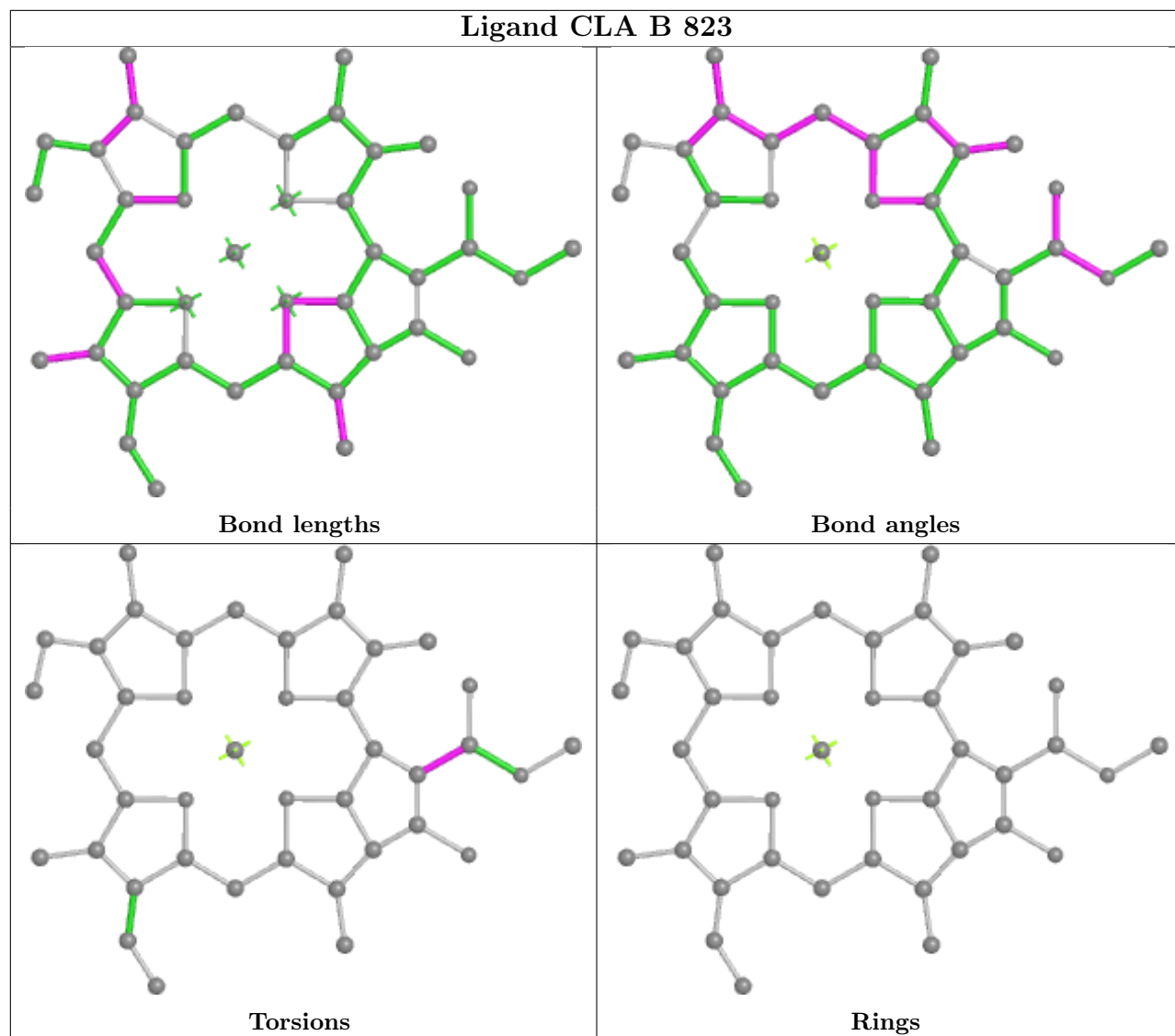


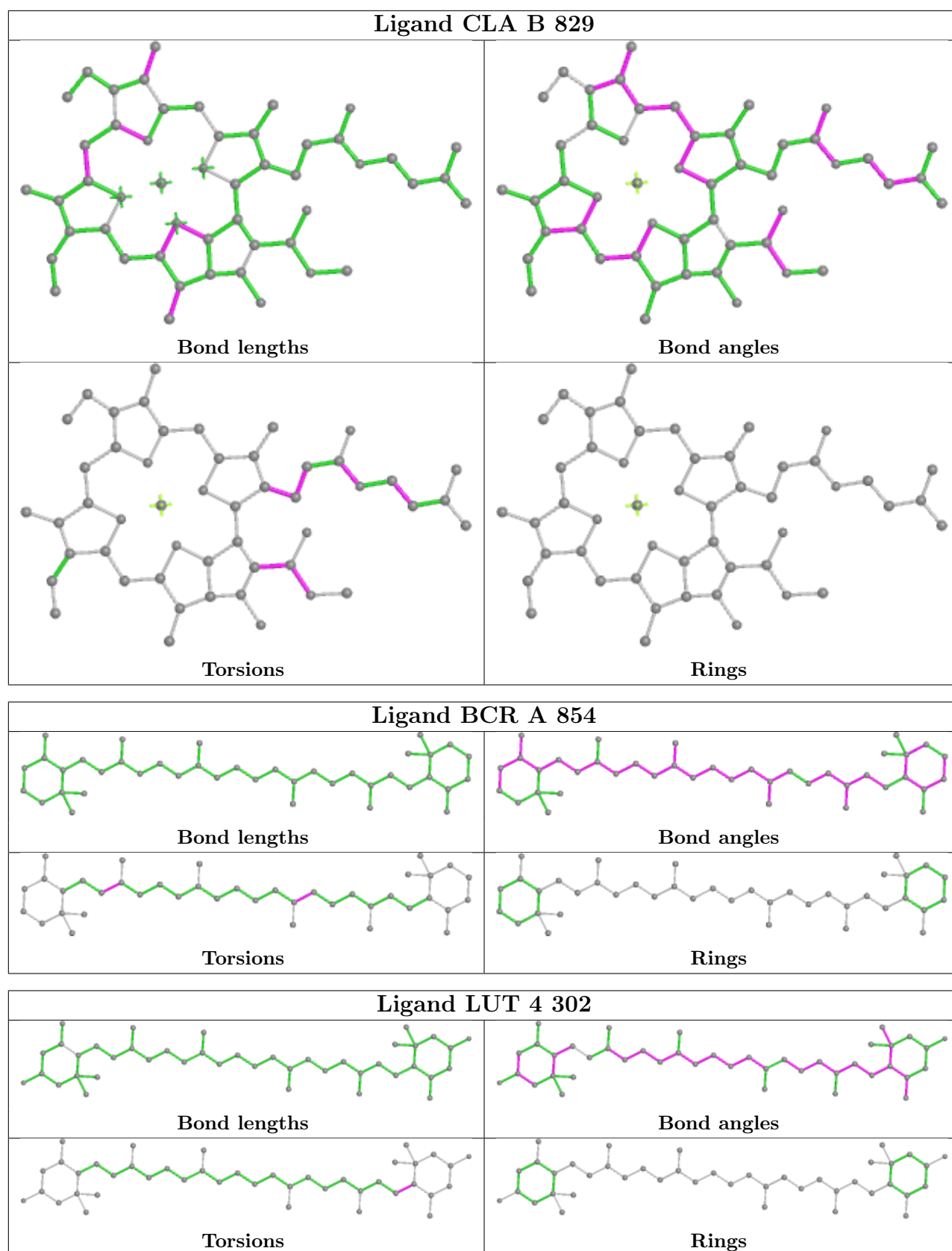


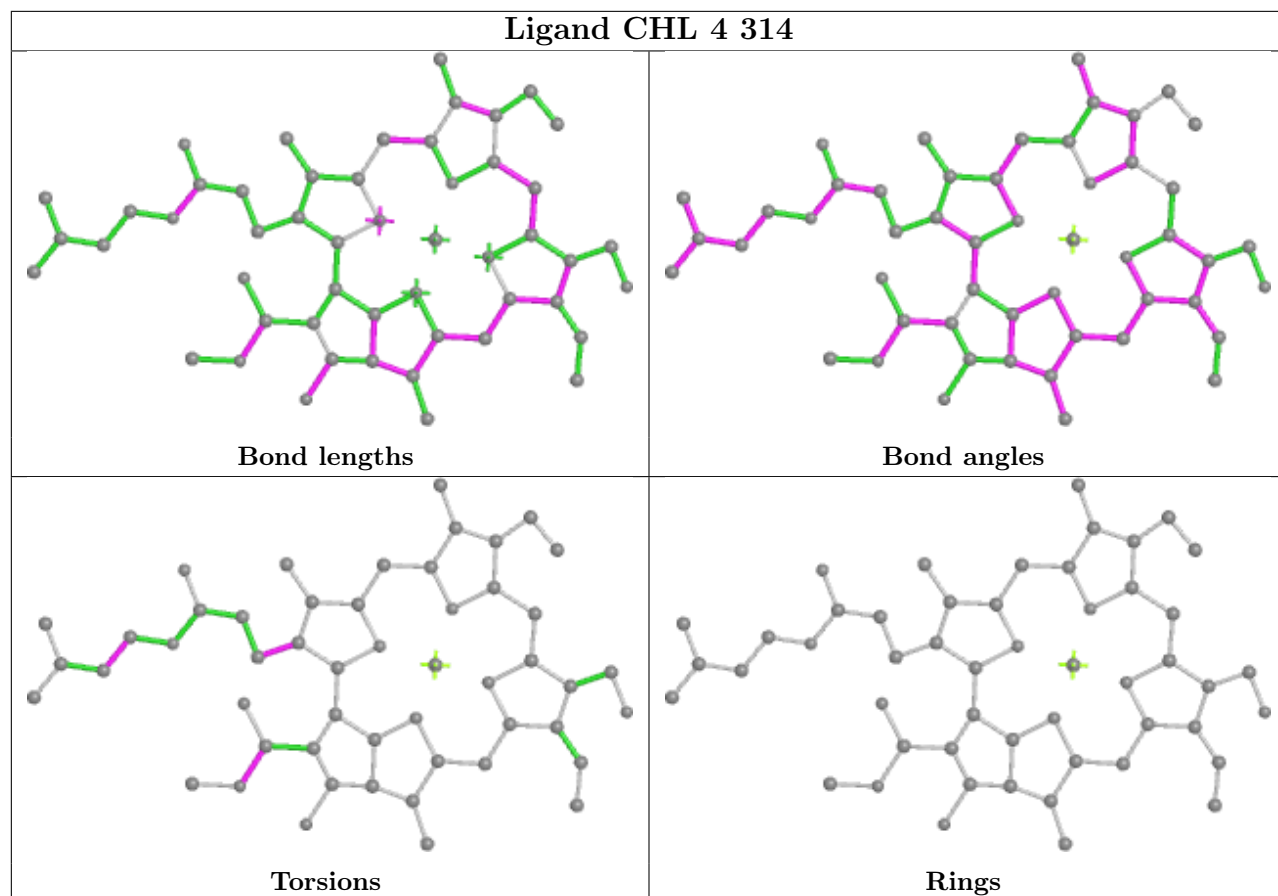


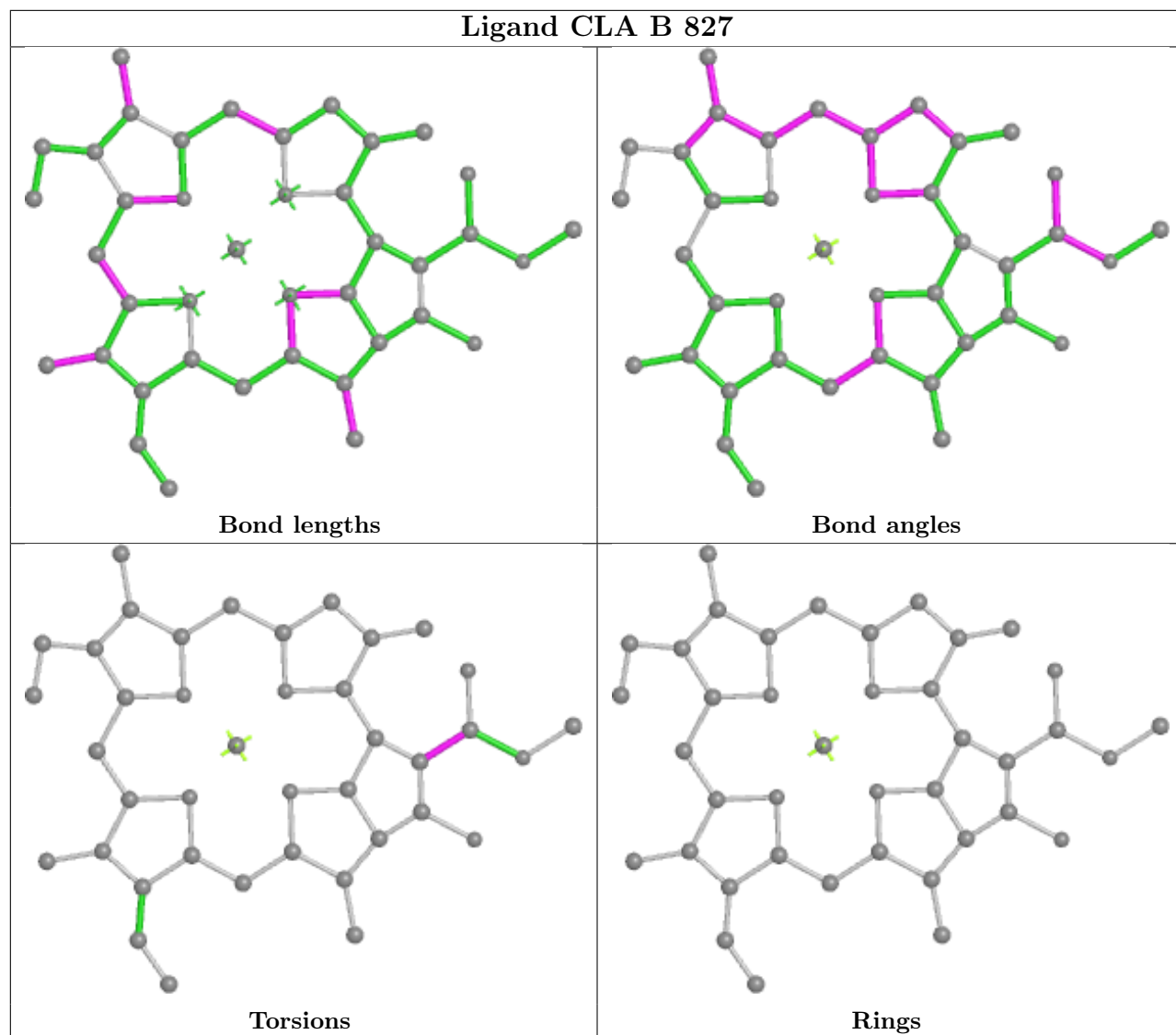


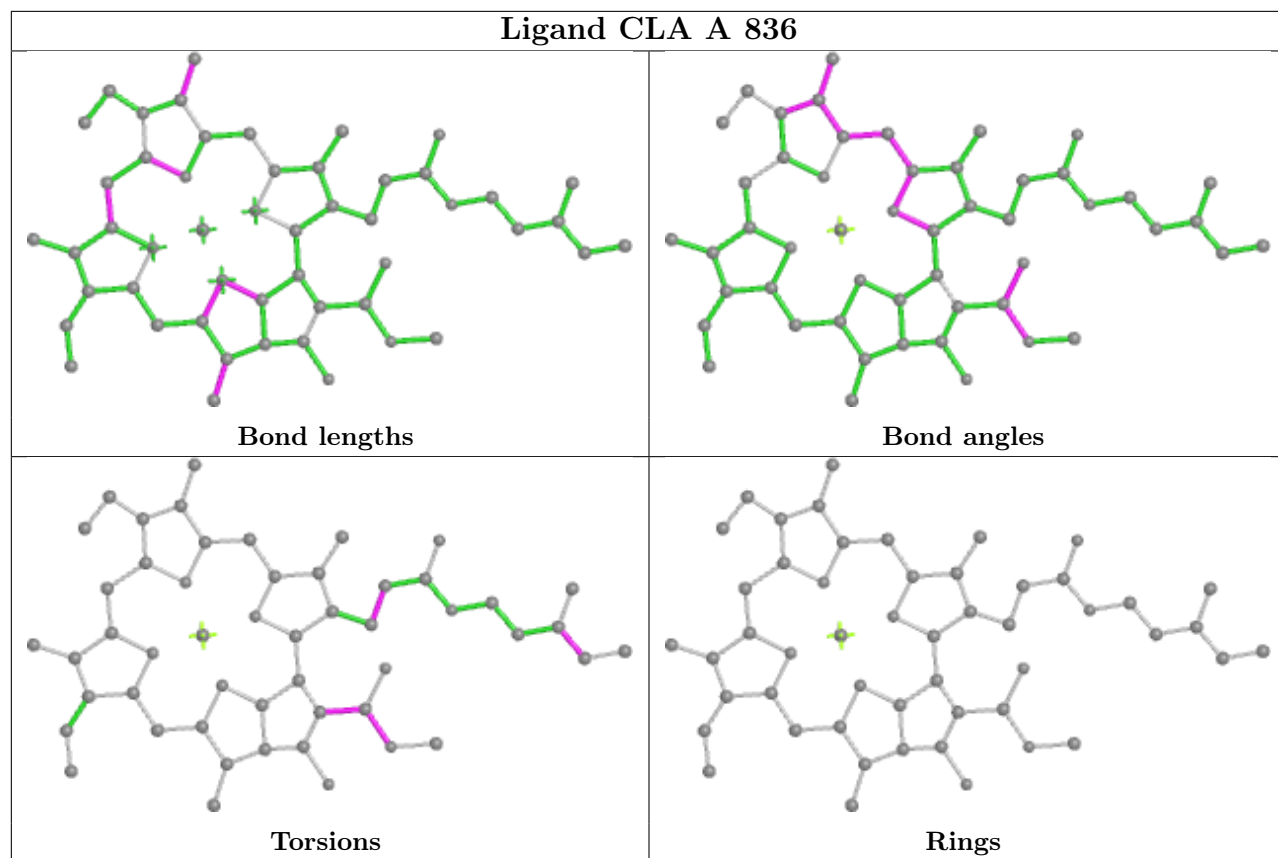


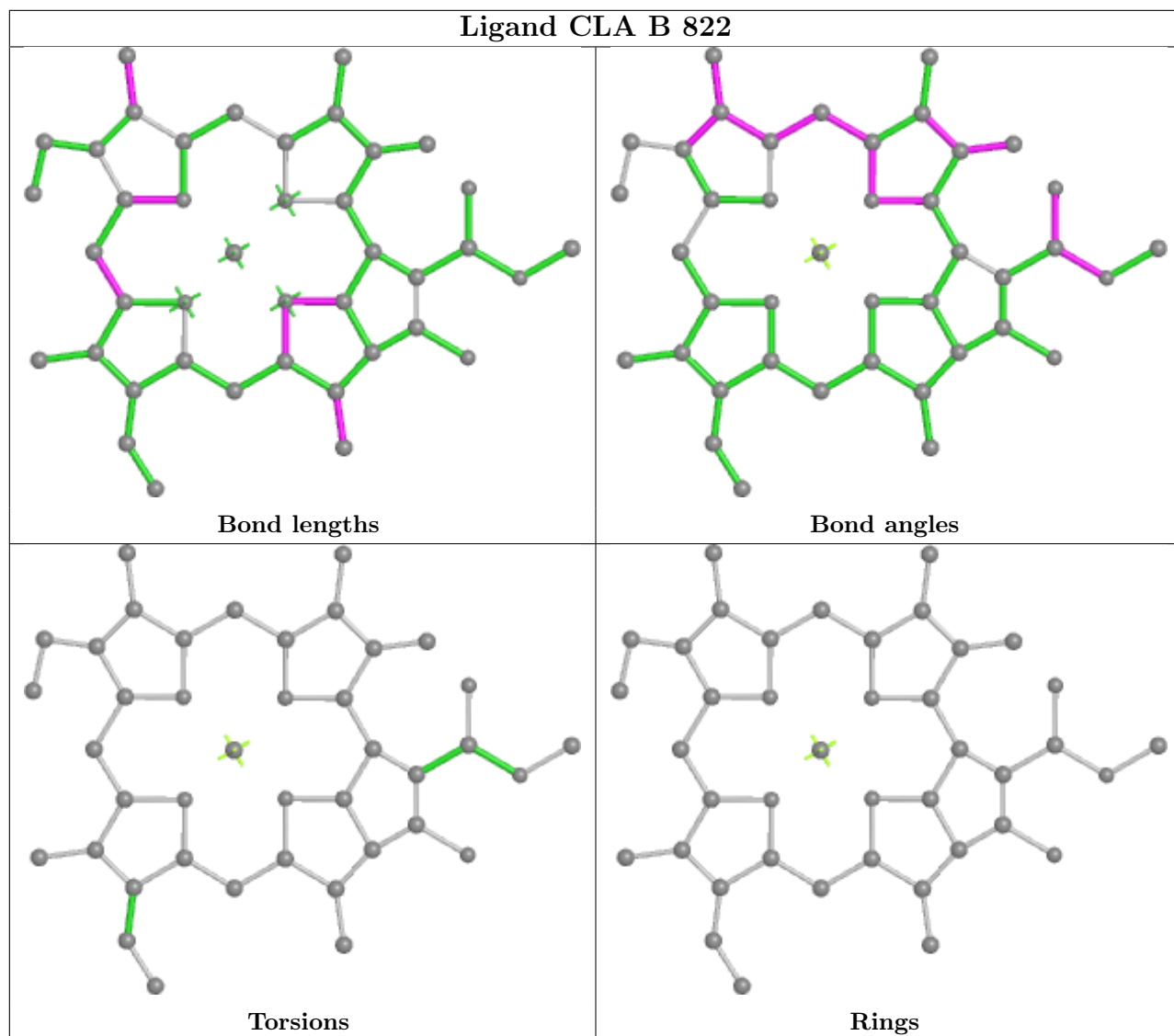


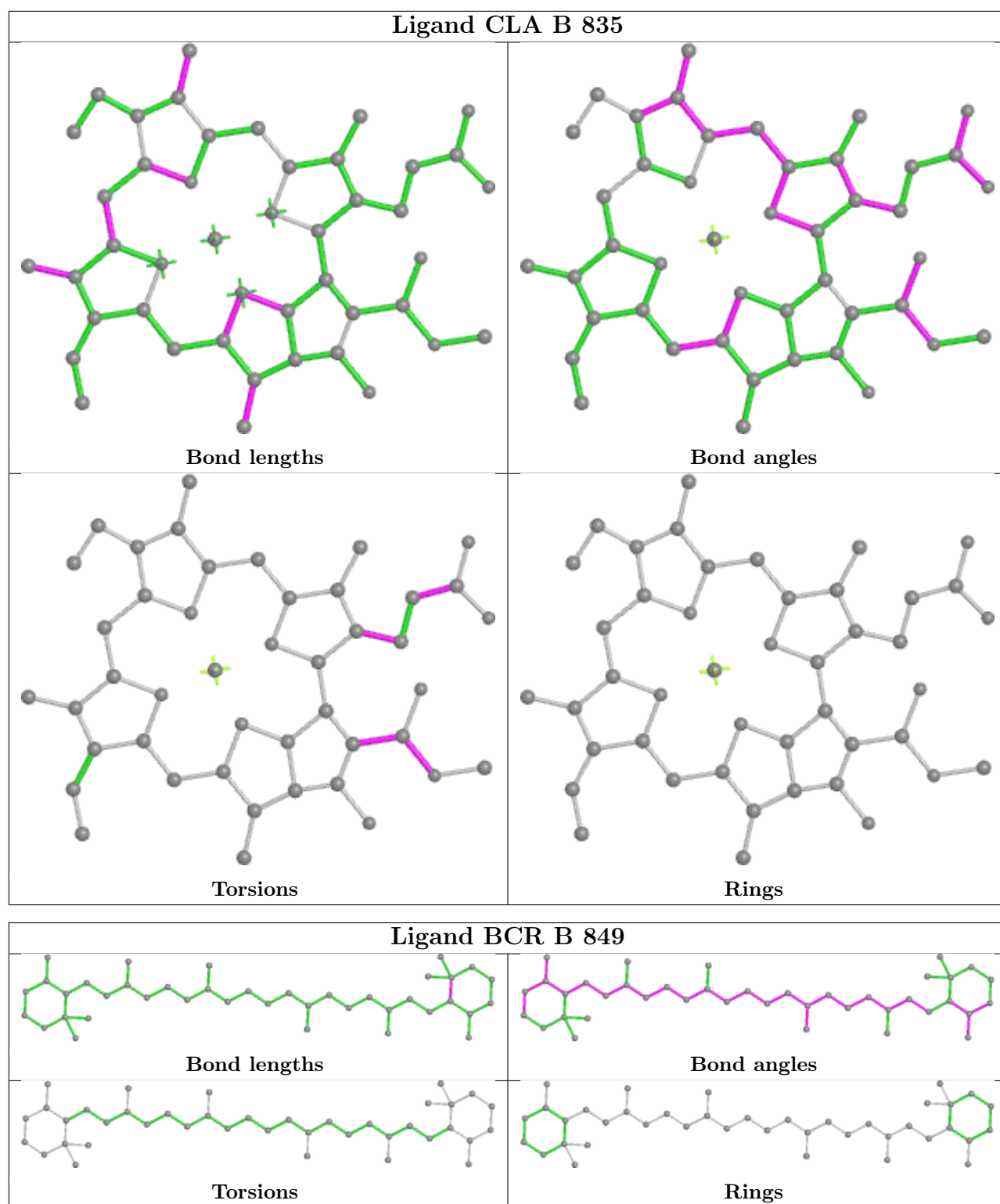


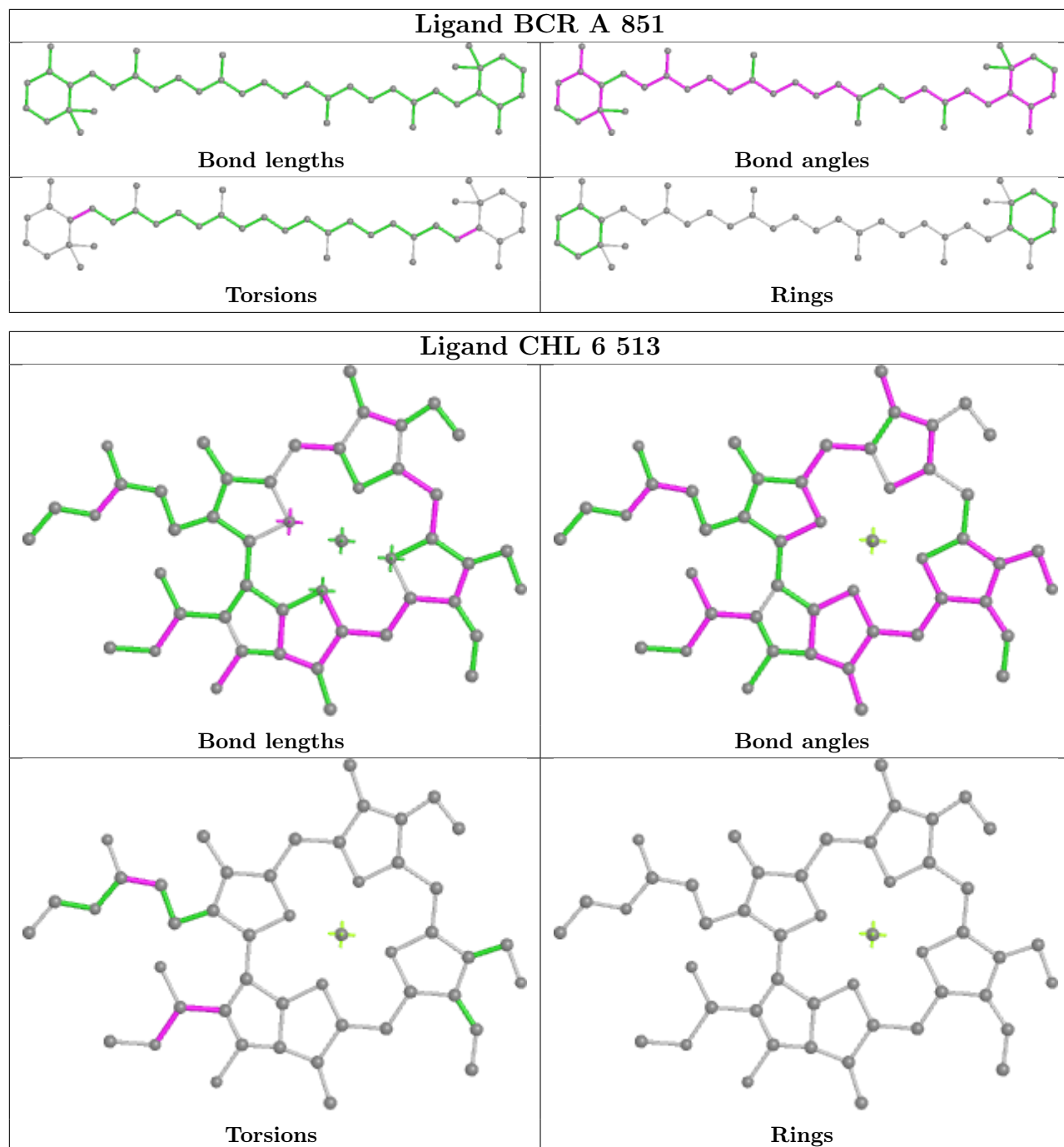


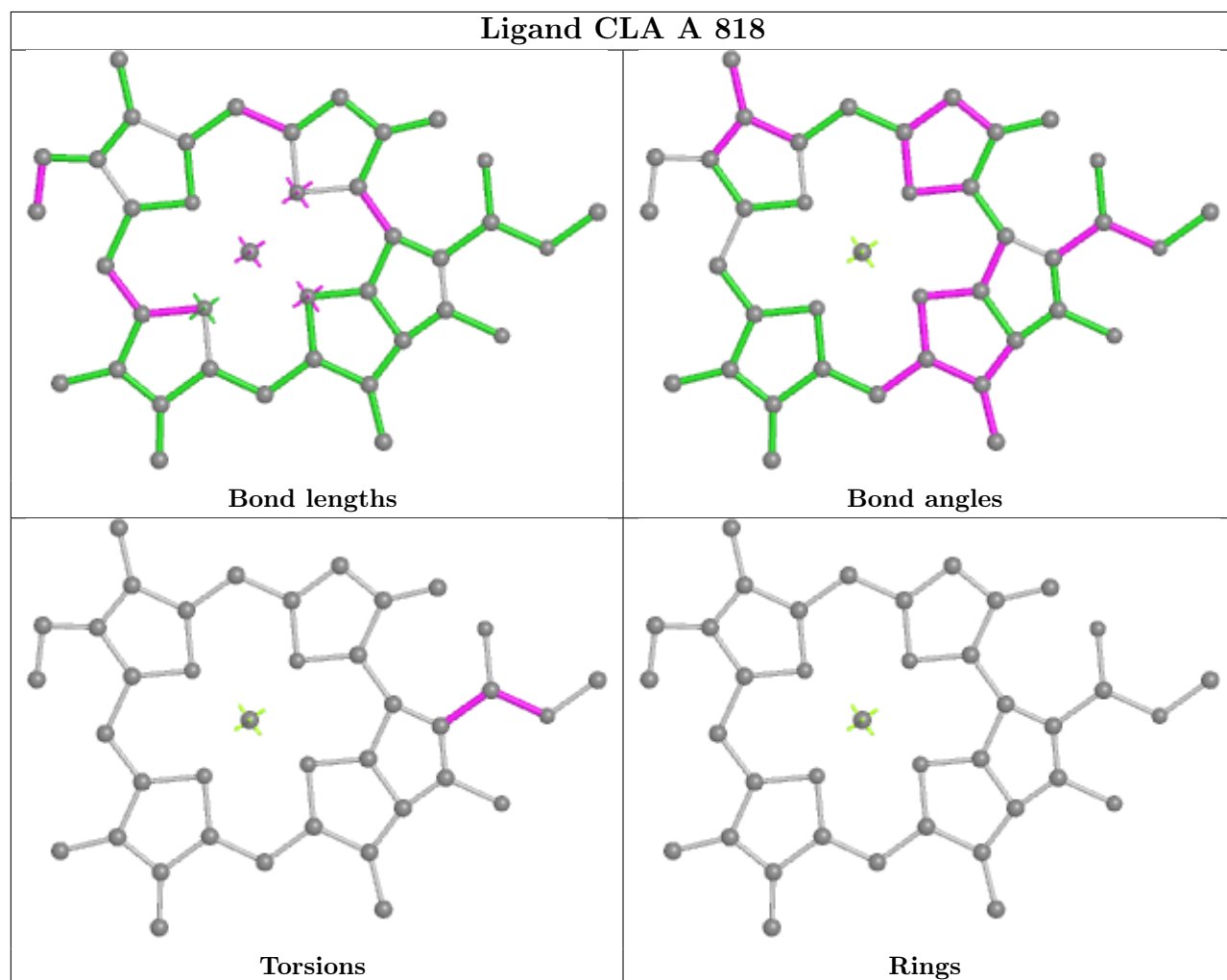
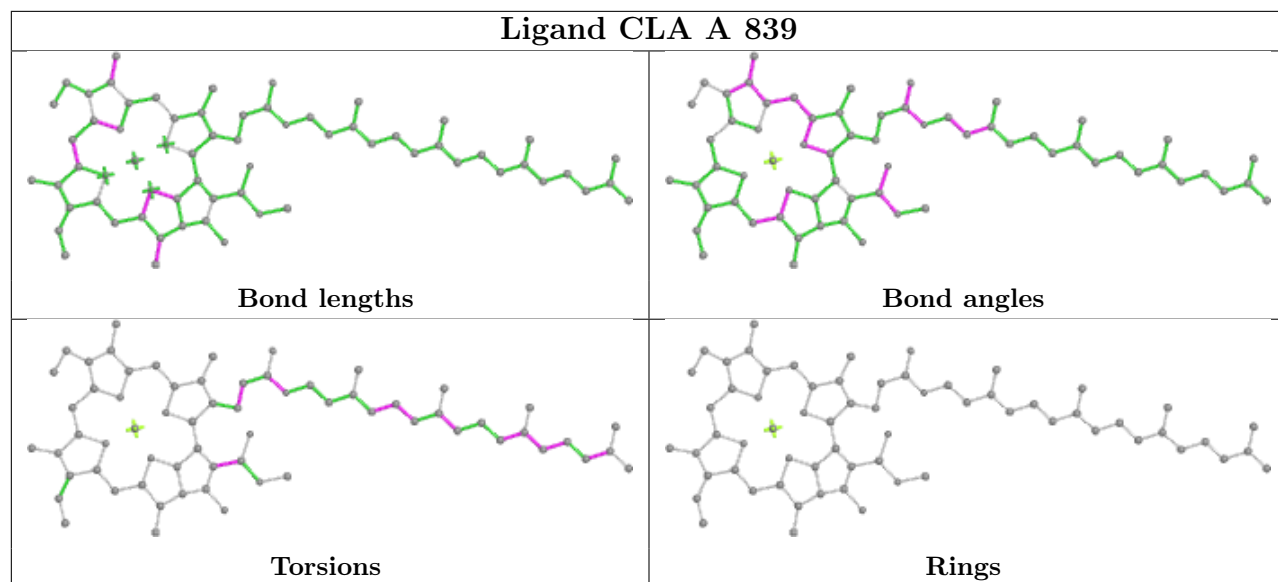


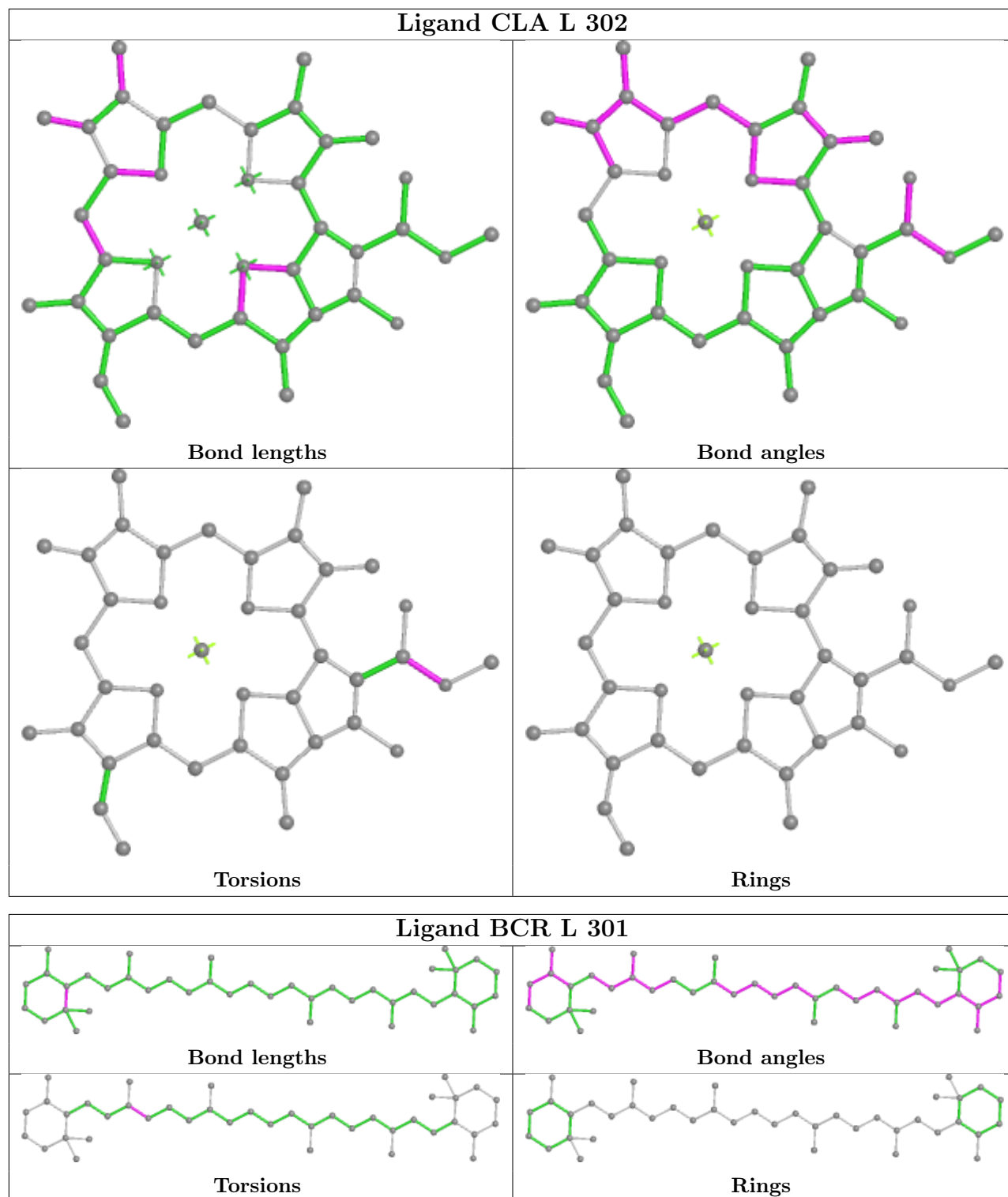


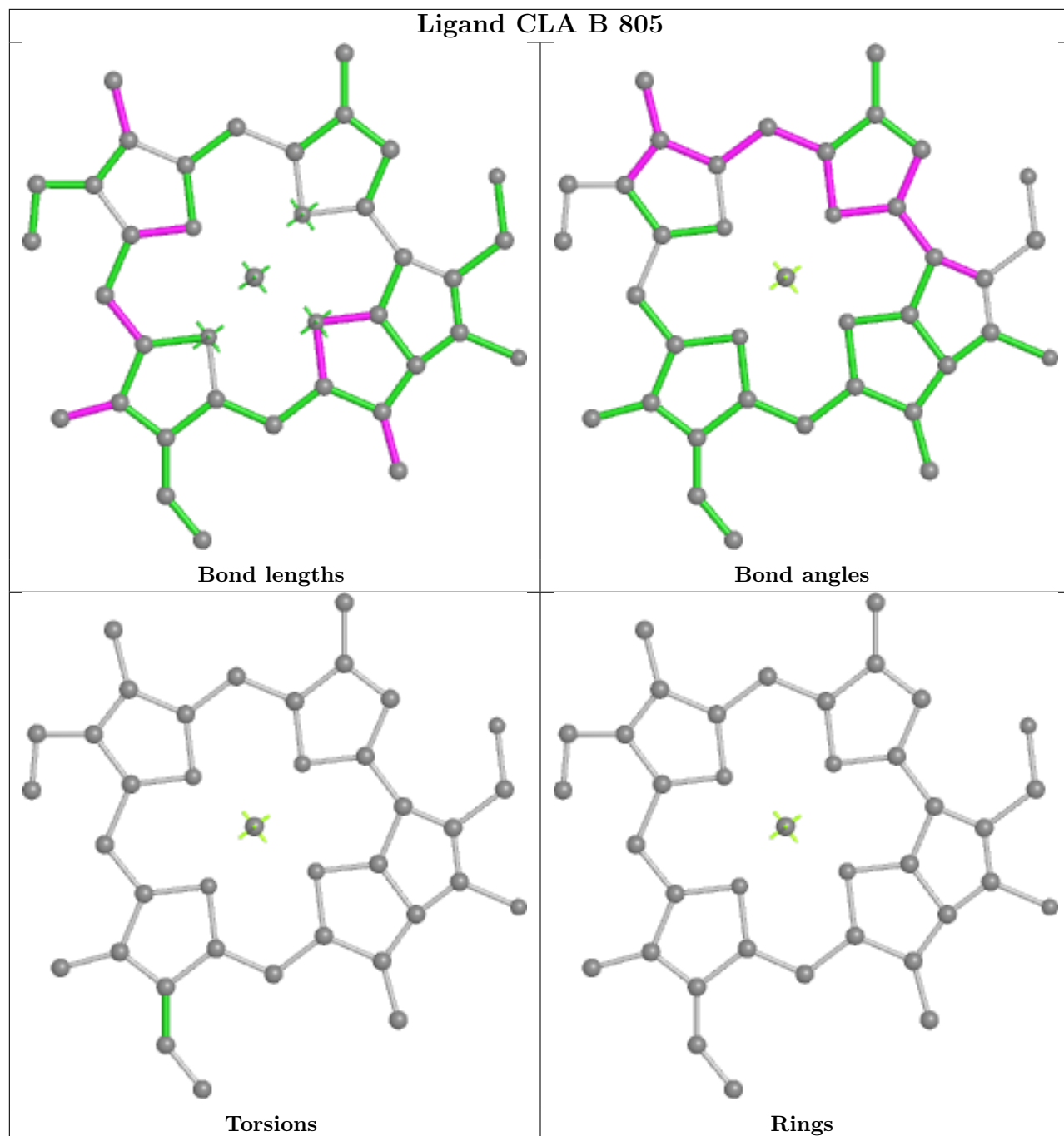


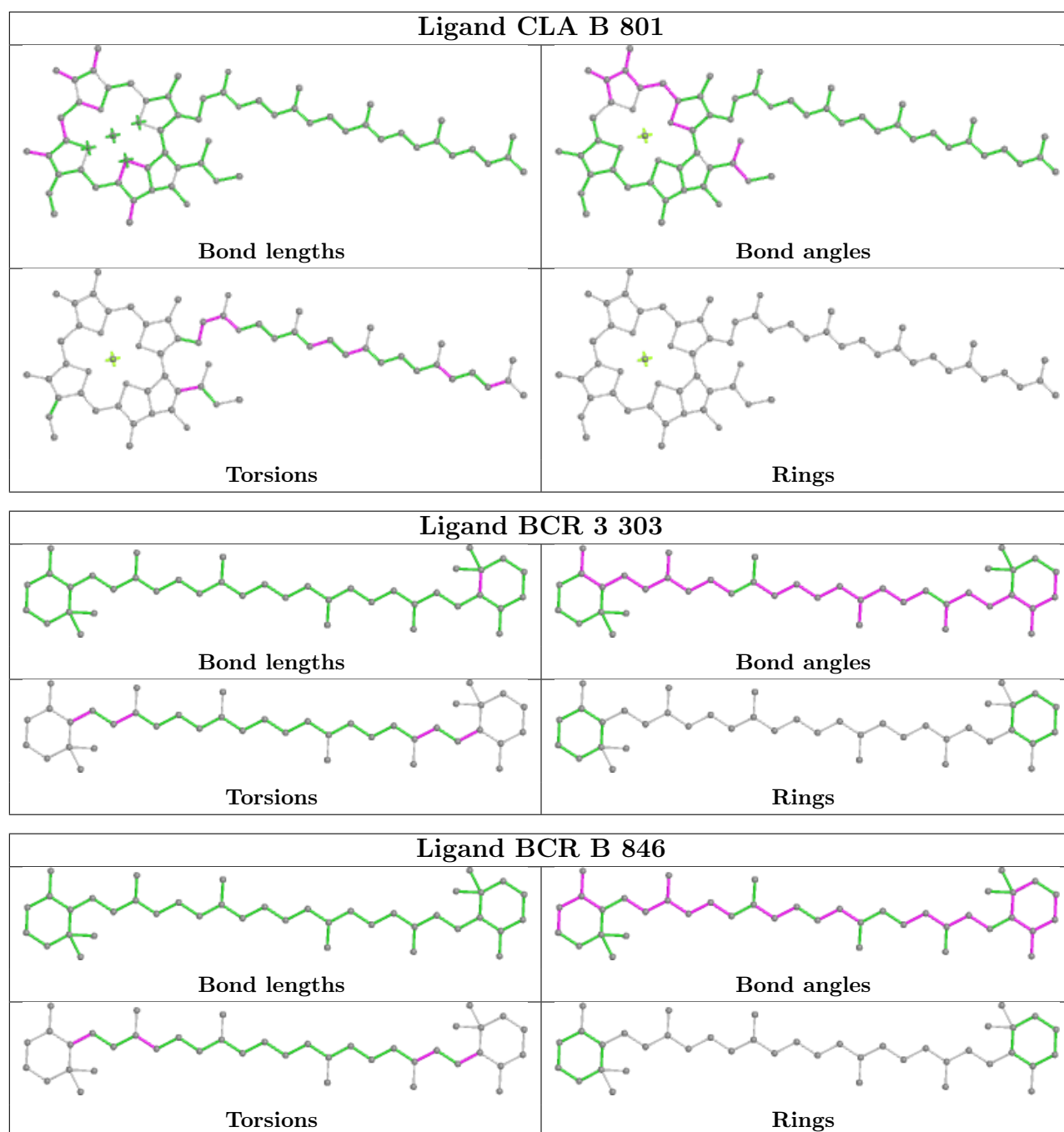


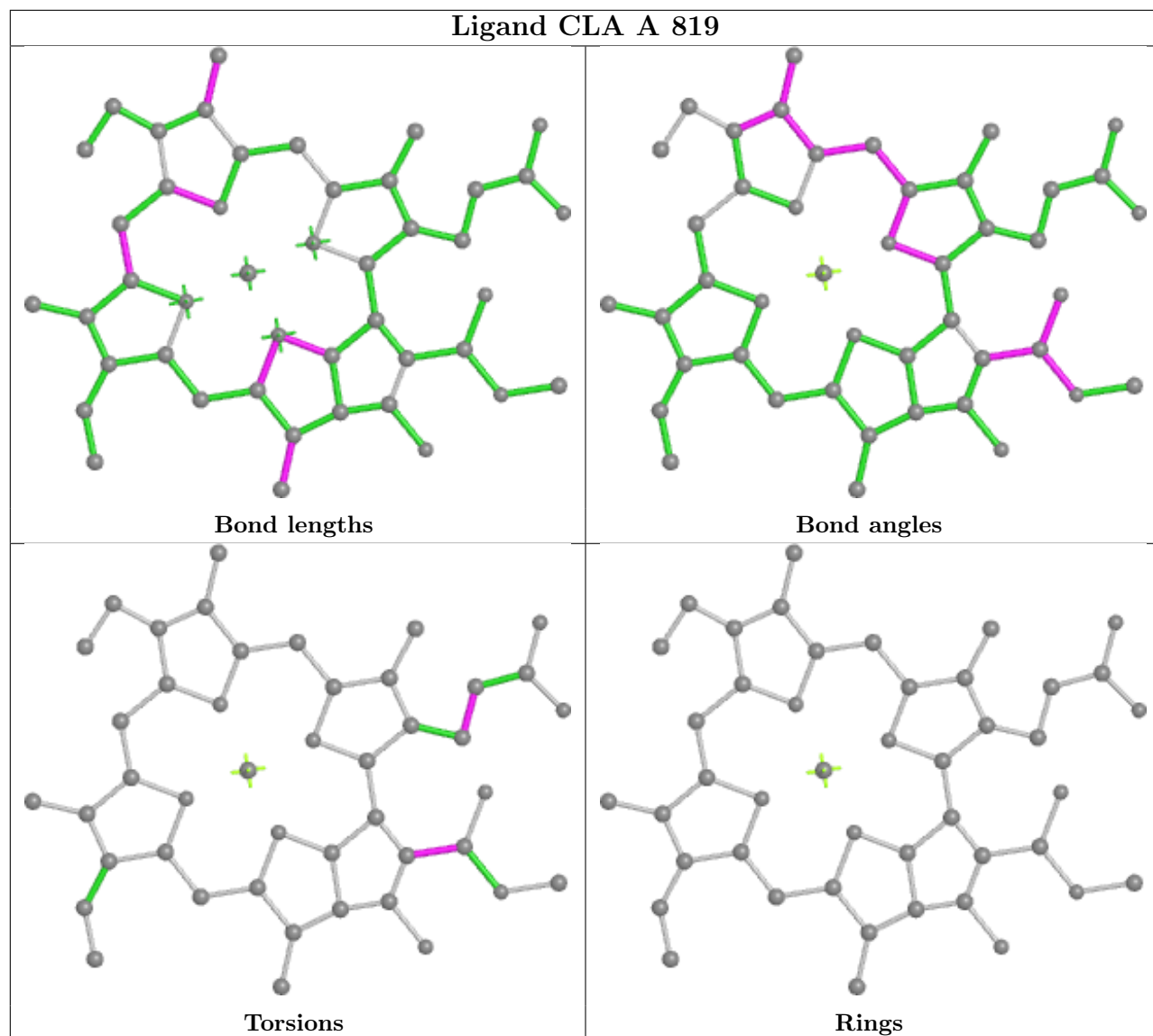


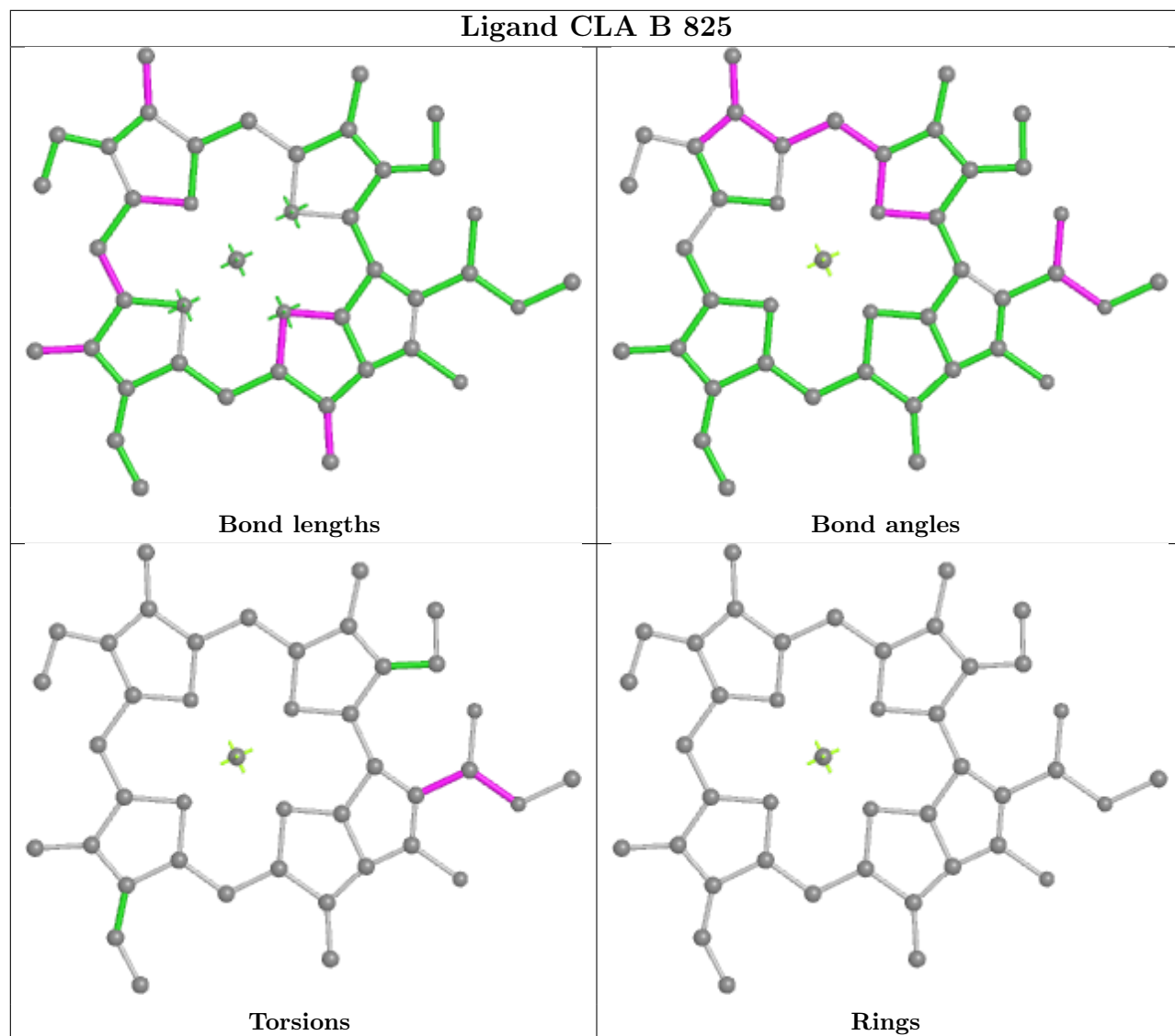


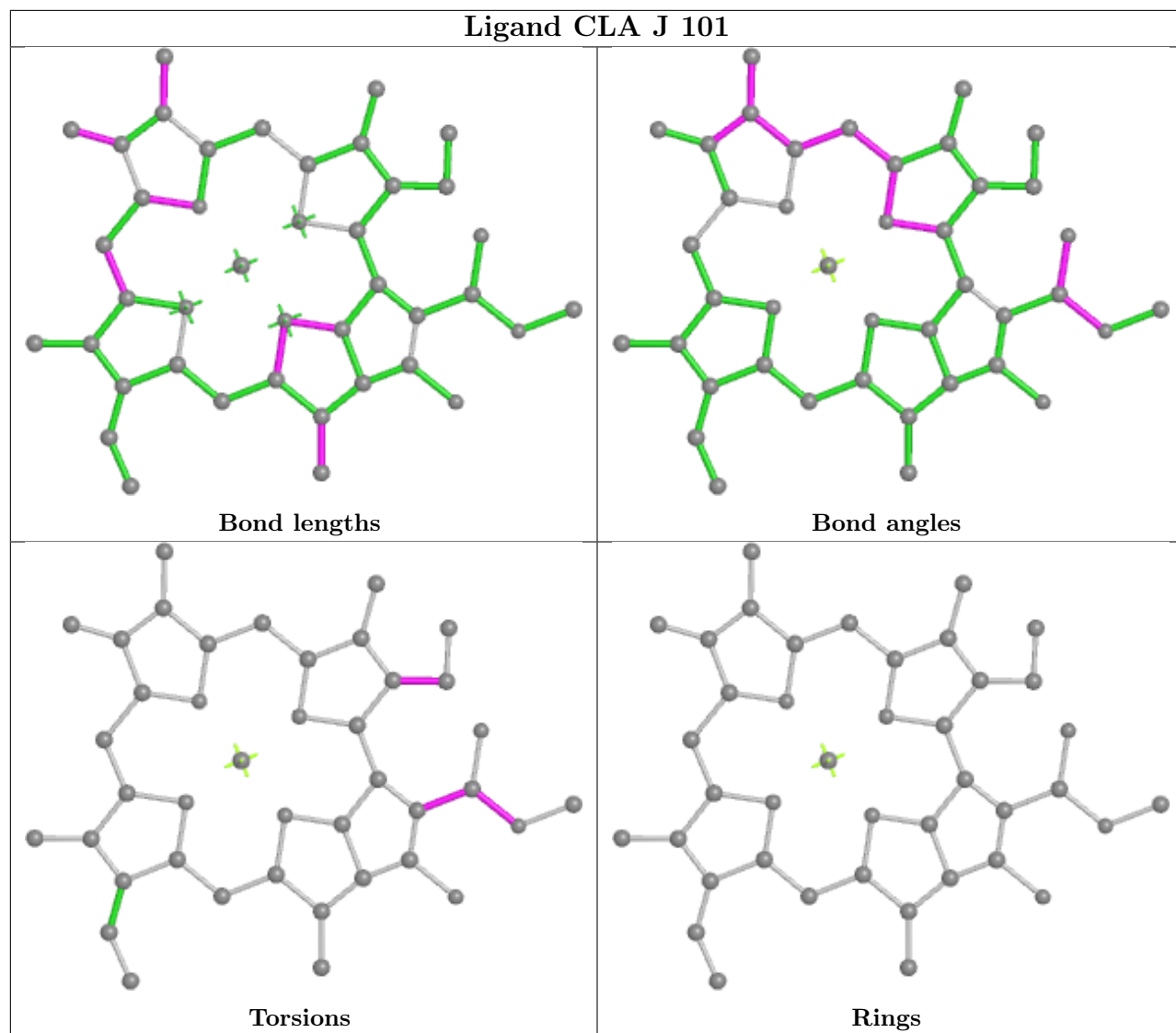


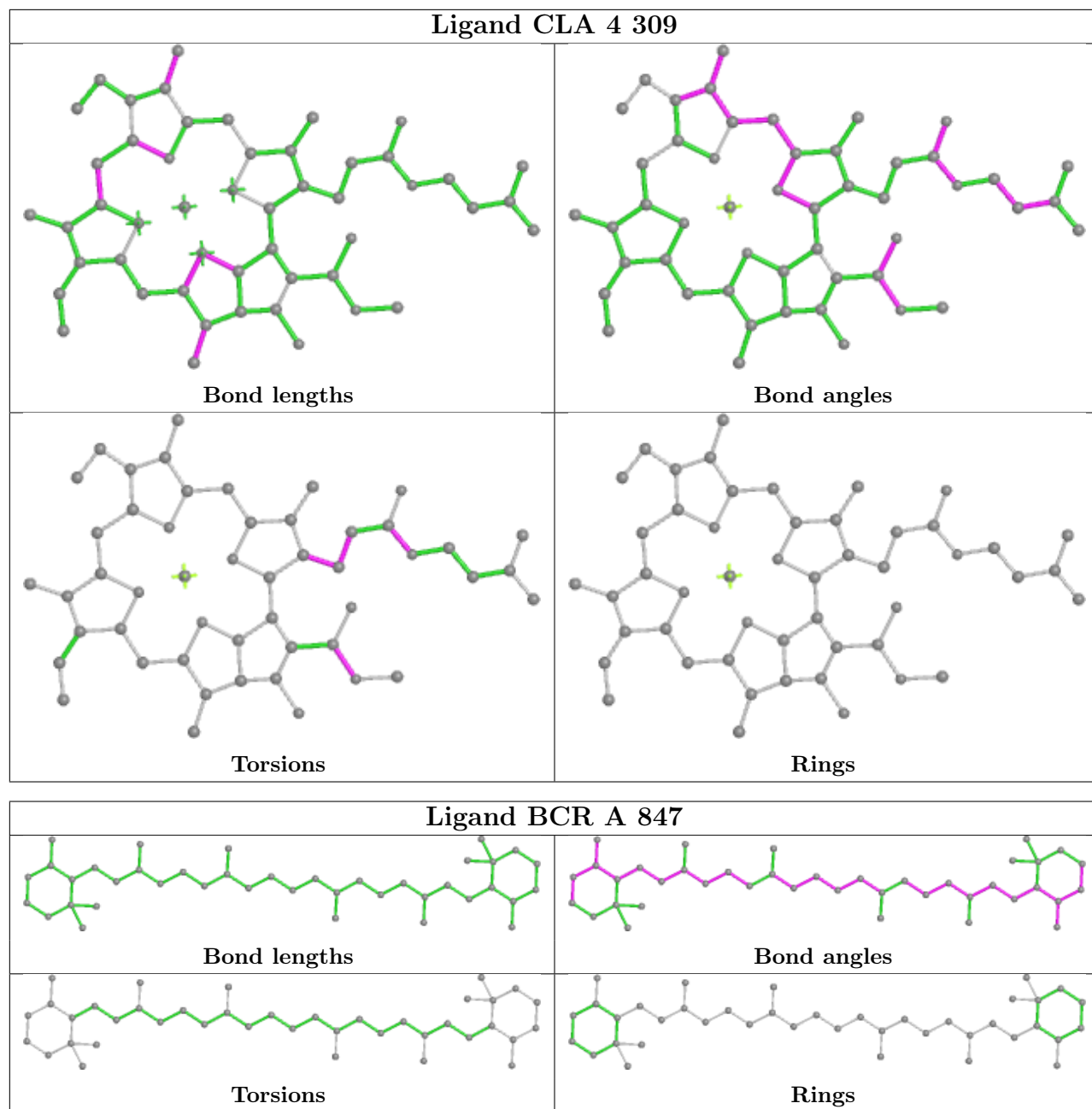


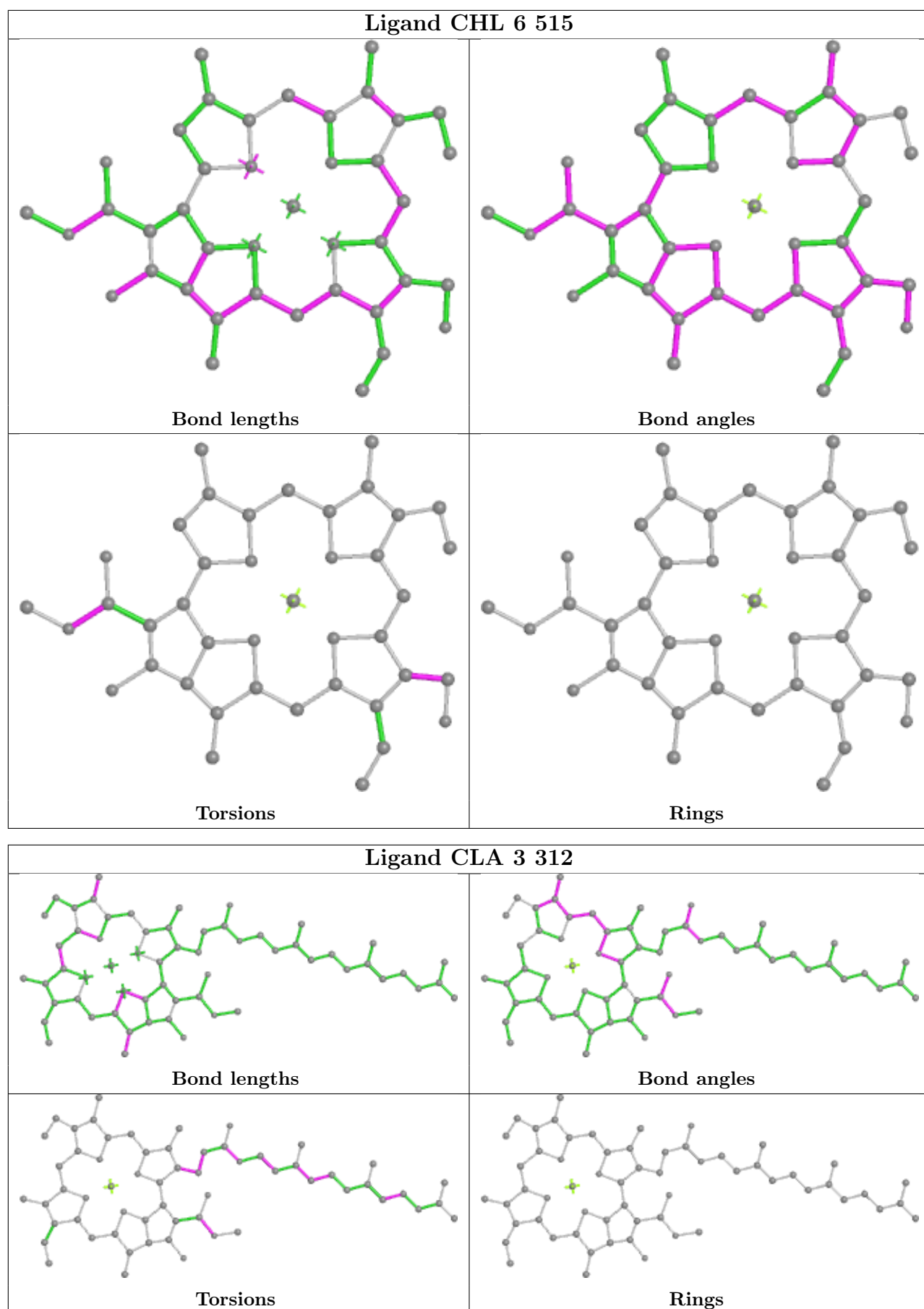


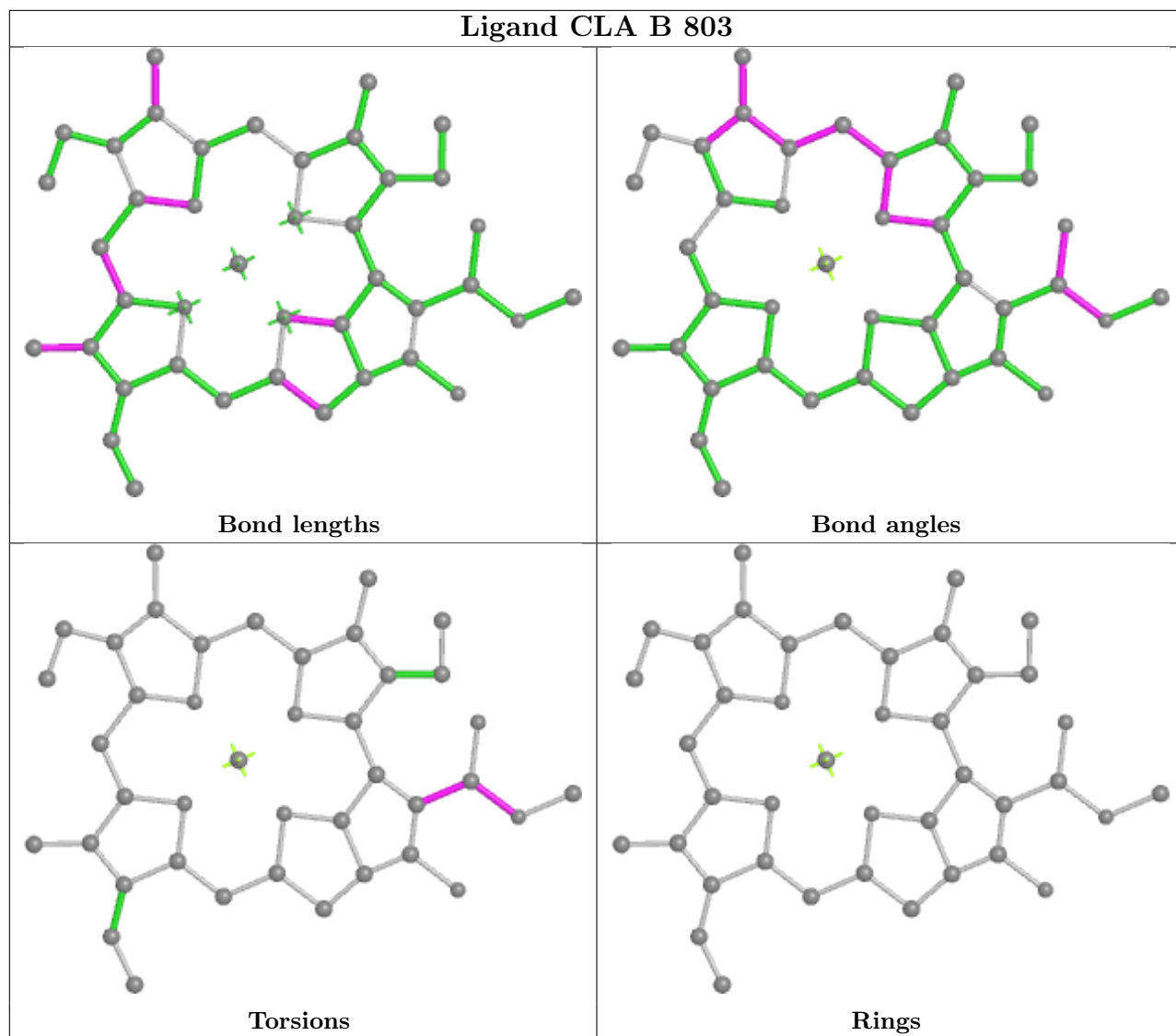


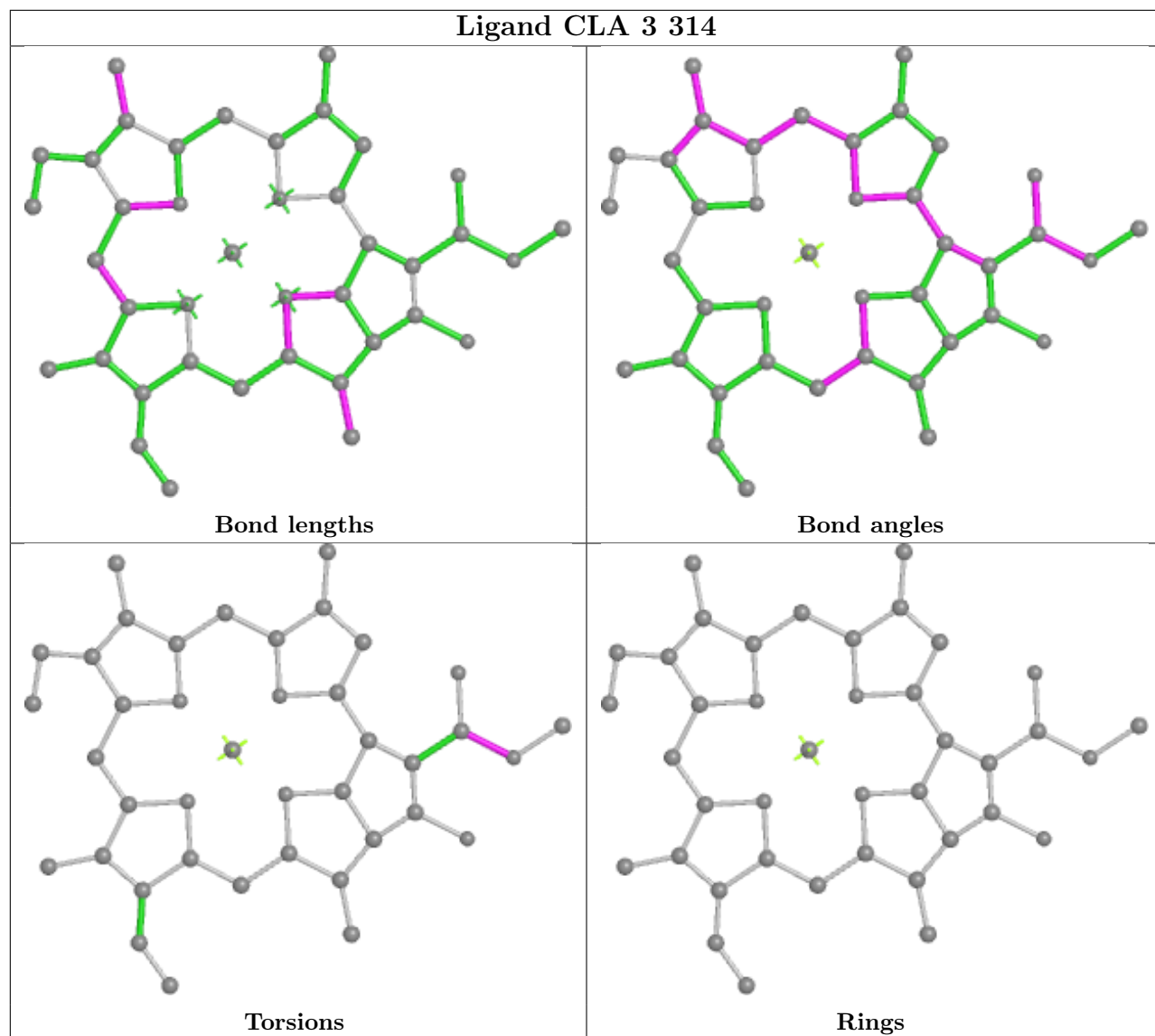


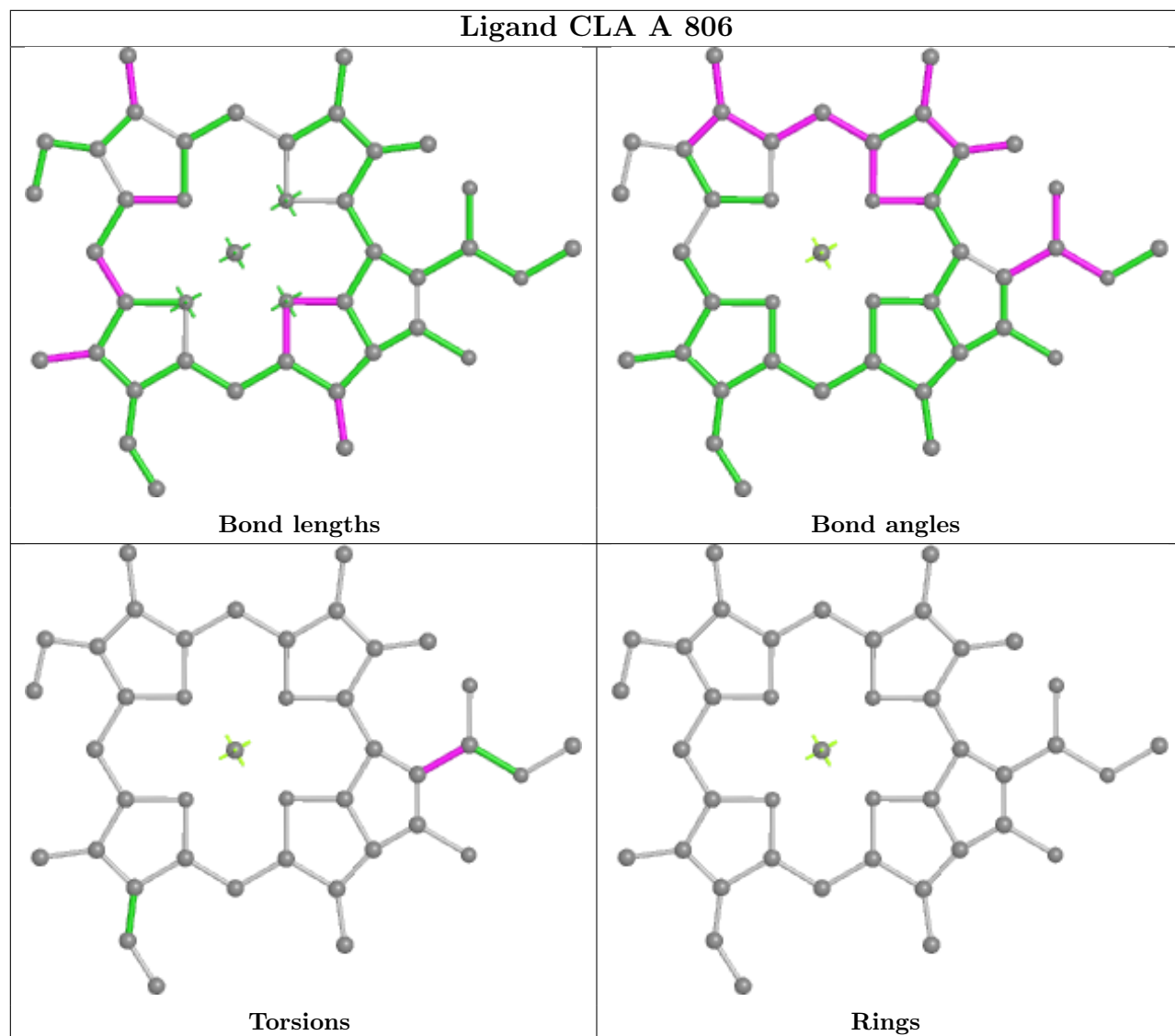


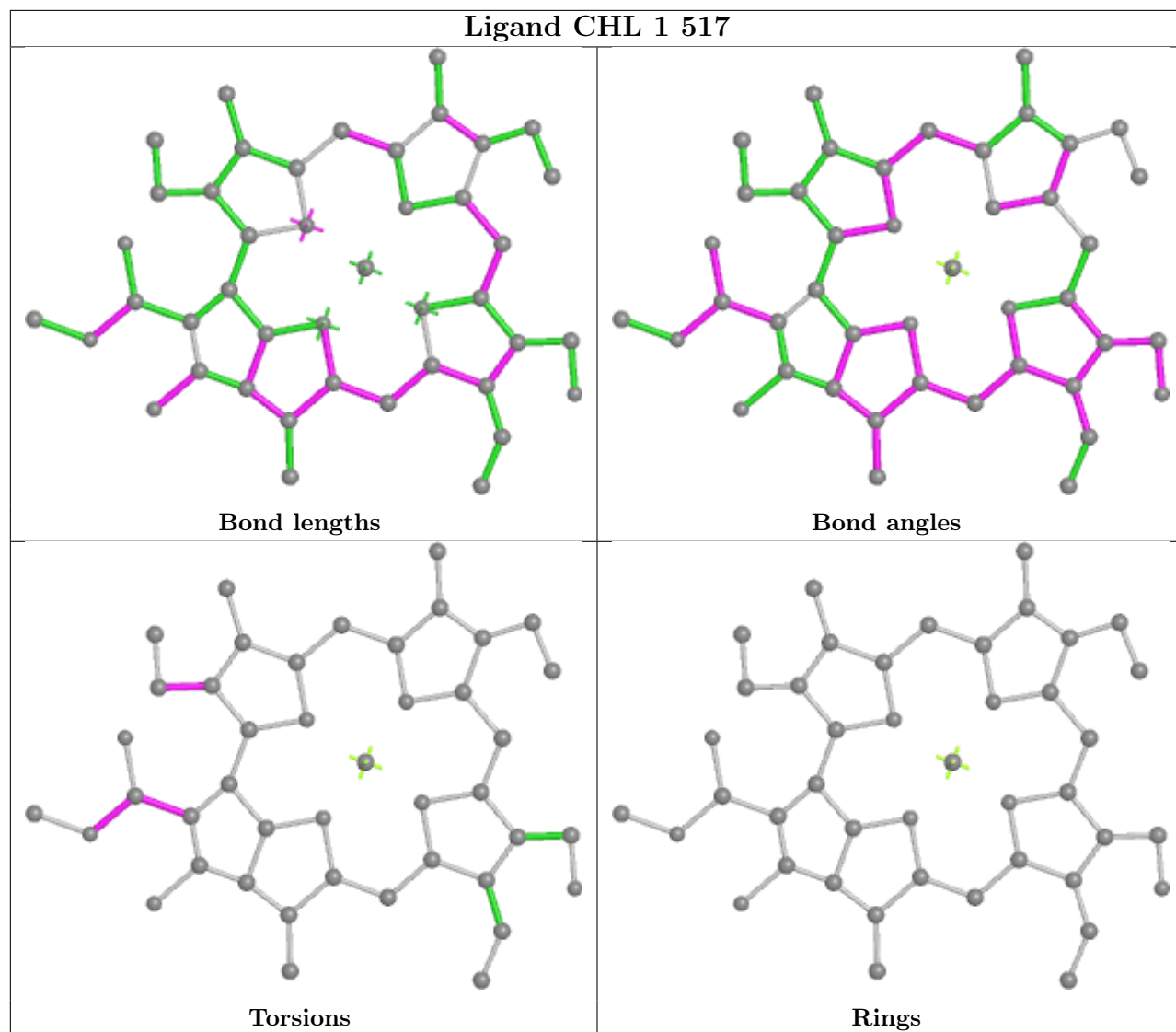


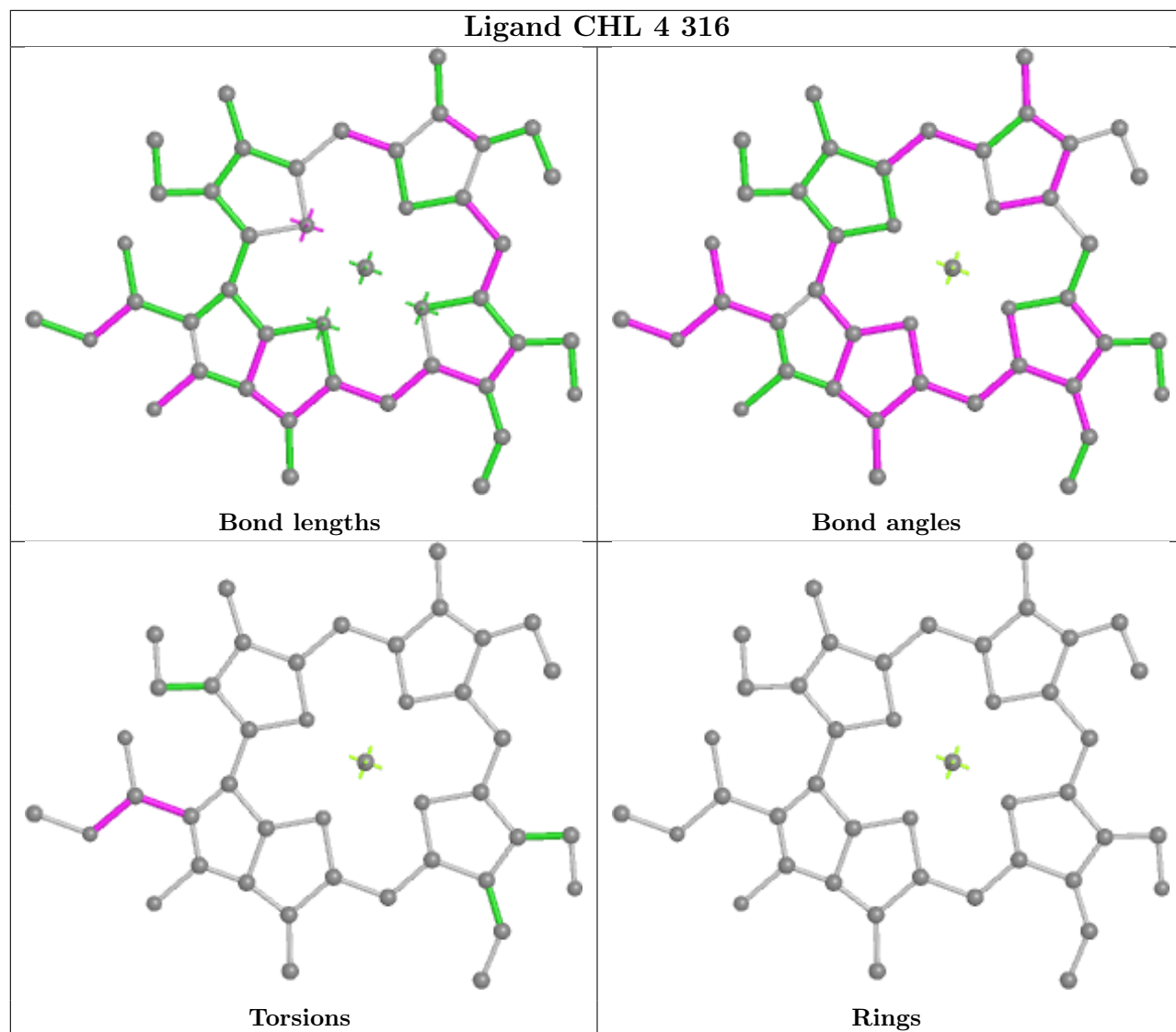


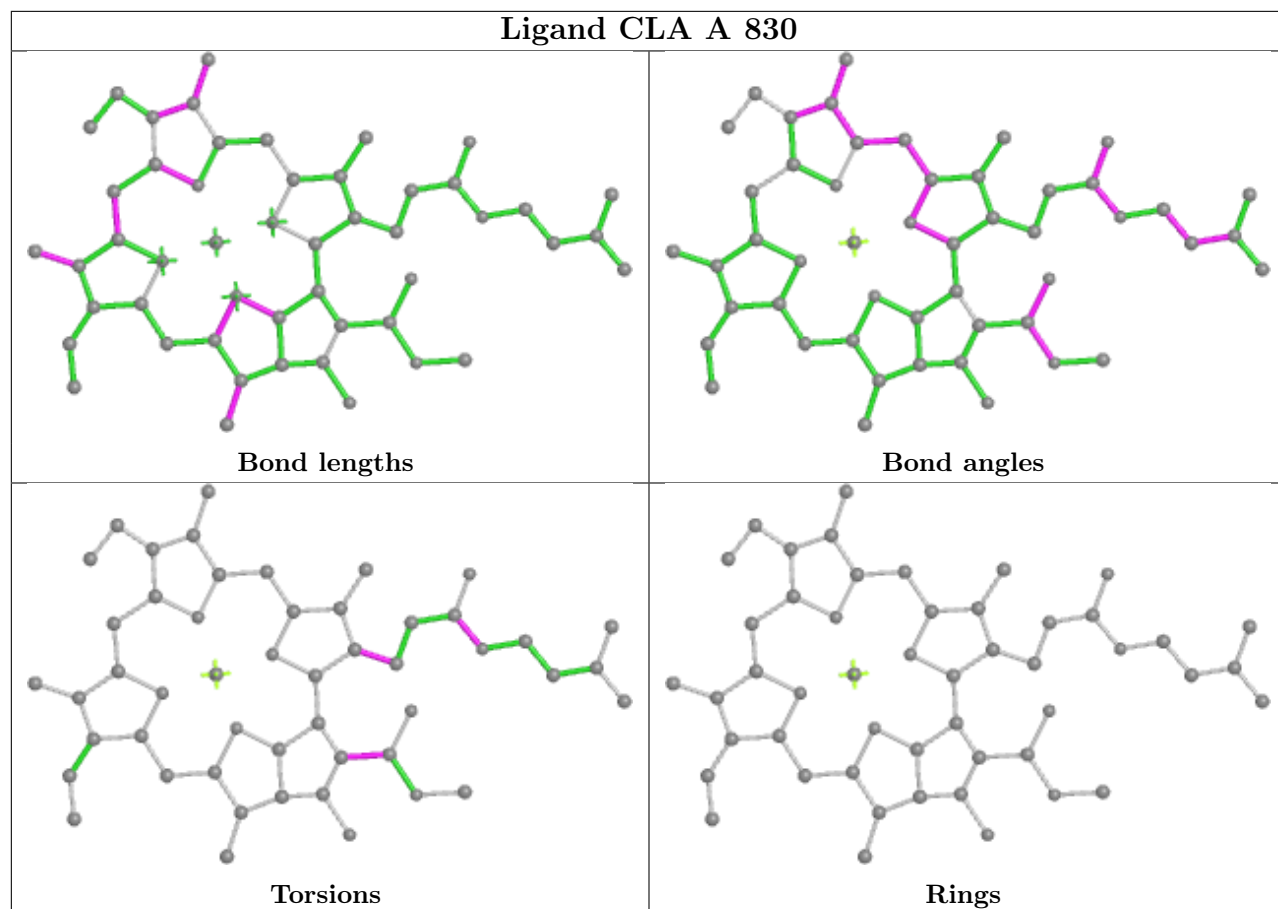


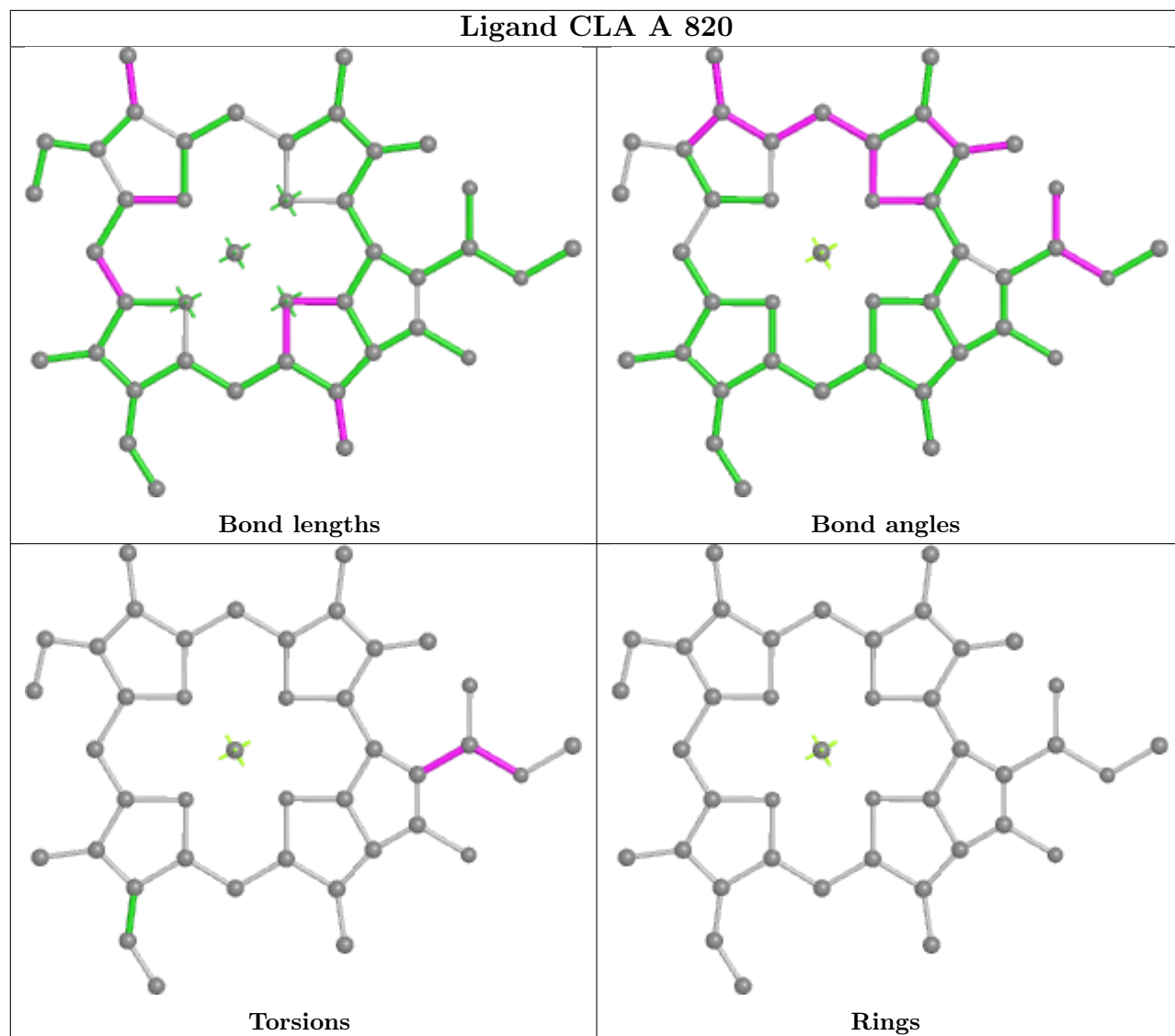


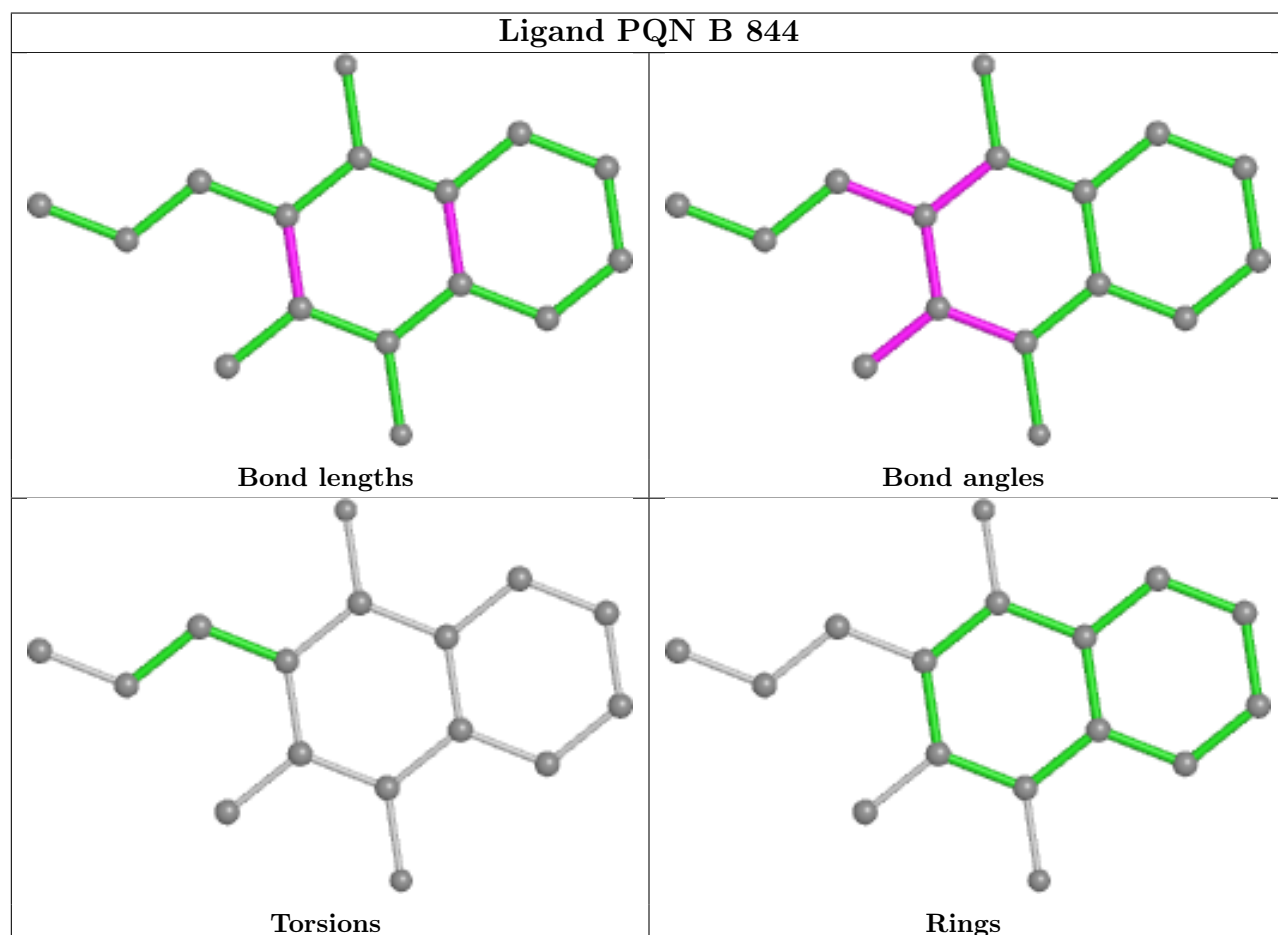
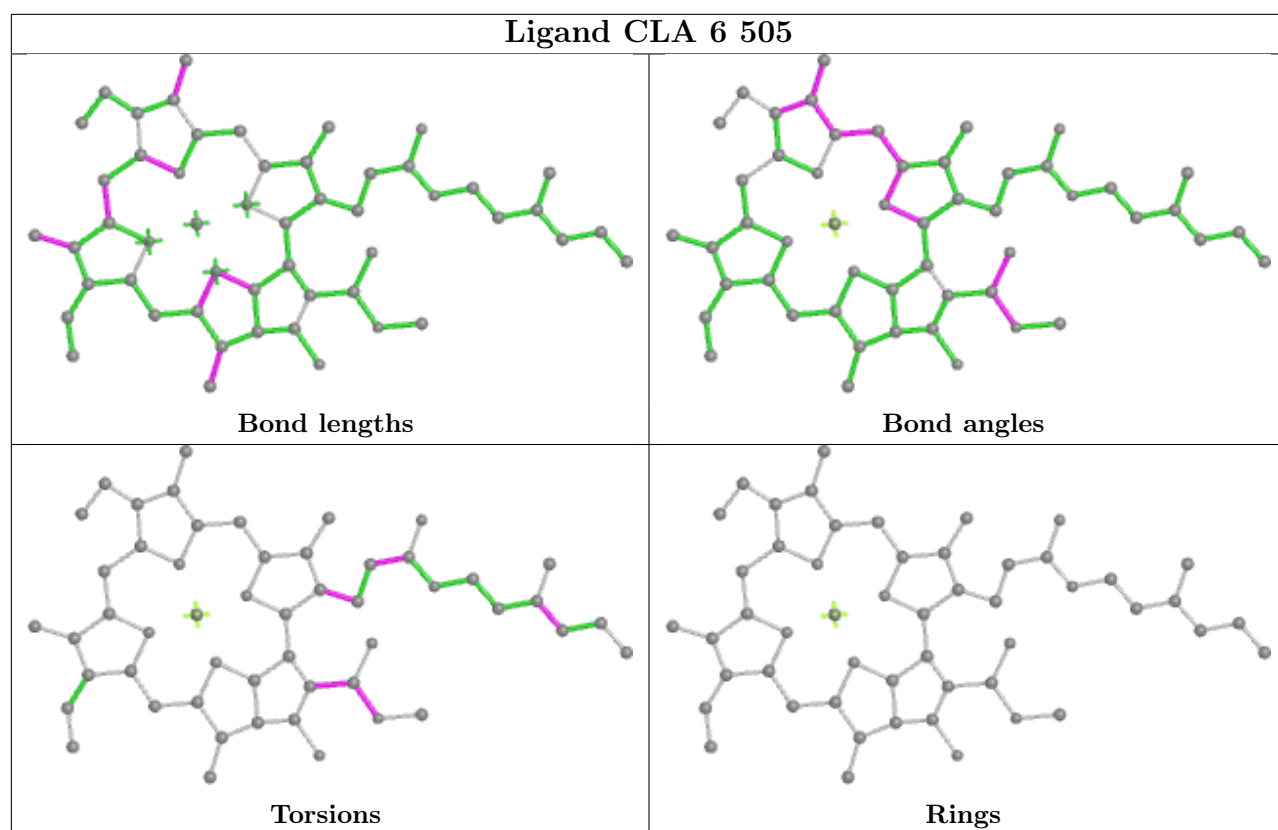


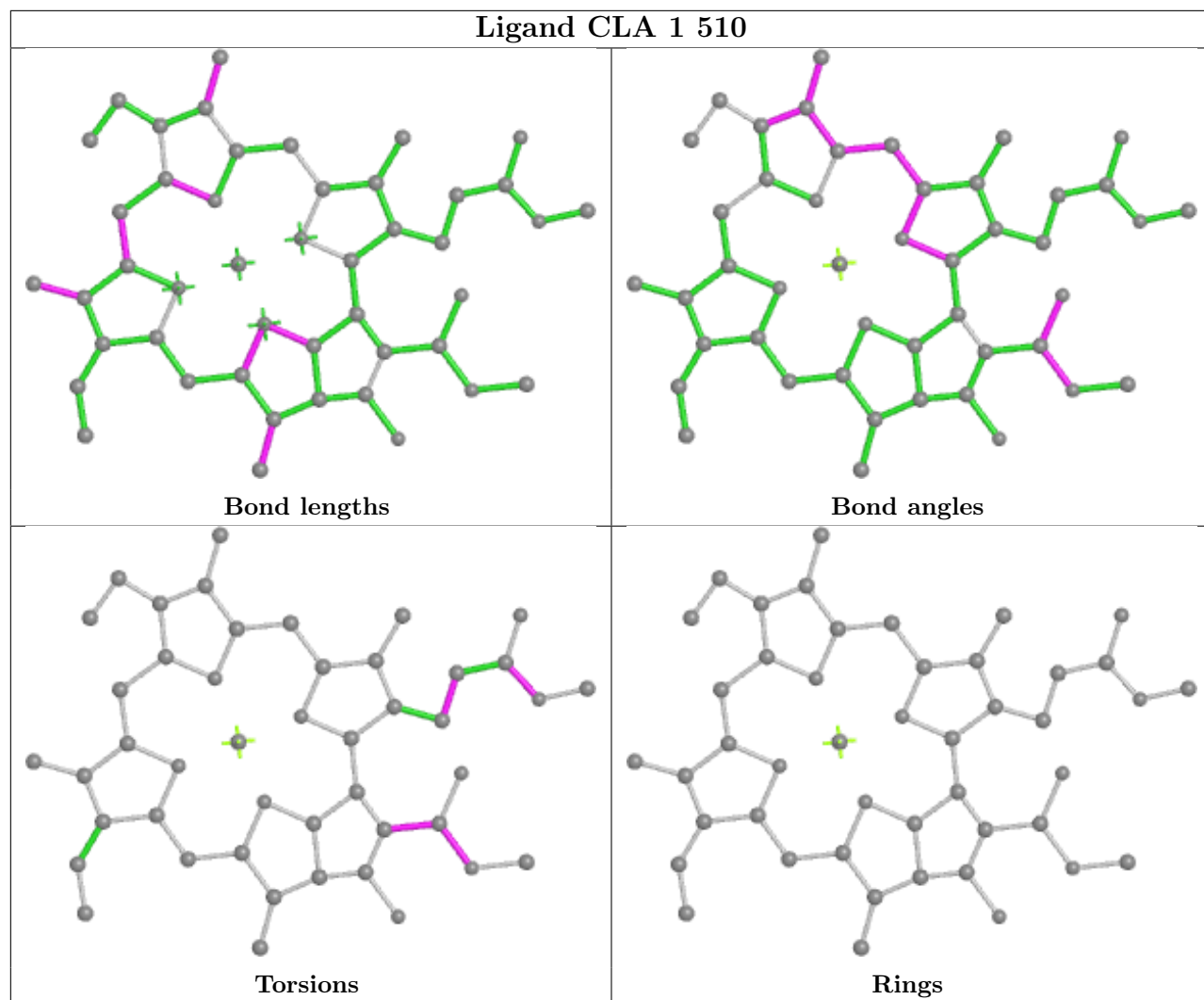


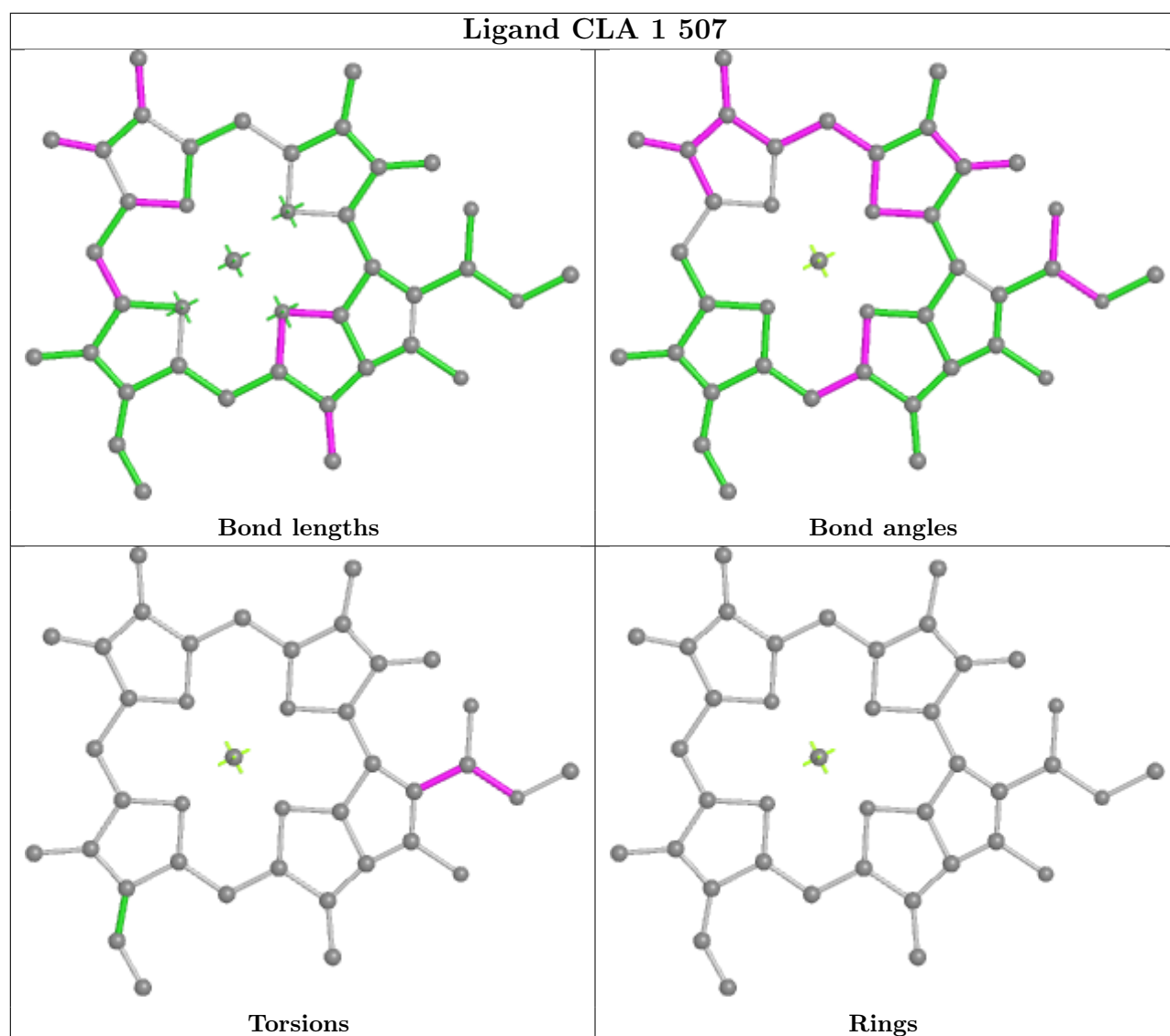












5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

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

There are no chain breaks in this entry.

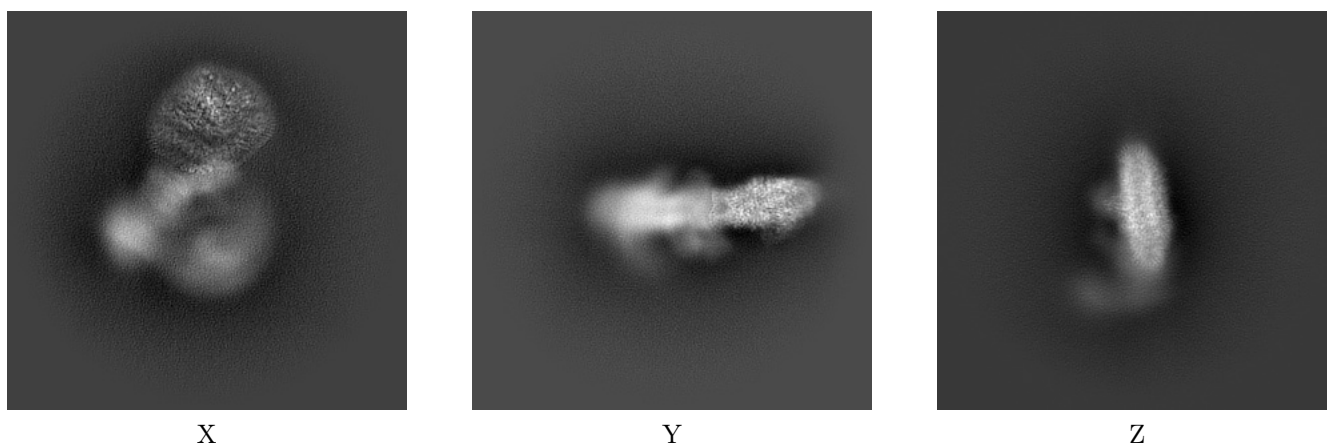
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-31350. These allow visual inspection of the internal detail of the map and identification of artifacts.

No raw map or half-maps were deposited for this entry and therefore no images, graphs, etc. pertaining to the raw map can be shown.

6.1 Orthogonal projections [i](#)

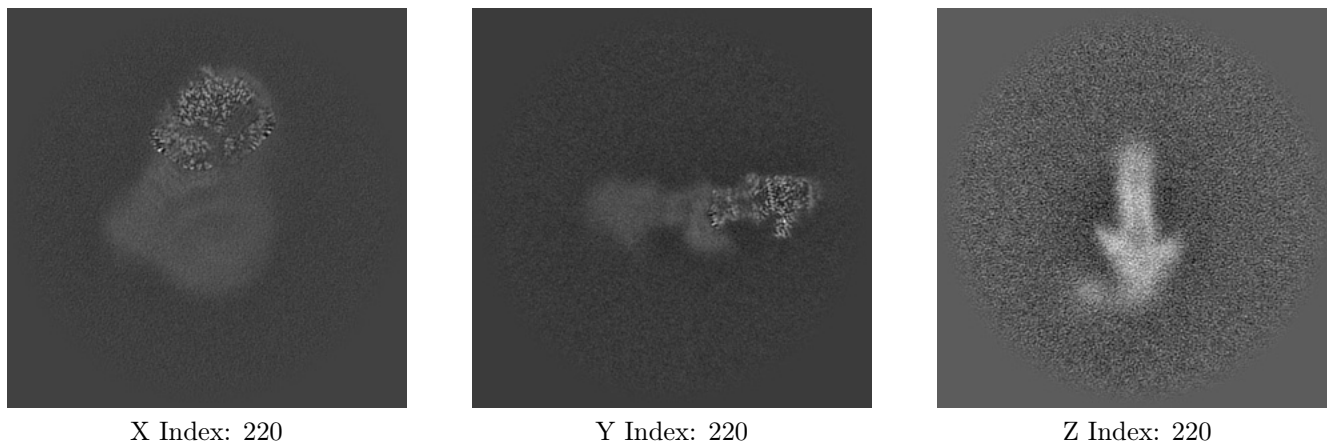
6.1.1 Primary map



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

6.2 Central slices [i](#)

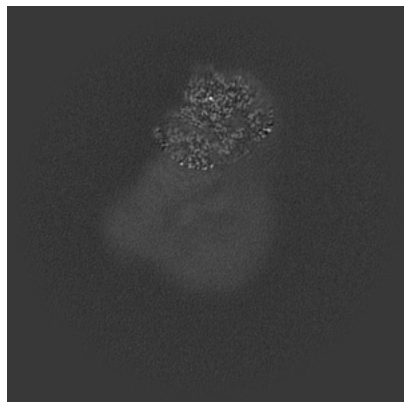
6.2.1 Primary map



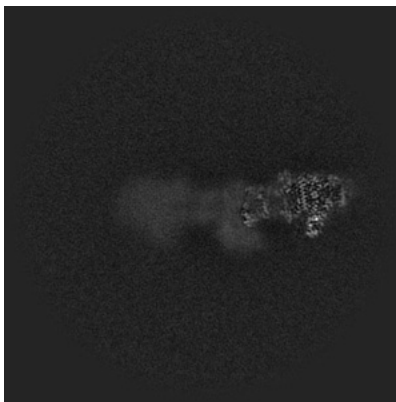
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

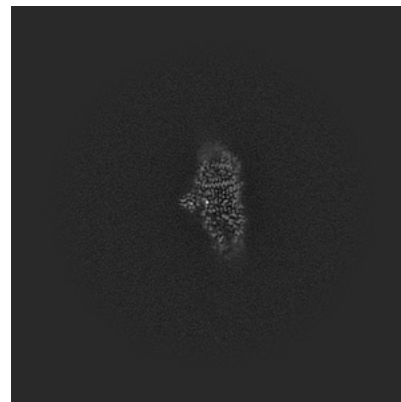
6.3.1 Primary map



X Index: 215



Y Index: 221

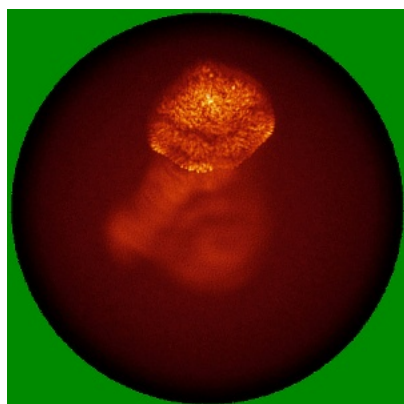


Z Index: 338

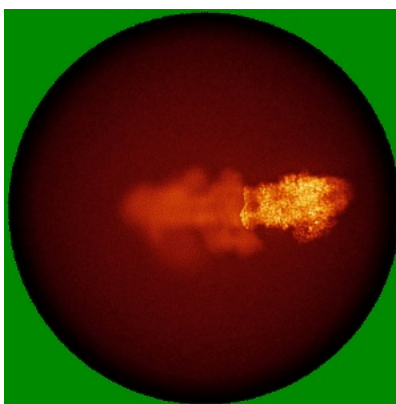
The images above show the largest variance slices of the map in three orthogonal directions.

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

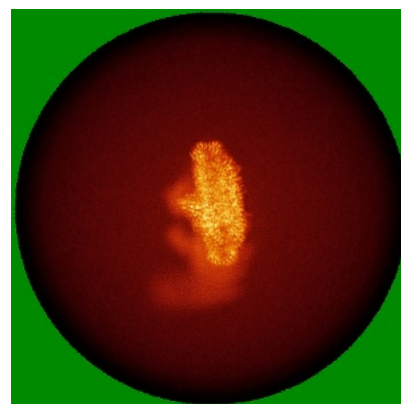
6.4.1 Primary map



X



Y

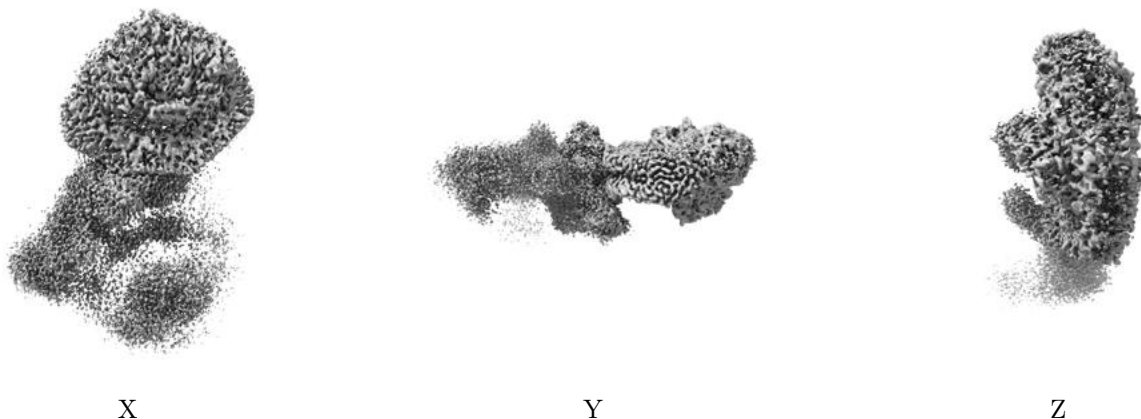


Z

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 0.43. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

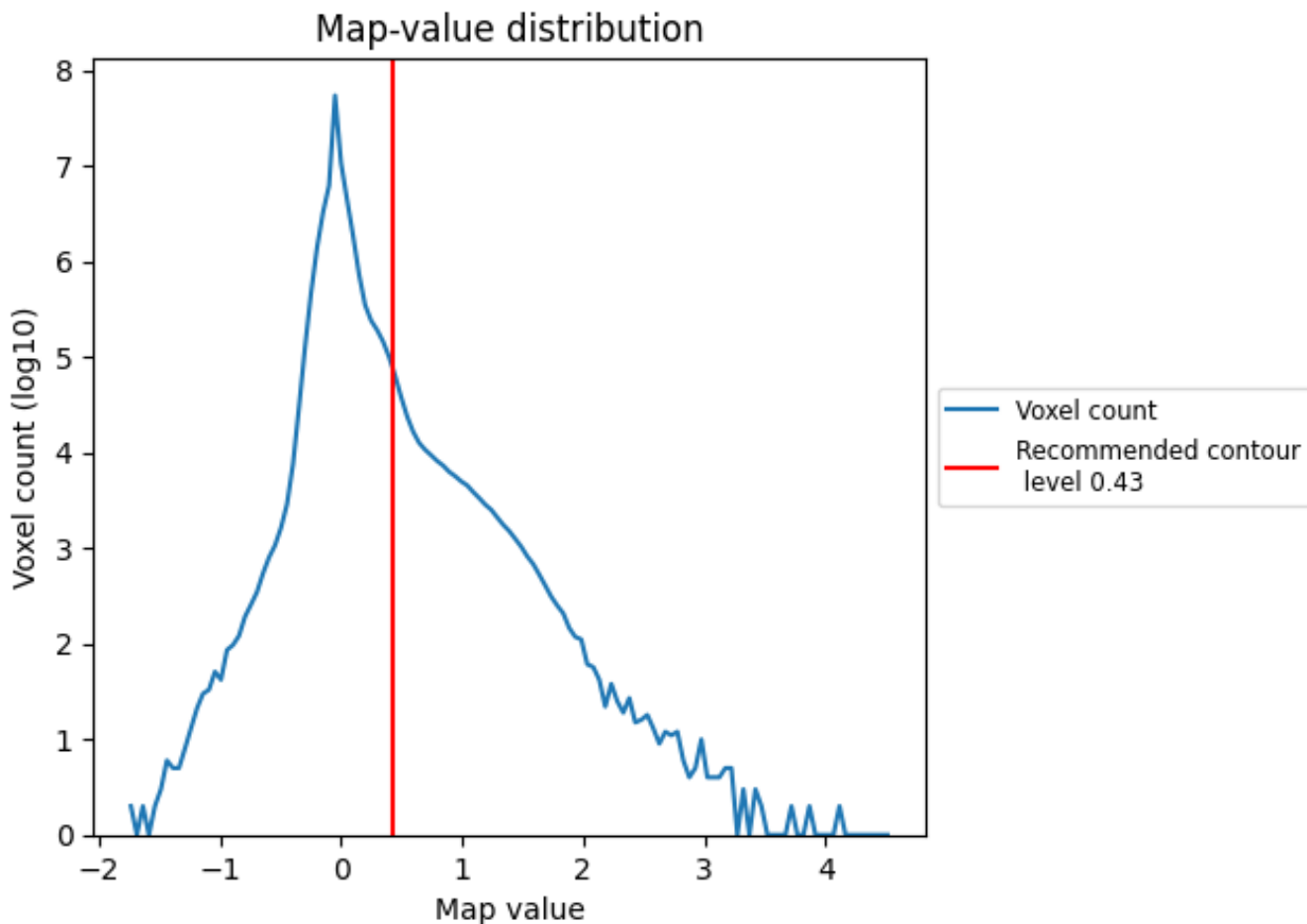
6.6 Mask visualisation [i](#)

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

7 Map analysis [i](#)

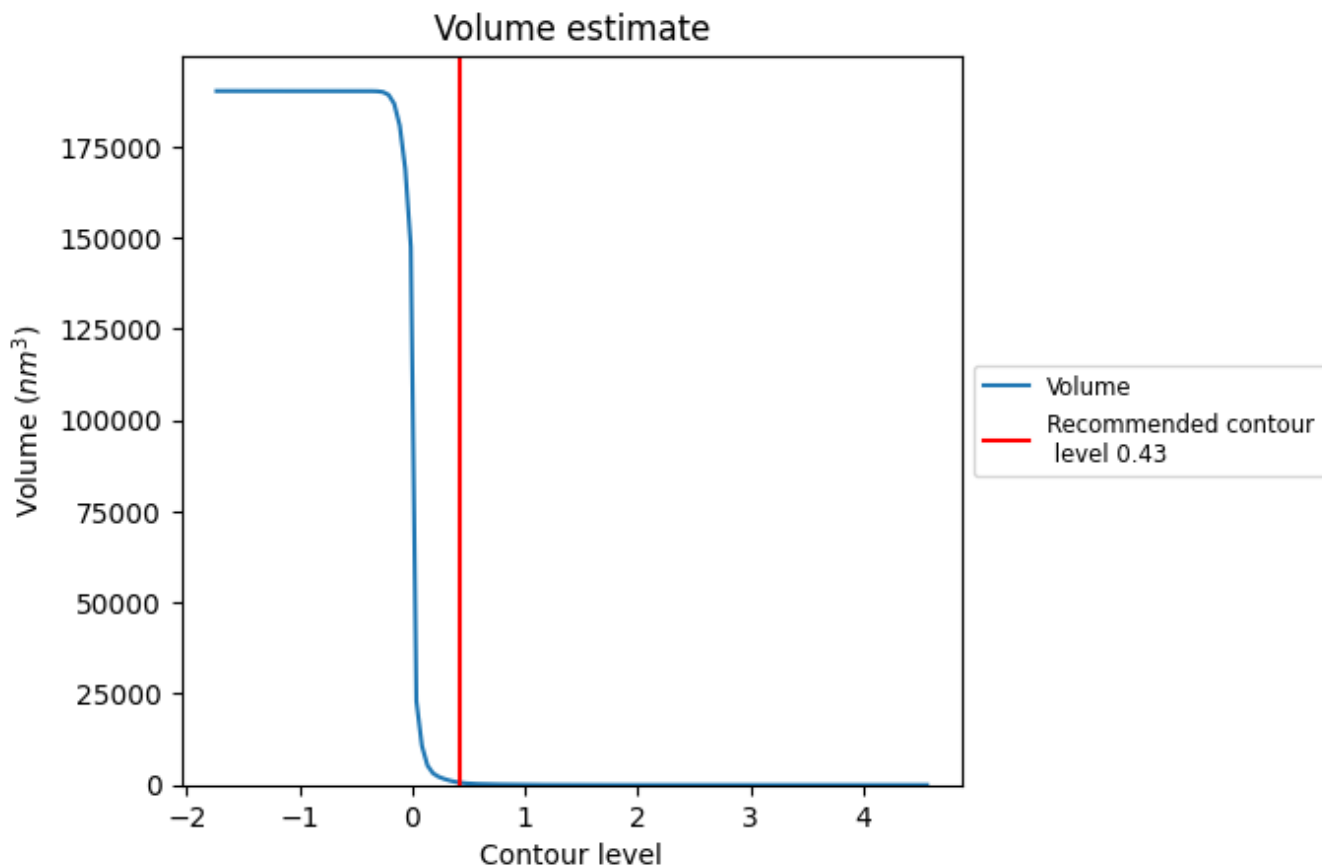
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

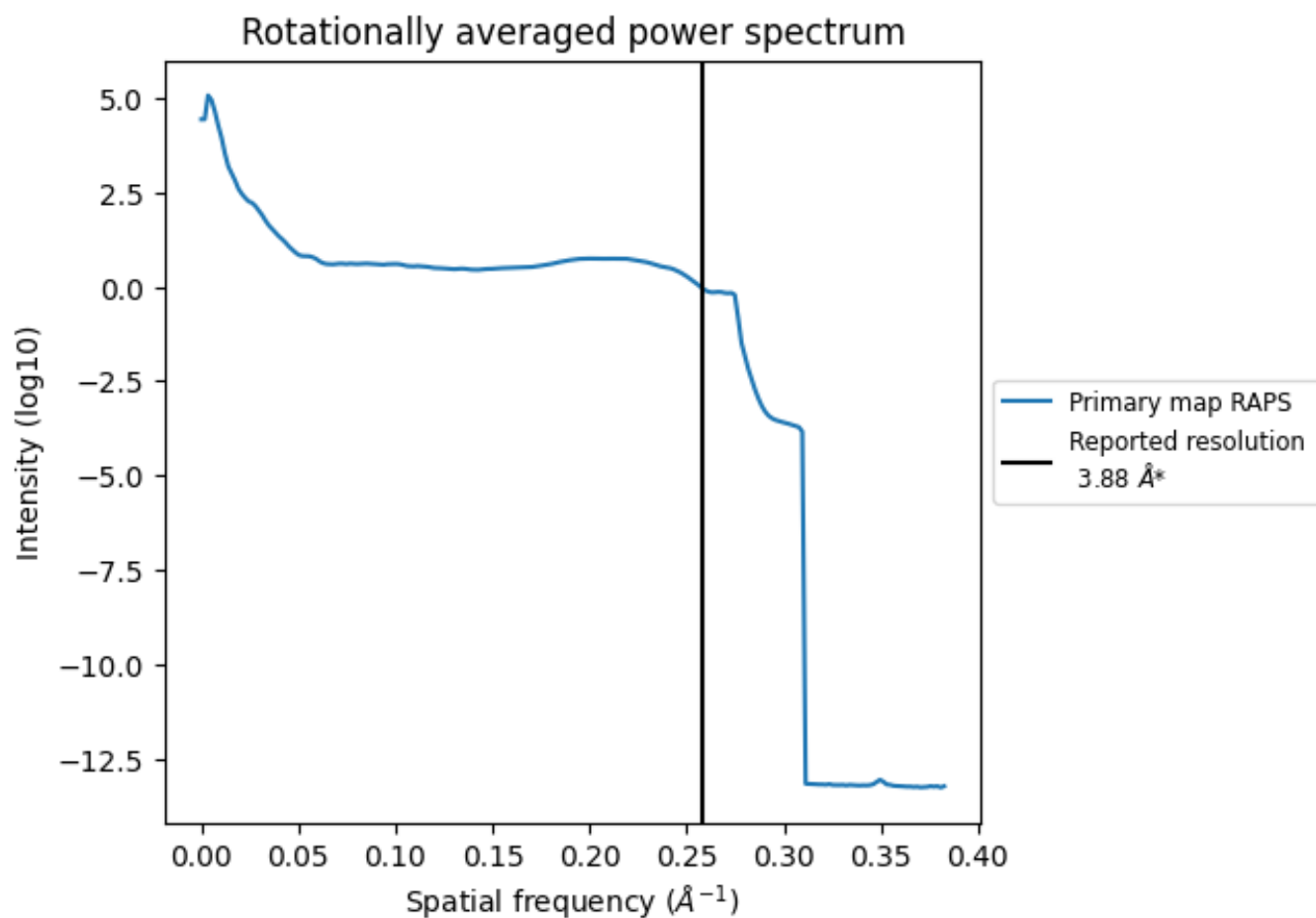
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 598 nm^3 ; this corresponds to an approximate mass of 540 kDa.

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

7.3 Rotationally averaged power spectrum [i](#)



*Reported resolution corresponds to spatial frequency of 0.258 Å⁻¹

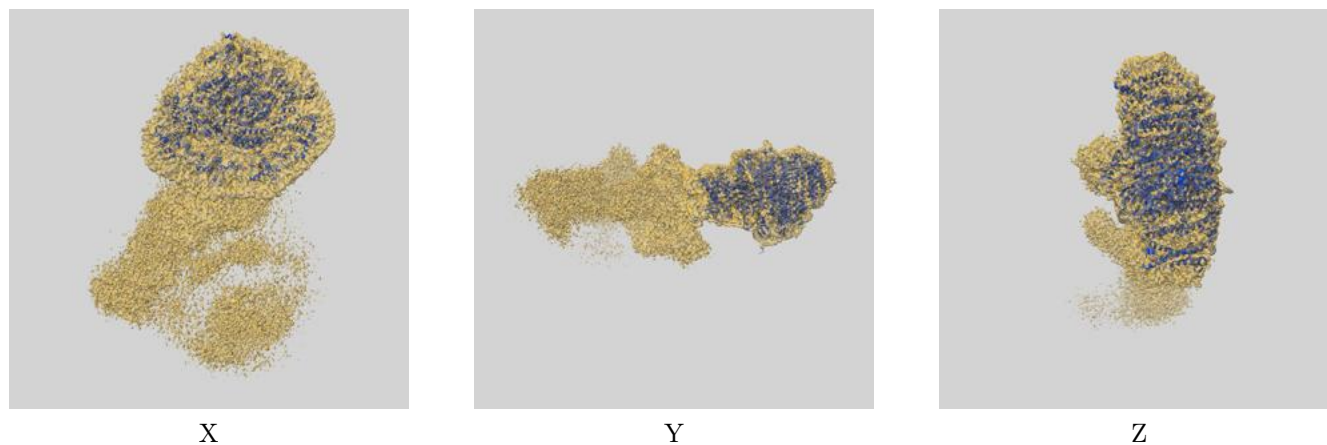
8 Fourier-Shell correlation

This section was not generated. No FSC curve or half-maps provided.

9 Map-model fit [i](#)

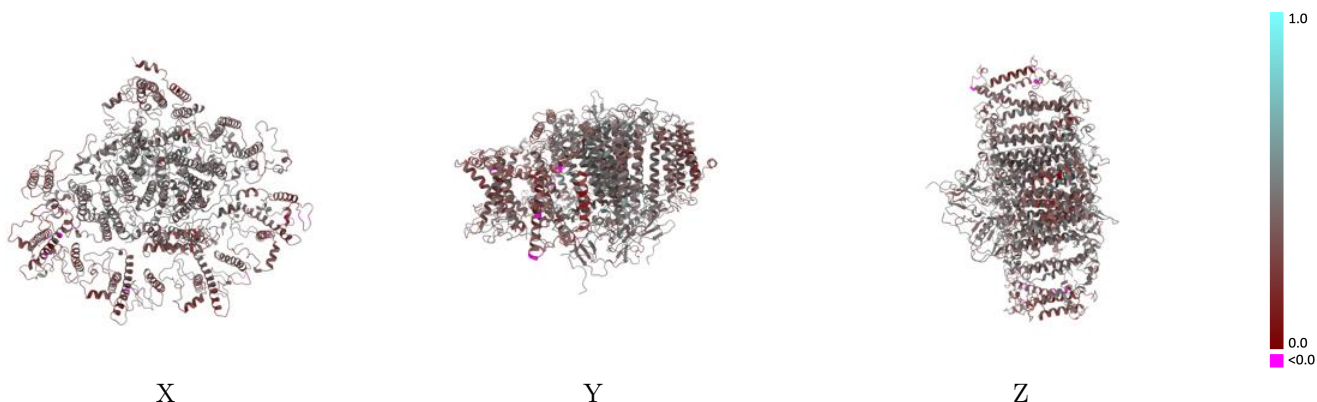
This section contains information regarding the fit between EMDB map EMD-31350 and PDB model 7EWK. Per-residue inclusion information can be found in section [3](#) on page [25](#).

9.1 Map-model overlay [i](#)



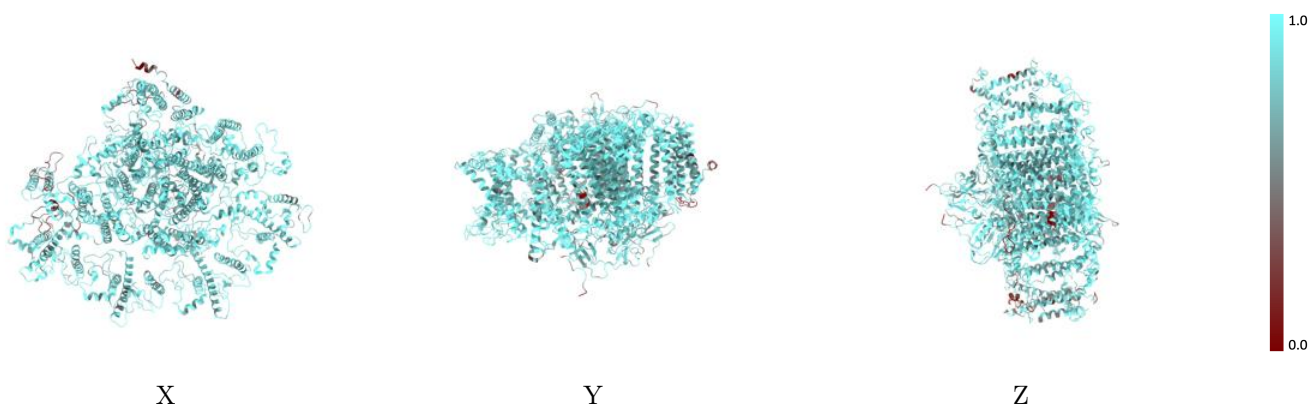
The images above show the 3D surface view of the map at the recommended contour level 0.43 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



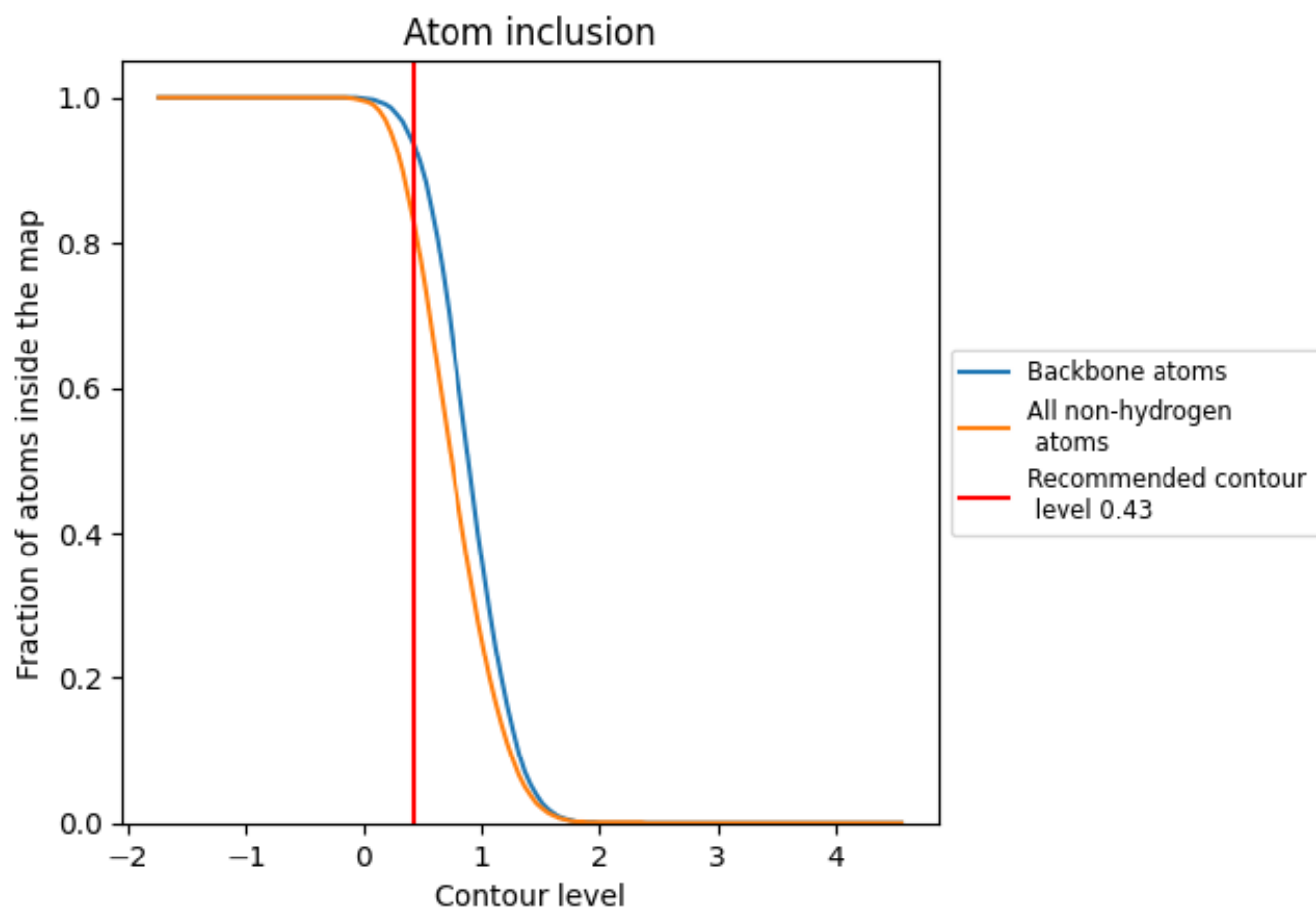
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.43).

































9.4 Atom inclusion [i](#)



At the recommended contour level, 93% of all backbone atoms, 82% 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 (0.43) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	 0.8240	 0.3960
1	 0.8130	 0.3100
3	 0.7240	 0.2960
4	 0.8440	 0.3590
6	 0.8200	 0.3440
A	 0.8500	 0.4450
B	 0.8690	 0.4350
C	 0.9140	 0.4250
D	 0.8180	 0.4190
E	 0.7990	 0.4320
F	 0.8380	 0.3750
H	 0.6350	 0.3430
I	 0.7740	 0.4220
J	 0.7400	 0.3940
K	 0.6770	 0.3280
L	 0.7240	 0.3620

