



wwPDB EM Validation Summary Report ⓘ

Apr 6, 2026 – 03:59 AM UTC

PDB ID : 9UOF / pdb_00009uof
EMDB ID : EMD-64378
Title : PSI-6 FCPI supercomplex from haptophyte *Chrysotila roscoffensis*
Authors : La Rocca, R.; Tsai, P.-C.; Kato, K.; Nakajima, Y.; Akita, F.; Shen, J.-R.
Deposited on : 2025-04-25
Resolution : 1.89 Å(reported)

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

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

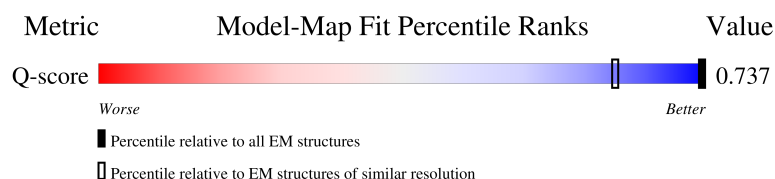
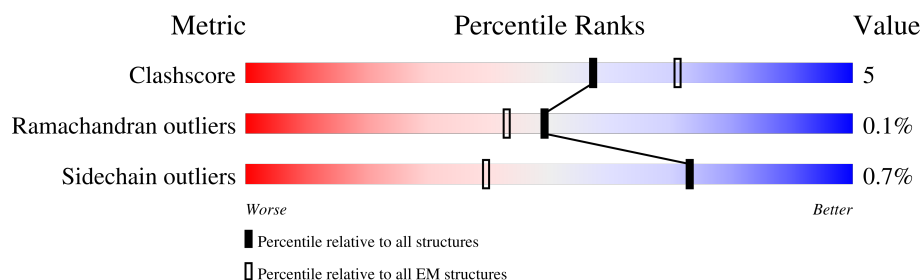
EMDB validation analysis : 0.0.1.dev132
Mogul : 2022.3.0, CSD as543be (2022)
MolProbity : 4-5-2 with Phenix2.0
Buster-report : wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics : 202505.v01 (Using data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

1 Overall quality at a glance

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

The reported resolution of this entry is 1.89 Å.

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



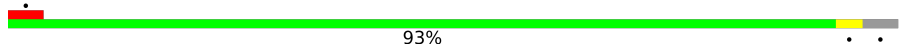

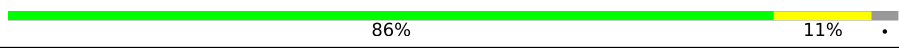
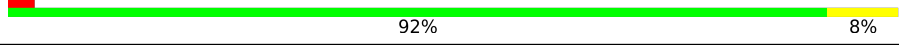

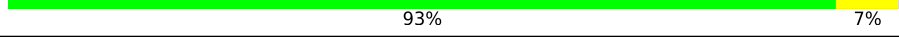

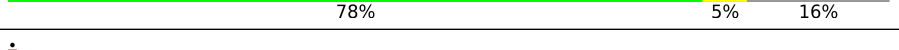
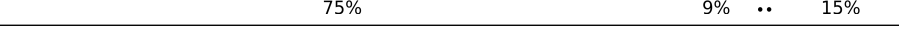
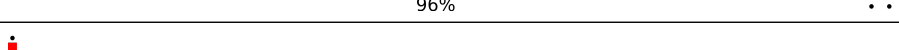


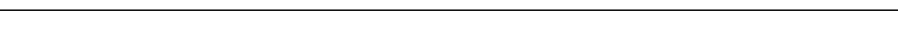
Metric	Whole archive (#Entries)	EM structures (#Entries)	Similar EM resolution (#Entries, resolution range(Å))
Clashscore	229148	23984	-
Ramachandran outliers	224038	23583	-
Sidechain outliers	223484	23102	-
Q-score	-	25397	1004 (1.39 - 2.38)

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	752	92% 6% .
2	B	734	94% 6%
3	C	81	96% ..
4	D	142	88% 9% .

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Mol	Chain	Length	Quality of chain
5	E	67	
6	F	184	
7	I	35	
8	J	39	
9	L	141	
10	M	29	
11	O	201	
12	P	231	
13	Q	197	
14	R	90	
15	S	215	
16	U	191	
17	T	202	

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
18	CLA	A	802	X	-	-	-
18	CLA	A	803	X	-	-	-
18	CLA	A	804	X	-	-	-
18	CLA	A	806	X	-	-	-
18	CLA	A	808	X	-	-	-
18	CLA	A	810	X	-	-	-
18	CLA	A	811	X	-	-	-
18	CLA	A	812	X	-	-	-
18	CLA	A	814	X	-	-	-
18	CLA	A	815	X	-	-	-
18	CLA	A	816	X	-	-	-
18	CLA	A	817	X	-	-	-
18	CLA	A	818	X	-	-	-
18	CLA	A	820	X	-	-	-
18	CLA	A	821	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
18	CLA	A	822	X	-	-	-
18	CLA	A	823	X	-	-	-
18	CLA	A	824	X	-	-	-
18	CLA	A	826	X	-	-	-
18	CLA	A	828	X	-	-	-
18	CLA	A	829	X	-	-	-
18	CLA	A	831	X	-	-	-
18	CLA	A	832	X	-	-	-
18	CLA	A	833	X	-	-	-
18	CLA	A	835	X	-	-	-
18	CLA	A	836	X	-	-	-
18	CLA	A	838	X	-	-	-
18	CLA	A	845	X	-	-	-
18	CLA	A	851	X	-	-	-
18	CLA	A	852	X	-	-	-
18	CLA	A	853	X	-	-	-
18	CLA	B	801	X	-	-	-
18	CLA	B	802	X	-	-	-
18	CLA	B	803	X	-	-	-
18	CLA	B	804	X	-	-	-
18	CLA	B	805	X	-	-	-
18	CLA	B	806	X	-	-	-
18	CLA	B	808	X	-	-	-
18	CLA	B	809	X	-	-	-
18	CLA	B	816	X	-	-	-
18	CLA	B	817	X	-	-	-
18	CLA	B	820	X	-	-	-
18	CLA	B	821	X	-	-	-
18	CLA	B	822	X	-	-	-
18	CLA	B	823	X	-	-	-
18	CLA	B	827	X	-	-	-
18	CLA	B	829	X	-	-	-
18	CLA	B	830	X	-	-	-
18	CLA	B	831	X	-	-	-
18	CLA	B	832	X	-	-	-
18	CLA	B	833	X	-	-	-
18	CLA	B	835	X	-	-	-
18	CLA	B	843	X	-	-	-
18	CLA	B	844	X	-	-	-
18	CLA	B	845	X	-	-	-
18	CLA	B	846	X	-	-	-
18	CLA	B	848	X	-	-	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
18	CLA	F	802	X	-	-	-
18	CLA	F	803	X	-	-	-
18	CLA	J	103	X	-	-	-
18	CLA	L	204	X	-	-	-
18	CLA	O	203	X	-	-	-
18	CLA	O	204	X	-	-	-
18	CLA	O	205	X	-	-	-
18	CLA	O	206	X	-	-	-
18	CLA	O	207	X	-	-	-
18	CLA	O	208	X	-	-	-
18	CLA	P	306	X	-	-	-
18	CLA	P	307	X	-	-	-
18	CLA	P	308	X	-	-	-
18	CLA	P	312	X	-	-	-
18	CLA	P	313	X	-	-	-
18	CLA	P	315	X	-	-	-
18	CLA	Q	204	X	-	-	-
18	CLA	Q	205	X	-	-	-
18	CLA	Q	206	X	-	-	-
18	CLA	Q	207	X	-	-	-
18	CLA	Q	208	X	-	-	-
18	CLA	Q	211	X	-	-	-
18	CLA	Q	213	X	-	-	-
18	CLA	R	202	X	-	-	-
18	CLA	R	205	X	-	-	-
18	CLA	S	204	X	-	-	-
18	CLA	S	205	X	-	-	-
18	CLA	S	206	X	-	-	-
18	CLA	S	214	X	-	-	-
18	CLA	S	215	X	-	-	-
18	CLA	T	202	X	-	-	-
18	CLA	T	203	X	-	-	-
18	CLA	T	204	X	-	-	-
18	CLA	T	205	X	-	-	-
18	CLA	T	206	X	-	-	-
18	CLA	T	207	X	-	-	-
18	CLA	T	212	X	-	-	-
18	CLA	U	204	X	-	-	-
18	CLA	U	206	X	-	-	-
18	CLA	U	208	X	-	-	-
18	CLA	U	211	X	-	-	-

2 Entry composition [i](#)

There are 30 unique types of molecules in this entry. The entry contains 36741 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 (psaA).

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	741	Total	C	N	O	S	0	0
			5813	3807	984	994	28		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
2	B	732	Total	C	N	O	S	0	0
			5805	3823	977	984	21		

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

Mol	Chain	Residues	Atoms					AltConf	Trace
3	C	80	Total	C	N	O	S	0	0
			599	366	106	116	11		

- Molecule 4 is a protein called Photosystem I reaction center subunit II (psaD).

Mol	Chain	Residues	Atoms					AltConf	Trace
4	D	138	Total	C	N	O	S	0	0
			1092	697	188	204	3		

- Molecule 5 is a protein called Photosystem I reaction center subunit IV (psaE).

Mol	Chain	Residues	Atoms					AltConf	Trace
5	E	64	Total	C	N	O	S	0	0
			494	314	86	93	1		

- Molecule 6 is a protein called Photosystem I reaction center subunit III (psaF).

Mol	Chain	Residues	Atoms					AltConf	Trace
6	F	161	Total	C	N	O	S	0	0
			1246	802	209	229	6		

- Molecule 7 is a protein called Photosystem I reaction center subunit VIII (psaI).

Mol	Chain	Residues	Atoms					AltConf	Trace
7	I	34	Total	C	N	O	S	0	0
			266	183	35	46	2		

- Molecule 8 is a protein called Photosystem I reaction center subunit IX (psaJ).

Mol	Chain	Residues	Atoms					AltConf	Trace
8	J	39	Total	C	N	O	S	0	0
			305	204	45	54	2		

- Molecule 9 is a protein called Photosystem I reaction center subunit XI (psaL).

Mol	Chain	Residues	Atoms					AltConf	Trace
9	L	140	Total	C	N	O	S	0	0
			1056	693	168	194	1		

- Molecule 10 is a protein called Photosystem I reaction center subunit XII (psaM).

Mol	Chain	Residues	Atoms					AltConf	Trace
10	M	29	Total	C	N	O	S	0	0
			216	144	34	37	1		

- Molecule 11 is a protein called Fucoxanthin chlorophyll a/c binding protein III (FCPI-3).

Mol	Chain	Residues	Atoms					AltConf	Trace
11	O	176	Total	C	N	O	S	0	0
			1341	872	217	244	8		

- Molecule 12 is a protein called Fucoxanthin chlorophyll a/c binding protein VI (FCPI-6), GLY-CYS-PRO.

Mol	Chain	Residues	Atoms					AltConf	Trace
12	P	193	Total	C	N	O	S	0	0
			1441	927	239	264	11		

- Molecule 13 is a protein called Fucoxanthin chlorophyll a/c binding protein IV (FCPI-4).

Mol	Chain	Residues	Atoms					AltConf	Trace
13	Q	167	Total	C	N	O	S	0	0
			1257	809	202	234	12		

- Molecule 14 is a protein called Photosystem I reaction center subunit psaR.

Mol	Chain	Residues	Atoms					AltConf	Trace
14	R	88	Total	C	N	O	S	0	0
			664	434	106	123	1		

- Molecule 15 is a protein called Fucoxanthin chlorophyll a/c binding protein II (FCPI-2).

Mol	Chain	Residues	Atoms					AltConf	Trace
15	S	165	Total	C	N	O	S	0	0
			1238	802	204	226	6		

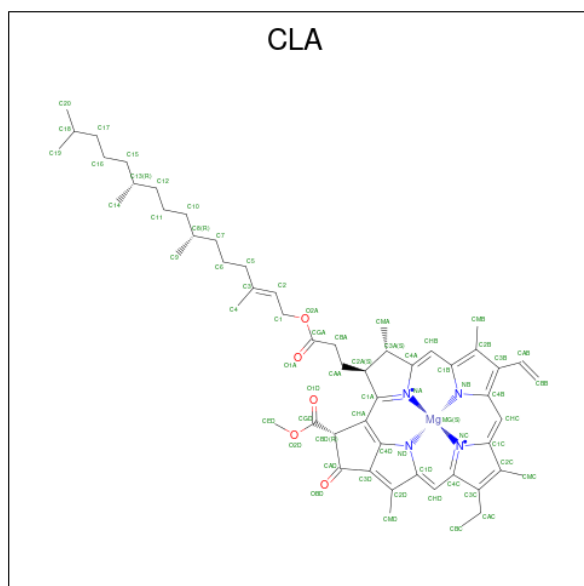
- Molecule 16 is a protein called Fucoxanthin chlorophyll a/c binding protein I (FCPI-1).

Mol	Chain	Residues	Atoms					AltConf	Trace
16	U	141	Total	C	N	O	S	0	0
			1082	692	183	198	9		

- Molecule 17 is a protein called Fucoxanthin chlorophyll a/c binding protein V (FCPI-5).

Mol	Chain	Residues	Atoms					AltConf	Trace
17	T	99	Total	C	N	O	S	0	0
			731	471	122	130	8		

- Molecule 18 is CHLOROPHYLL A (CCD ID: CLA) (formula: $C_{55}H_{72}MgN_4O_5$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms					AltConf
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			55	45	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			49	39	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			62	52	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			54	44	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			50	40	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			43	35	1	4	3	
18	A	1	Total	C	Mg	N	O	0
			51	41	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	A	1	Total	C	Mg	N	O	0
			65	55	1	4	5	

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Mol	Chain	Residues	Atoms					AltConf
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 62	C 52	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 50	C 40	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 50	C 40	Mg 1	N 4	O 5	0
18	A	1	Total 45	C 35	Mg 1	N 4	O 5	0
18	A	1	Total 51	C 41	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 52	C 42	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 60	C 50	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0

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Mol	Chain	Residues	Atoms					AltConf
18	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 45	C 35	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 54	C 44	Mg 1	N 4	O 5	0
18	B	1	Total 55	C 45	Mg 1	N 4	O 5	0
18	B	1	Total 54	C 44	Mg 1	N 4	O 5	0
18	B	1	Total 49	C 39	Mg 1	N 4	O 5	0
18	B	1	Total 55	C 45	Mg 1	N 4	O 5	0
18	B	1	Total 59	C 49	Mg 1	N 4	O 5	0
18	B	1	Total 60	C 50	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 46	C 36	Mg 1	N 4	O 5	0
18	B	1	Total 53	C 43	Mg 1	N 4	O 5	0
18	B	1	Total 63	C 53	Mg 1	N 4	O 5	0

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Mol	Chain	Residues	Atoms					AltConf
18	B	1	Total 64	C 54	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 50	C 40	Mg 1	N 4	O 5	0
18	B	1	Total 49	C 39	Mg 1	N 4	O 5	0
18	B	1	Total 58	C 48	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 58	C 48	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 47	C 37	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	B	1	Total 50	C 40	Mg 1	N 4	O 5	0
18	F	1	Total 48	C 38	Mg 1	N 4	O 5	0

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Mol	Chain	Residues	Atoms					AltConf
18	F	1	Total 46	C 36	Mg 1	N 4	O 5	0
18	J	1	Total 42	C 34	Mg 1	N 4	O 3	0
18	L	1	Total 49	C 39	Mg 1	N 4	O 5	0
18	L	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	L	1	Total 50	C 40	Mg 1	N 4	O 5	0
18	O	1	Total 43	C 35	Mg 1	N 4	O 3	0
18	O	1	Total 45	C 35	Mg 1	N 4	O 5	0
18	O	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	O	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	O	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	O	1	Total 46	C 36	Mg 1	N 4	O 5	0
18	O	1	Total 60	C 50	Mg 1	N 4	O 5	0
18	O	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	O	1	Total 41	C 33	Mg 1	N 4	O 3	0
18	P	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	P	1	Total 56	C 46	Mg 1	N 4	O 5	0
18	P	1	Total 52	C 42	Mg 1	N 4	O 5	0
18	P	1	Total 47	C 37	Mg 1	N 4	O 5	0
18	P	1	Total 50	C 40	Mg 1	N 4	O 5	0
18	P	1	Total 41	C 33	Mg 1	N 4	O 3	0
18	P	1	Total 45	C 35	Mg 1	N 4	O 5	0

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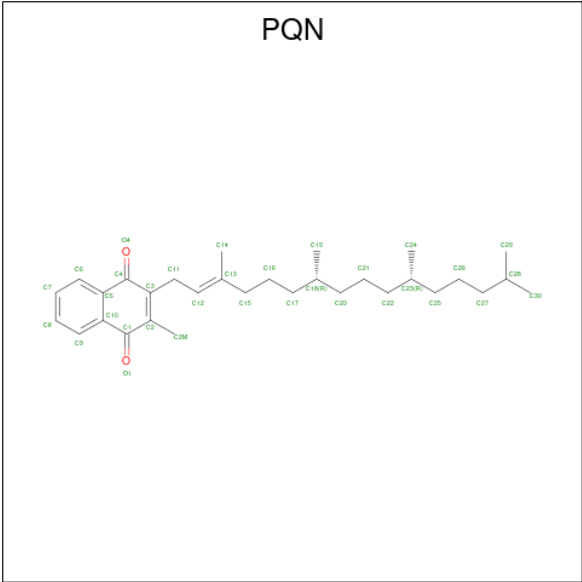
Mol	Chain	Residues	Atoms					AltConf
18	P	1	Total 47	C 37	Mg 1	N 4	O 5	0
18	Q	1	Total 48	C 38	Mg 1	N 4	O 5	0
18	Q	1	Total 61	C 51	Mg 1	N 4	O 5	0
18	Q	1	Total 60	C 50	Mg 1	N 4	O 5	0
18	Q	1	Total 51	C 41	Mg 1	N 4	O 5	0
18	Q	1	Total 46	C 36	Mg 1	N 4	O 5	0
18	Q	1	Total 50	C 40	Mg 1	N 4	O 5	0
18	Q	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	Q	1	Total 41	C 33	Mg 1	N 4	O 3	0
18	Q	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	Q	1	Total 57	C 47	Mg 1	N 4	O 5	0
18	Q	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	R	1	Total 45	C 35	Mg 1	N 4	O 5	0
18	R	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	S	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	S	1	Total 46	C 36	Mg 1	N 4	O 5	0
18	S	1	Total 46	C 36	Mg 1	N 4	O 5	0
18	S	1	Total 45	C 35	Mg 1	N 4	O 5	0
18	S	1	Total 52	C 42	Mg 1	N 4	O 5	0
18	S	1	Total 65	C 55	Mg 1	N 4	O 5	0
18	S	1	Total 65	C 55	Mg 1	N 4	O 5	0

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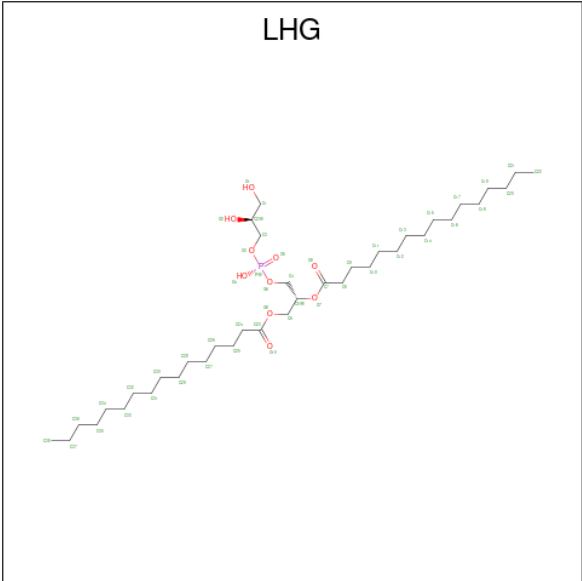
Mol	Chain	Residues	Atoms					AltConf
18	U	1	Total	C	Mg	N	O	0
			61	51	1	4	5	
18	U	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	U	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
18	U	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	U	1	Total	C	Mg	N	O	0
			46	36	1	4	5	
18	U	1	Total	C	Mg	N	O	0
			42	34	1	4	3	
18	U	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	U	1	Total	C	Mg	N	O	0
			52	42	1	4	5	
18	T	1	Total	C	Mg	N	O	0
			42	34	1	4	3	
18	T	1	Total	C	Mg	N	O	0
			41	33	1	4	3	
18	T	1	Total	C	Mg	N	O	0
			46	36	1	4	5	
18	T	1	Total	C	Mg	N	O	0
			57	47	1	4	5	
18	T	1	Total	C	Mg	N	O	0
			46	36	1	4	5	
18	T	1	Total	C	Mg	N	O	0
			42	34	1	4	3	
18	T	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
18	T	1	Total	C	Mg	N	O	0
			41	33	1	4	3	
18	T	1	Total	C	Mg	N	O	0
			46	36	1	4	5	
18	T	1	Total	C	Mg	N	O	0
			47	37	1	4	5	

- Molecule 19 is PHYLLOQUINONE (CCD ID: PQN) (formula: C₃₁H₄₆O₂) (labeled as "Ligand of Interest" by depositor).



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

- Molecule 20 is 1,2-DIPALMITOYL-PHOSPHATIDYL-GLYCEROLE (CCD ID: LHG) (formula: C₃₈H₇₅O₁₀P) (labeled as "Ligand of Interest" by depositor).



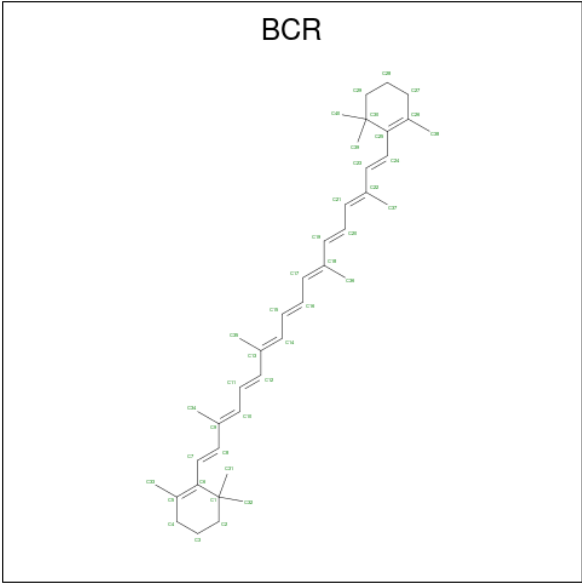
Mol	Chain	Residues	Atoms				AltConf
20	A	1	Total	C	O	P	0
			48	37	10	1	

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Mol	Chain	Residues	Atoms				AltConf
20	A	1	Total	C	O	P	0
			27	16	10	1	
20	R	1	Total	C	O	P	0
			49	38	10	1	

- Molecule 21 is BETA-CAROTENE (CCD ID: BCR) (formula: C₄₀H₅₆) (labeled as "Ligand of Interest" by depositor).



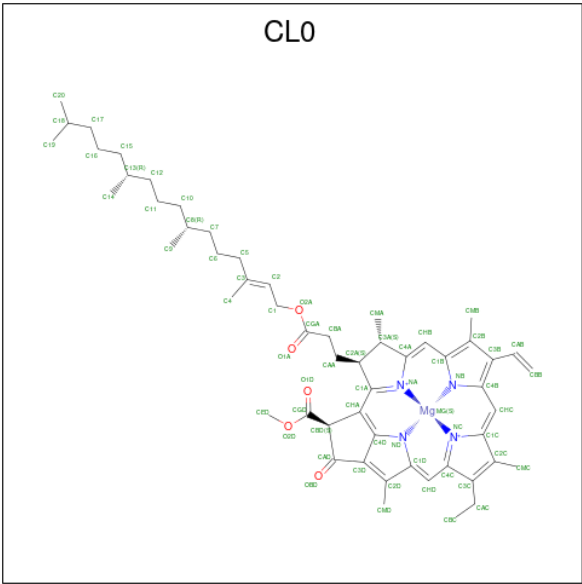
Mol	Chain	Residues	Atoms		AltConf
21	A	1	Total	C	0
			40	40	
21	A	1	Total	C	0
			40	40	
21	A	1	Total	C	0
			40	40	
21	A	1	Total	C	0
			40	40	
21	B	1	Total	C	0
			40	40	
21	B	1	Total	C	0
			40	40	
21	B	1	Total	C	0
			40	40	
21	B	1	Total	C	0
			40	40	

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Mol	Chain	Residues	Atoms		AltConf
21	F	1	Total	C	0
			40	40	
21	F	1	Total	C	0
			40	40	
21	I	1	Total	C	0
			40	40	
21	I	1	Total	C	0
			40	40	
21	J	1	Total	C	0
			40	40	
21	L	1	Total	C	0
			40	40	
21	L	1	Total	C	0
			40	40	
21	M	1	Total	C	0
			40	40	
21	R	1	Total	C	0
			39	39	

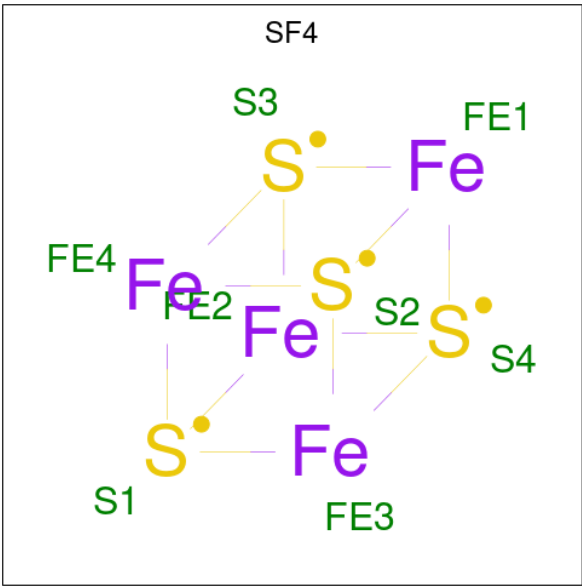
- Molecule 22 is CHLOROPHYLL A ISOMER (CCD ID: CL0) (formula: C₅₅H₇₂MgN₄O₅) (labeled as "Ligand of Interest" by depositor).



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

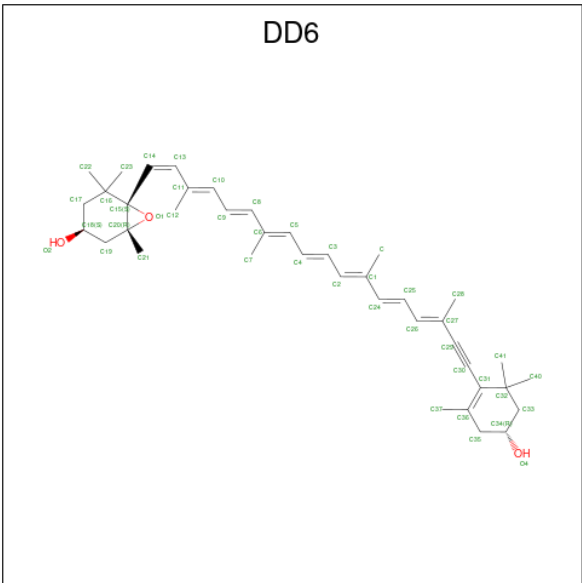
- Molecule 23 is IRON/SULFUR CLUSTER (CCD ID: SF4) (formula: Fe₄S₄) (labeled as

"Ligand of Interest" by depositor).



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

- Molecule 24 is (3S,3'R,5R,6S,7cis)-7',8'-didehydro-5,6-dihydro-5,6-epoxy-beta,beta-carotene-3,3'-diol (CCD ID: DD6) (formula: C₄₀H₅₄O₃) (labeled as "Ligand of Interest" by depositor).



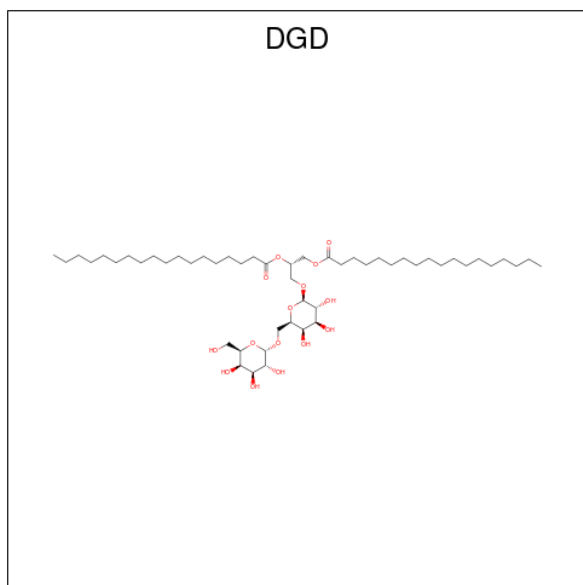
Mol	Chain	Residues	Atoms			AltConf
24	A	1	Total 43	C 40	O 3	0
24	J	1	Total 43	C 40	O 3	0
24	O	1	Total 43	C 40	O 3	0
24	O	1	Total 43	C 40	O 3	0
24	O	1	Total 43	C 40	O 3	0
24	O	1	Total 43	C 40	O 3	0
24	O	1	Total 43	C 40	O 3	0
24	P	1	Total 43	C 40	O 3	0
24	P	1	Total 43	C 40	O 3	0
24	P	1	Total 43	C 40	O 3	0
24	P	1	Total 23	C 21	O 2	0
24	Q	1	Total 43	C 40	O 3	0
24	Q	1	Total 43	C 40	O 3	0
24	S	1	Total 43	C 40	O 3	0
24	S	1	Total 43	C 40	O 3	0
24	S	1	Total 43	C 40	O 3	0
24	S	1	Total 43	C 40	O 3	0
24	S	1	Total 43	C 40	O 3	0
24	U	1	Total 43	C 40	O 3	0
24	U	1	Total 43	C 40	O 3	0
24	U	1	Total 26	C 25	O 1	0
24	T	1	Total 43	C 40	O 3	0

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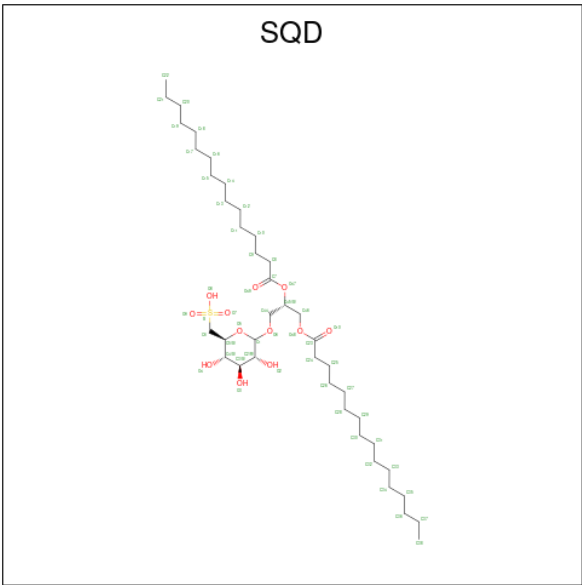
Mol	Chain	Residues	Atoms			AltConf
24	T	1	Total	C	O	0
			43	40	3	

- Molecule 25 is DIGALACTOSYL DIACYL GLYCEROL (DGDG) (CCD ID: DGD) (formula: $C_{51}H_{96}O_{15}$) (labeled as "Ligand of Interest" by depositor).



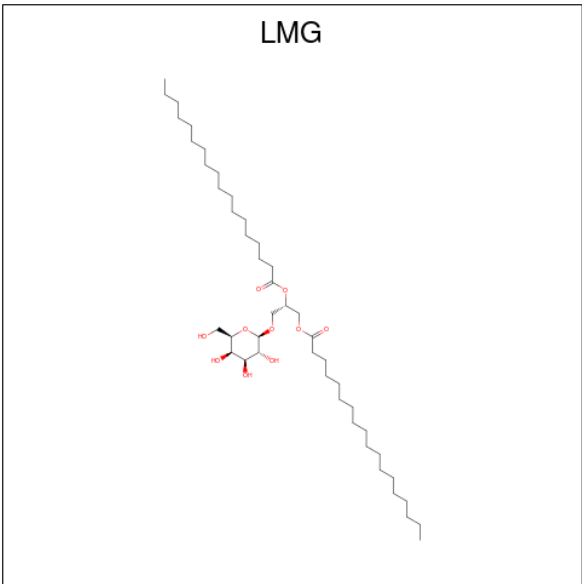
Mol	Chain	Residues	Atoms			AltConf
25	B	1	Total	C	O	0
			60	45	15	

- Molecule 26 is 1,2-DI-O-ACYL-3-O-[6-DEOXY-6-SULFO-ALPHA-D-GLUCOPYRANOSYL]-SN-GLYCEROL (CCD ID: SQD) (formula: $C_{41}H_{78}O_{12}S$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				AltConf
26	B	1	Total	C	O	S	0
			50	37	12	1	
26	M	1	Total	C	O	S	0
			46	33	12	1	

- Molecule 27 is 1,2-DISTEAROYL-MONOGALACTOSYL-DIGLYCERIDE (CCD ID: LMG) (formula: C₄₅H₈₆O₁₀) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			AltConf
27	B	1	Total	C	O	0
			25	15	10	

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Mol	Chain	Residues	Atoms			AltConf
27	J	1	Total 39	C 29	O 10	0
27	P	1	Total 34	C 24	O 10	0
27	Q	1	Total 55	C 45	O 10	0
27	S	1	Total 49	C 39	O 10	0
27	U	1	Total 32	C 22	O 10	0

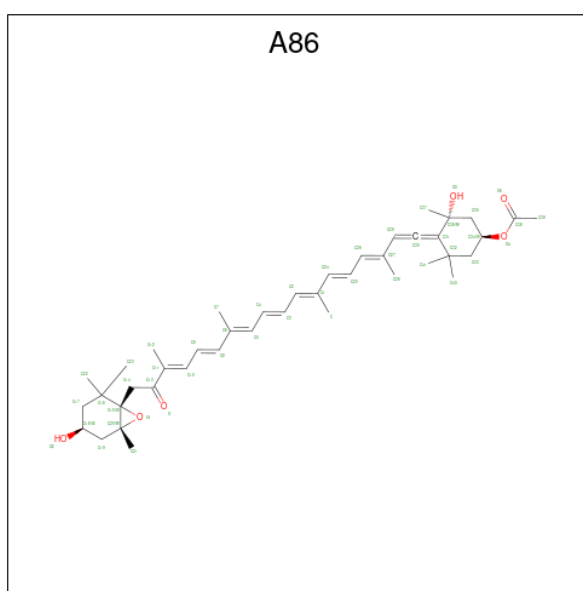
- # KC1

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Mol	Chain	Residues	Atoms					AltConf
28	S	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
28	U	1	Total	C	Mg	N	O	0
			45	35	1	4	5	
28	T	1	Total	C	Mg	N	O	0
			45	35	1	4	5	

- Molecule 29 is (3S,3'S,5R,5'R,6S,6'R,8'R)-3,5'-dihydroxy-8-oxo-6',7'-didehydro-5,5',6,6',7,8-hexahydro-5,6-epoxy-beta,beta-caroten-3'-yl acetate (CCD ID: A86) (formula: $C_{42}H_{58}O_6$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms			AltConf
29	P	1	Total	C	O	0
			48	42	6	
29	Q	1	Total	C	O	0
			48	42	6	
29	Q	1	Total	C	O	0
			48	42	6	
29	R	1	Total	C	O	0
			44	40	4	
29	R	1	Total	C	O	0
			48	42	6	
29	U	1	Total	C	O	0
			48	42	6	
29	T	1	Total	C	O	0
			48	42	6	

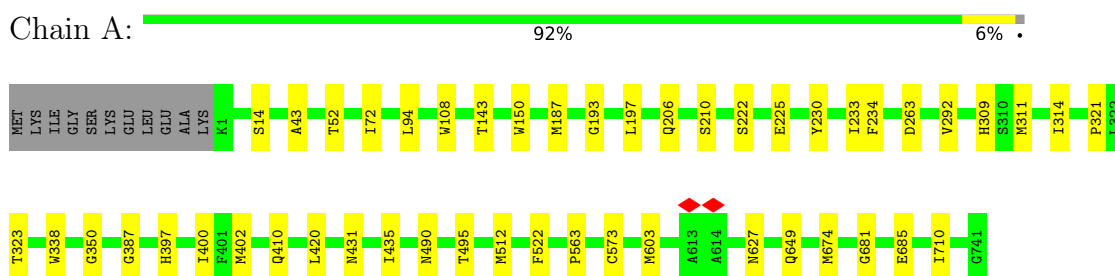
- Molecule 30 is water.

Mol	Chain	Residues	Atoms		AltConf
30	A	189	Total 189	O 189	0
30	B	286	Total 286	O 286	0
30	C	59	Total 59	O 59	0
30	D	39	Total 39	O 39	0
30	E	13	Total 13	O 13	0
30	F	33	Total 33	O 33	0
30	I	2	Total 2	O 2	0
30	J	2	Total 2	O 2	0
30	L	23	Total 23	O 23	0
30	M	5	Total 5	O 5	0
30	O	22	Total 22	O 22	0
30	P	30	Total 30	O 30	0
30	Q	19	Total 19	O 19	0
30	R	27	Total 27	O 27	0
30	S	25	Total 25	O 25	0
30	U	9	Total 9	O 9	0
30	T	7	Total 7	O 7	0

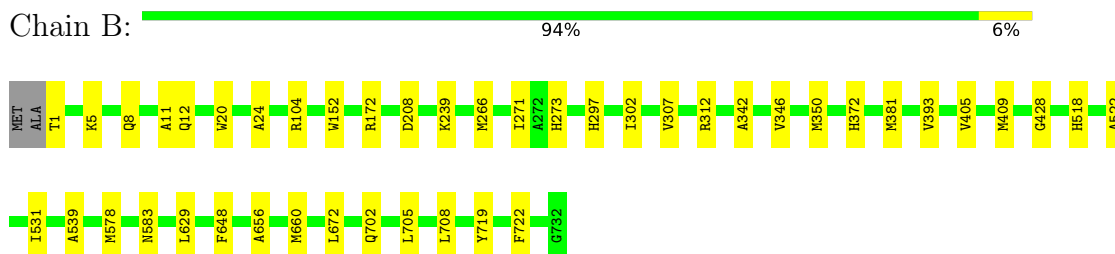
3 Residue-property plots

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.

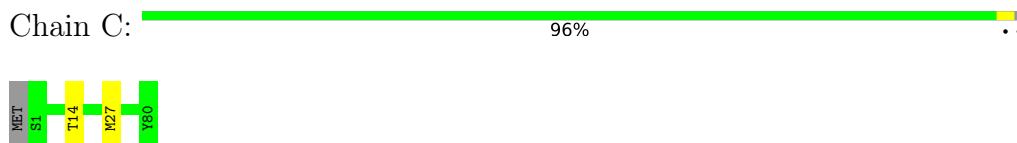
- Molecule 1: Photosystem I P700 chlorophyll a apoprotein A1 (psaA)



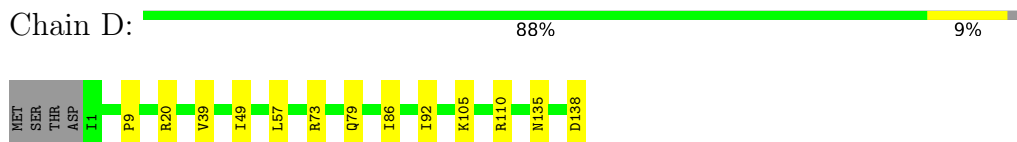
- Molecule 2: Photosystem I P700 chlorophyll a apoprotein A2 (psaB)



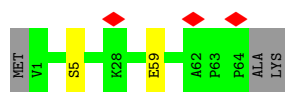
- Molecule 3: Photosystem I iron-sulfur center (psaC)



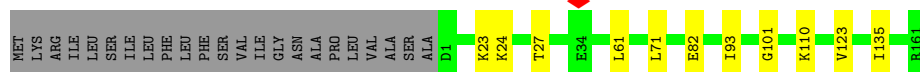
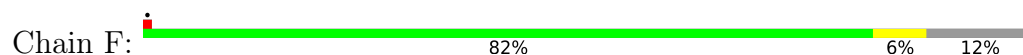
- Molecule 4: Photosystem I reaction center subunit II (psaD)



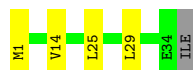
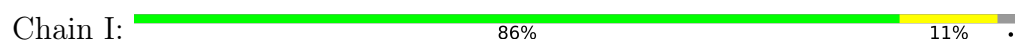
- Molecule 5: Photosystem I reaction center subunit IV (psaE)



- Molecule 6: Photosystem I reaction center subunit III (psaF)



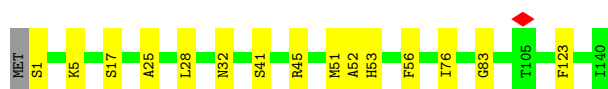
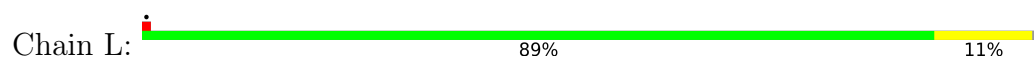
- Molecule 7: Photosystem I reaction center subunit VIII (psaI)



- Molecule 8: Photosystem I reaction center subunit IX (psaJ)



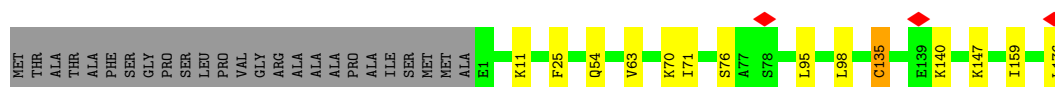
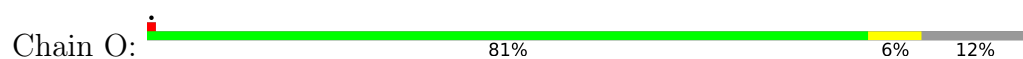
- Molecule 9: Photosystem I reaction center subunit XI (psaL)




- Molecule 10: Photosystem I reaction center subunit XII (psaM)

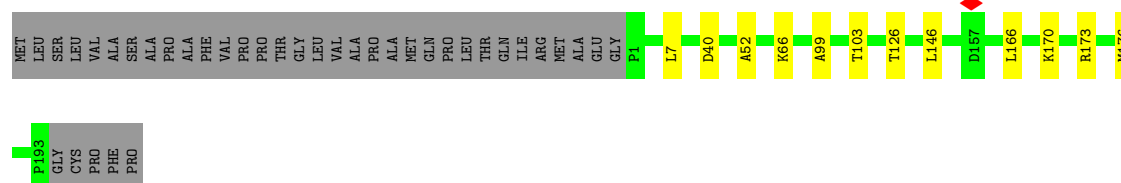


- Molecule 11: Fucoxanthin chlorophyll a/c binding protein III (FCPI-3)



- Molecule 12: Fucoxanthin chlorophyll a/c binding protein VI (FCPI-6), GLY-CYS-PRO

Chain P:  78% 5% 16%



- Molecule 13: Fucoxanthin chlorophyll a/c binding protein IV (FCPI-4)

Chain Q:  75% 9% 15%



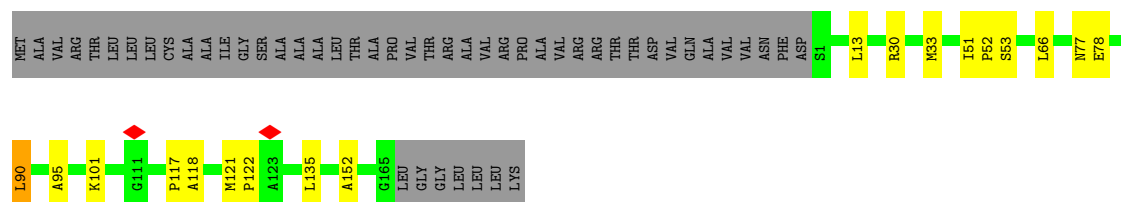
- Molecule 14: Photosystem I reaction center subunit psaR

Chain R:  96% 4% 0%



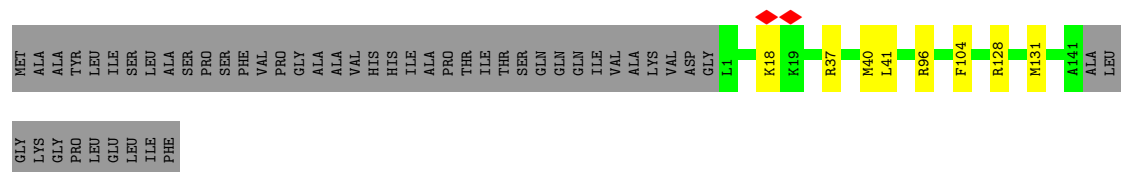
- Molecule 15: Fucoxanthin chlorophyll a/c binding protein II (FCPI-2)

Chain S:  68% 8% 23%



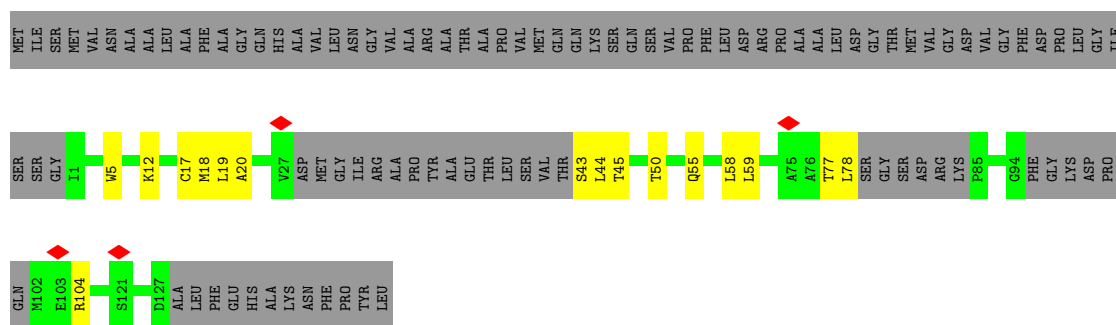
- Molecule 16: Fucoxanthin chlorophyll a/c binding protein I (FCPI-1)

Chain U:  70% 1% 26%



- Molecule 17: Fucoxanthin chlorophyll a/c binding protein V (FCPI-5)

Frequency	Percentage
Daily	41%
Weekly	8%
Monthly	51%



4 Experimental information

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C1	Depositor
Number of particles used	114380	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS KRIOS	Depositor
Voltage (kV)	300	Depositor
Electron dose ($e^-/\text{\AA}^2$)	50	Depositor
Minimum defocus (nm)	800	Depositor
Maximum defocus (nm)	1800	Depositor
Magnification	165000	Depositor
Image detector	FEI FALCON IV (4k x 4k)	Depositor
Maximum map value	1.114	Depositor
Minimum map value	-0.354	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.017	Depositor
Recommended contour level	0.08	Depositor
Map size (Å)	436.2, 436.2, 436.2	wwPDB
Map dimensions	600, 600, 600	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	0.727, 0.727, 0.727	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: DD6, DGD, A86, KC1, CLA, SQD, BCR, CL0, LMG, LHG, SF4, PQN

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	0.26	0/6007	0.48	0/8185
2	B	0.25	0/6015	0.47	0/8205
3	C	0.21	0/609	0.45	0/826
4	D	0.21	0/1116	0.50	0/1503
5	E	0.22	0/505	0.41	0/689
6	F	0.24	0/1275	0.47	0/1728
7	I	0.26	0/273	0.53	0/373
8	J	0.26	0/313	0.55	0/427
9	L	0.24	0/1081	0.52	0/1470
10	M	0.21	0/218	0.34	0/295
11	O	0.24	0/1376	0.51	0/1865
12	P	0.22	0/1480	0.49	0/2010
13	Q	0.30	0/1285	0.58	2/1736 (0.1%)
14	R	0.26	0/681	0.47	0/930
15	S	0.20	0/1272	0.41	0/1732
16	U	0.24	0/1109	0.49	0/1499
17	T	0.28	0/741	0.60	1/999 (0.1%)
All	All	0.25	0/25356	0.49	3/34472 (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
13	Q	0	1

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
17	T	19	LEU	N-CA-C	-5.92	106.70	114.04
13	Q	91	MET	CB-CG-SD	5.04	127.83	112.70
13	Q	165	PRO	CA-N-CD	-5.00	105.00	112.00

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
13	Q	164	PHE	Peptide

5.2 Too-close contacts ⓘ

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	5813	0	5698	32	0
2	B	5805	0	5634	36	0
3	C	599	0	577	1	0
4	D	1092	0	1096	7	0
5	E	494	0	488	1	0
6	F	1246	0	1256	8	0
7	I	266	0	278	4	0
8	J	305	0	310	3	0
9	L	1056	0	1068	14	0
10	M	216	0	234	3	0
11	O	1341	0	1347	10	0
12	P	1441	0	1421	8	0
13	Q	1257	0	1260	11	0
14	R	664	0	668	2	0
15	S	1238	0	1217	12	0
16	U	1082	0	1058	6	0
17	T	731	0	749	11	0
18	A	2615	0	2651	59	0
18	B	2399	0	2438	70	0
18	F	94	0	69	1	0
18	J	42	0	31	2	0
18	L	164	0	150	4	0
18	O	495	0	475	13	0
18	P	403	0	337	4	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
18	Q	609	0	566	13	0
18	R	110	0	105	4	0
18	S	384	0	358	13	0
18	T	473	0	379	13	0
18	U	441	0	417	13	0
19	A	33	0	46	2	0
19	B	33	0	46	1	0
20	A	75	0	93	2	0
20	R	49	0	74	2	0
21	A	160	0	224	5	0
21	B	200	0	280	7	0
21	F	80	0	112	4	0
21	I	80	0	112	4	0
21	J	40	0	56	1	0
21	L	80	0	112	6	0
21	M	40	0	56	4	0
21	R	39	0	53	3	0
22	A	65	0	72	3	0
23	A	8	0	0	0	0
23	C	16	0	0	0	0
24	A	43	0	0	0	0
24	J	43	0	0	0	0
24	O	215	0	0	0	0
24	P	152	0	0	1	0
24	Q	86	0	0	0	0
24	S	215	0	0	0	0
24	T	86	0	0	0	0
24	U	112	0	0	0	0
25	B	60	0	81	5	0
26	B	50	0	67	3	0
26	M	46	0	56	3	0
27	B	25	0	20	2	0
27	J	39	0	48	2	0
27	P	34	0	38	1	0
27	Q	55	0	86	6	0
27	S	49	0	68	6	0
27	U	32	0	34	2	0
28	O	45	0	0	1	0
28	P	134	0	0	0	0
28	Q	45	0	0	0	0
28	S	90	0	0	0	0
28	T	45	0	0	2	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
28	U	45	0	0	0	0
29	P	48	0	0	0	0
29	Q	96	0	0	0	0
29	R	92	0	0	2	0
29	T	48	0	0	0	0
29	U	48	0	0	1	0
30	A	189	0	0	1	0
30	B	286	0	0	1	0
30	C	59	0	0	0	0
30	D	39	0	0	1	0
30	E	13	0	0	0	0
30	F	33	0	0	1	0
30	I	2	0	0	0	0
30	J	2	0	0	0	0
30	L	23	0	0	0	0
30	M	5	0	0	0	0
30	O	22	0	0	0	0
30	P	30	0	0	0	0
30	Q	19	0	0	0	0
30	R	27	0	0	0	0
30	S	25	0	0	0	0
30	T	7	0	0	0	0
30	U	9	0	0	0	0
All	All	36741	0	34169	336	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 5.

The worst 5 of 336 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:193:GLY:O	1:A:197:LEU:HB2	1.87	0.75
27:U:201:LMG:H321	18:U:211:CLA:H42	1.74	0.69
1:A:522:PHE:HA	18:A:833:CLA:HED1	1.75	0.69
18:A:838:CLA:HED1	18:U:210:CLA:HAA2	1.76	0.68
18:Q:206:CLA:H42	27:Q:217:LMG:H182	1.76	0.67

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	739/752 (98%)	722 (98%)	17 (2%)	0	100	100
2	B	730/734 (100%)	714 (98%)	16 (2%)	0	100	100
3	C	78/81 (96%)	78 (100%)	0	0	100	100
4	D	136/142 (96%)	132 (97%)	4 (3%)	0	100	100
5	E	62/67 (92%)	61 (98%)	1 (2%)	0	100	100
6	F	159/184 (86%)	157 (99%)	2 (1%)	0	100	100
7	I	32/35 (91%)	32 (100%)	0	0	100	100
8	J	37/39 (95%)	36 (97%)	1 (3%)	0	100	100
9	L	138/141 (98%)	137 (99%)	1 (1%)	0	100	100
10	M	27/29 (93%)	27 (100%)	0	0	100	100
11	O	174/201 (87%)	171 (98%)	3 (2%)	0	100	100
12	P	191/231 (83%)	190 (100%)	1 (0%)	0	100	100
13	Q	165/197 (84%)	157 (95%)	6 (4%)	2 (1%)	10	4
14	R	86/90 (96%)	86 (100%)	0	0	100	100
15	S	163/215 (76%)	161 (99%)	2 (1%)	0	100	100
16	U	139/191 (73%)	138 (99%)	1 (1%)	0	100	100
17	T	91/202 (45%)	90 (99%)	1 (1%)	0	100	100
All	All	3147/3531 (89%)	3089 (98%)	56 (2%)	2 (0%)	49	40

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
13	Q	165	PRO
13	Q	33	LEU

5.3.2 Protein sidechains ⓘ

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	603/612 (98%)	598 (99%)	5 (1%)	73	75
2	B	590/591 (100%)	587 (100%)	3 (0%)	81	84
3	C	68/69 (99%)	68 (100%)	0	100	100
4	D	118/122 (97%)	118 (100%)	0	100	100
5	E	53/55 (96%)	53 (100%)	0	100	100
6	F	133/152 (88%)	132 (99%)	1 (1%)	73	75
7	I	31/32 (97%)	31 (100%)	0	100	100
8	J	32/32 (100%)	32 (100%)	0	100	100
9	L	111/112 (99%)	111 (100%)	0	100	100
10	M	21/21 (100%)	21 (100%)	0	100	100
11	O	145/161 (90%)	143 (99%)	2 (1%)	59	59
12	P	144/173 (83%)	144 (100%)	0	100	100
13	Q	133/157 (85%)	131 (98%)	2 (2%)	57	56
14	R	71/73 (97%)	71 (100%)	0	100	100
15	S	125/162 (77%)	122 (98%)	3 (2%)	43	38
16	U	110/148 (74%)	110 (100%)	0	100	100
17	T	73/153 (48%)	71 (97%)	2 (3%)	39	34
All	All	2561/2825 (91%)	2543 (99%)	18 (1%)	73	78

5 of 18 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
15	S	90	LEU
17	T	77	THR
17	T	43	SER
6	F	135	ILE
15	S	53	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 19

such sidechains are listed below:

Mol	Chain	Res	Type
14	R	49	ASN
15	S	72	ASN
16	U	120	GLN
15	S	46	GLN
6	F	15	ASN

5.3.3 RNA [i](#)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates [i](#)

There are no oligosaccharides in this entry.

5.6 Ligand geometry [i](#)

220 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	$\# Z > 2$
18	CLA	A	846	1	64,68,73	1.38	7 (10%)	76,107,113	0.88	4 (5%)
21	BCR	F	804	-	41,41,41	1.03	2 (4%)	56,56,56	1.27	4 (7%)
18	CLA	A	826	1	69,73,73	1.30	8 (11%)	82,113,113	0.86	3 (3%)
18	CLA	A	845	30	69,73,73	1.17	8 (11%)	82,113,113	0.92	3 (3%)
18	CLA	O	209	11	69,73,73	1.18	7 (10%)	82,113,113	1.19	8 (9%)
18	CLA	A	853	30	69,73,73	1.19	7 (10%)	82,113,113	1.02	6 (7%)
24	DD6	Q	215	-	40,45,45	1.38	8 (20%)	51,67,67	1.53	10 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
18	CLA	T	210	17	45,49,73	1.41	5 (11%)	54,84,113	1.25	6 (11%)
23	SF4	C	102	3	0,12,12	-	-	-		
18	CLA	B	807	2	69,73,73	1.19	8 (11%)	82,113,113	0.93	4 (4%)
18	CLA	B	805	2	69,73,73	1.20	9 (13%)	82,113,113	0.93	5 (6%)
18	CLA	O	205	11	69,73,73	1.19	7 (10%)	82,113,113	0.88	4 (4%)
18	CLA	U	206	16	49,53,73	1.40	6 (12%)	58,89,113	1.05	4 (6%)
18	CLA	U	207	-	69,73,73	1.14	6 (8%)	82,113,113	1.09	8 (9%)
18	CLA	A	838	20	56,60,73	1.27	8 (14%)	65,97,113	1.03	4 (6%)
18	CLA	P	312	12	45,49,73	1.43	6 (13%)	54,84,113	1.21	6 (11%)
18	CLA	A	851	1	69,73,73	1.20	8 (11%)	82,113,113	0.92	3 (3%)
18	CLA	U	204	-	65,69,73	1.21	7 (10%)	77,108,113	0.92	5 (6%)
18	CLA	S	215	15	69,73,73	1.16	7 (10%)	82,113,113	0.91	3 (3%)
18	CLA	O	203	-	49,53,73	1.42	7 (14%)	58,89,113	1.08	4 (6%)
18	CLA	U	209	16	46,50,73	1.46	7 (15%)	53,85,113	1.12	4 (7%)
18	CLA	B	828	2	62,66,73	1.30	8 (12%)	73,104,113	1.00	5 (6%)
18	CLA	P	315	12	51,55,73	1.39	8 (15%)	60,91,113	1.15	5 (8%)
18	CLA	U	210	16	69,73,73	1.15	8 (11%)	82,113,113	0.94	5 (6%)
18	CLA	B	833	-	69,73,73	1.20	8 (11%)	82,113,113	1.00	6 (7%)
24	DD6	P	317	-	21,24,45	1.41	6 (28%)	27,37,67	1.85	6 (22%)
27	LMG	S	211	-	49,49,55	0.74	1 (2%)	57,57,63	1.25	4 (7%)
27	LMG	B	849	-	25,25,55	1.38	5 (20%)	33,33,63	1.25	4 (12%)
24	DD6	O	215	-	40,45,45	1.27	7 (17%)	51,67,67	1.61	11 (21%)
18	CLA	B	811	2	59,63,73	1.26	7 (11%)	70,101,113	0.95	3 (4%)
18	CLA	B	803	-	69,73,73	1.15	7 (10%)	82,113,113	0.95	5 (6%)
26	SQD	M	102	-	44,46,54	0.99	3 (6%)	54,57,65	1.56	9 (16%)
18	CLA	O	206	-	69,73,73	1.15	8 (11%)	82,113,113	0.88	4 (4%)
18	CLA	A	803	1	69,73,73	1.19	8 (11%)	82,113,113	0.96	4 (4%)
18	CLA	T	203	17	45,49,73	1.42	7 (15%)	54,84,113	1.14	4 (7%)
18	CLA	O	202	11	47,51,73	1.42	7 (14%)	55,86,113	1.10	4 (7%)
18	CLA	A	828	1	69,73,73	1.23	6 (8%)	82,113,113	0.92	4 (4%)
18	CLA	A	832	1	55,59,73	1.39	8 (14%)	64,96,113	1.07	4 (6%)
18	CLA	S	214	30	69,73,73	1.15	6 (8%)	82,113,113	0.97	5 (6%)
18	CLA	B	848	2	54,58,73	1.35	8 (14%)	64,95,113	1.09	5 (7%)
18	CLA	A	827	1	54,58,73	1.43	7 (12%)	64,95,113	1.02	4 (6%)
18	CLA	A	807	1	69,73,73	1.20	8 (11%)	82,113,113	0.98	5 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
24	DD6	S	212	-	40,45,45	1.30	7 (17%)	51,67,67	1.44	8 (15%)
18	CLA	A	830	1	54,58,73	1.32	8 (14%)	64,95,113	1.01	5 (7%)
24	DD6	O	213	-	40,45,45	1.35	6 (15%)	51,67,67	1.37	7 (13%)
18	CLA	R	205	14	69,73,73	1.26	9 (13%)	82,113,113	0.94	5 (6%)
18	CLA	A	808	1	60,64,73	1.32	8 (13%)	71,102,113	1.18	6 (8%)
18	CLA	Q	211	13	45,49,73	1.47	7 (15%)	54,84,113	1.21	6 (11%)
24	DD6	O	214	-	40,45,45	1.24	7 (17%)	51,67,67	1.48	10 (19%)
18	CLA	A	821	-	69,73,73	1.22	8 (11%)	82,113,113	0.99	6 (7%)
18	CLA	T	207	17	46,50,73	1.42	6 (13%)	53,85,113	1.12	4 (7%)
28	KC1	P	311	12	46,52,53	1.68	8 (17%)	59,87,89	1.32	7 (11%)
20	LHG	R	201	-	48,48,48	0.67	2 (4%)	51,54,54	1.27	5 (9%)
23	SF4	C	101	3	0,12,12	-	-	-	-	-
21	BCR	R	203	-	40,40,41	1.19	4 (10%)	54,54,56	1.46	11 (20%)
24	DD6	U	203	-	40,45,45	1.36	9 (22%)	51,67,67	1.84	11 (21%)
29	A86	R	204	-	44,46,50	1.40	4 (9%)	49,70,76	1.87	15 (30%)
18	CLA	A	829	1	69,73,73	1.21	8 (11%)	82,113,113	0.91	4 (4%)
18	CLA	J	103	8	46,50,73	1.47	7 (15%)	53,85,113	1.08	3 (5%)
18	CLA	U	208	16	50,54,73	1.35	6 (12%)	59,90,113	1.11	5 (8%)
19	PQN	A	837	-	34,34,34	0.41	0	43,45,45	0.61	1 (2%)
24	DD6	P	316	-	40,45,45	1.28	6 (15%)	51,67,67	1.76	12 (23%)
18	CLA	P	310	12	54,58,73	1.31	7 (12%)	64,95,113	1.08	3 (4%)
18	CLA	B	814	2	59,63,73	1.42	8 (13%)	70,101,113	1.00	4 (5%)
18	CLA	B	816	2	64,68,73	1.26	8 (12%)	76,107,113	0.96	4 (5%)
21	BCR	F	801	-	41,41,41	1.05	2 (4%)	56,56,56	1.31	6 (10%)
18	CLA	B	829	30	69,73,73	1.17	7 (10%)	82,113,113	0.99	5 (6%)
18	CLA	P	307	12	60,64,73	1.27	7 (11%)	71,102,113	1.00	6 (8%)
18	CLA	B	815	2	63,67,73	1.23	7 (11%)	74,105,113	1.04	5 (6%)
18	CLA	S	204	15	50,54,73	1.36	7 (14%)	59,90,113	1.06	4 (6%)
18	CLA	B	818	2	50,54,73	1.40	8 (16%)	59,90,113	1.15	5 (8%)
24	DD6	T	213	-	40,45,45	1.33	7 (17%)	51,67,67	1.70	10 (19%)
21	BCR	A	844	-	41,41,41	1.09	2 (4%)	56,56,56	1.24	7 (12%)
18	CLA	B	808	2	69,73,73	1.24	8 (11%)	82,113,113	0.89	2 (2%)
18	CLA	R	202	30	49,53,73	1.41	8 (16%)	58,89,113	1.14	7 (12%)
24	DD6	O	212	-	40,45,45	1.39	8 (20%)	51,67,67	1.68	10 (19%)
28	KC1	U	213	16	49,53,53	1.60	9 (18%)	61,89,89	0.96	3 (4%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
27	LMG	P	301	-	34,34,55	1.13	3 (8%)	42,42,63	1.31	2 (4%)
29	A86	T	201	-	47,50,50	1.40	5 (10%)	51,76,76	1.40	9 (17%)
18	CLA	A	814	30	49,53,73	1.41	8 (16%)	58,89,113	1.10	4 (6%)
18	CLA	A	801	-	69,73,73	1.17	7 (10%)	82,113,113	0.90	4 (4%)
18	CLA	B	822	2	69,73,73	1.25	8 (11%)	82,113,113	0.96	4 (4%)
18	CLA	T	205	-	61,65,73	1.21	6 (9%)	72,103,113	0.94	3 (4%)
29	A86	Q	201	-	47,50,50	1.37	5 (10%)	51,76,76	1.67	11 (21%)
18	CLA	B	817	-	69,73,73	1.23	7 (10%)	82,113,113	0.99	3 (3%)
18	CLA	B	821	-	68,72,73	1.15	8 (11%)	80,111,113	0.92	5 (6%)
21	BCR	B	840	-	41,41,41	1.03	2 (4%)	56,56,56	1.27	6 (10%)
18	CLA	B	845	2	69,73,73	1.27	8 (11%)	82,113,113	0.98	4 (4%)
21	BCR	L	205	-	41,41,41	1.07	2 (4%)	56,56,56	1.25	4 (7%)
28	KC1	S	210	15	49,53,53	1.58	9 (18%)	61,89,89	1.12	5 (8%)
18	CLA	L	204	-	54,58,73	1.34	8 (14%)	64,95,113	1.10	5 (7%)
18	CLA	P	309	30	51,55,73	1.34	7 (13%)	60,91,113	1.00	4 (6%)
18	CLA	S	205	15	50,54,73	1.45	7 (14%)	59,90,113	1.05	4 (6%)
27	LMG	Q	217	-	55,55,55	0.75	0	63,63,63	1.33	8 (12%)
18	CLA	A	811	1	69,73,73	1.22	7 (10%)	82,113,113	0.94	5 (6%)
18	CLA	Q	208	13	54,58,73	1.31	7 (12%)	64,95,113	1.02	5 (7%)
18	CLA	S	206	15	49,53,73	1.33	6 (12%)	58,89,113	1.02	5 (8%)
18	CLA	B	846	2	69,73,73	1.20	8 (11%)	82,113,113	1.19	6 (7%)
18	CLA	B	813	2	53,57,73	1.40	8 (15%)	61,93,113	1.10	4 (6%)
29	A86	Q	214	-	47,50,50	1.43	4 (8%)	51,76,76	1.39	9 (17%)
18	CLA	Q	212	13	69,73,73	1.23	7 (10%)	82,113,113	1.01	7 (8%)
18	CLA	O	211	11	45,49,73	1.47	7 (15%)	54,84,113	1.19	6 (11%)
18	CLA	A	816	1	69,73,73	1.22	8 (11%)	82,113,113	0.99	5 (6%)
18	CLA	Q	207	13	50,54,73	1.39	7 (14%)	59,90,113	1.17	5 (8%)
27	LMG	U	201	-	32,32,55	1.00	0	40,40,63	1.16	4 (10%)
18	CLA	Q	204	-	65,69,73	1.19	5 (7%)	77,108,113	0.97	6 (7%)
18	CLA	B	823	2	69,73,73	1.24	7 (10%)	82,113,113	0.92	2 (2%)
18	CLA	B	801	-	69,73,73	1.19	8 (11%)	82,113,113	0.94	5 (6%)
18	CLA	B	835	30	69,73,73	1.22	7 (10%)	82,113,113	1.03	7 (8%)
18	CLA	A	820	1	55,59,73	1.38	7 (12%)	64,96,113	1.15	5 (7%)
21	BCR	B	841	-	41,41,41	1.08	2 (4%)	56,56,56	1.11	5 (8%)
24	DD6	P	314	-	40,45,45	1.36	8 (20%)	51,67,67	1.94	10 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
25	DGD	B	842	-	61,61,67	0.97	1 (1%)	75,75,81	1.34	7 (9%)
29	A86	R	206	-	47,50,50	1.39	4 (8%)	51,76,76	1.30	8 (15%)
18	CLA	B	832	2	51,55,73	1.42	8 (15%)	60,91,113	1.11	5 (8%)
18	CLA	A	819	1	47,51,73	1.43	8 (17%)	55,86,113	1.06	5 (9%)
21	BCR	B	838	-	41,41,41	1.07	2 (4%)	56,56,56	1.20	6 (10%)
21	BCR	B	839	-	41,41,41	1.06	3 (7%)	56,56,56	1.19	6 (10%)
21	BCR	I	102	-	41,41,41	1.04	2 (4%)	56,56,56	1.20	4 (7%)
18	CLA	B	843	2	69,73,73	1.29	8 (11%)	82,113,113	0.93	3 (3%)
18	CLA	A	834	1	69,73,73	1.33	9 (13%)	82,113,113	0.93	4 (4%)
24	DD6	T	214	-	40,45,45	1.31	7 (17%)	51,67,67	1.84	10 (19%)
18	CLA	B	819	2	57,61,73	1.29	7 (12%)	67,98,113	1.01	5 (7%)
18	CLA	Q	206	13	54,59,73	1.38	7 (12%)	62,96,113	0.95	3 (4%)
21	BCR	A	843	-	41,41,41	1.09	2 (4%)	56,56,56	1.24	4 (7%)
18	CLA	T	204	17	50,54,73	1.44	7 (14%)	59,90,113	0.96	3 (5%)
18	CLA	B	827	2	53,57,73	1.38	8 (15%)	61,93,113	1.05	4 (6%)
18	CLA	A	806	1	69,73,73	1.14	7 (10%)	82,113,113	0.89	4 (4%)
18	CLA	A	831	1	49,53,73	1.36	7 (14%)	58,89,113	1.11	4 (6%)
28	KC1	Q	210	13	49,53,53	1.59	8 (16%)	61,89,89	1.21	6 (9%)
18	CLA	U	211	16	56,60,73	1.31	7 (12%)	65,97,113	1.02	5 (7%)
20	LHG	A	840	18	26,26,48	0.85	1 (3%)	29,32,54	1.32	3 (10%)
26	SQD	B	847	-	48,50,54	1.01	4 (8%)	58,61,65	1.49	9 (15%)
18	CLA	T	208	17	69,73,73	1.16	7 (10%)	82,113,113	1.10	5 (6%)
21	BCR	L	201	-	41,41,41	1.10	2 (4%)	56,56,56	1.19	5 (8%)
18	CLA	A	813	1	54,58,73	1.39	7 (12%)	64,95,113	1.11	6 (9%)
18	CLA	A	805	1	53,57,73	1.41	7 (13%)	61,93,113	1.06	4 (6%)
18	CLA	T	206	-	50,54,73	1.35	7 (14%)	59,90,113	1.15	6 (10%)
18	CLA	B	824	2	69,73,73	1.26	8 (11%)	82,113,113	1.66	11 (13%)
18	CLA	O	207	11	50,54,73	1.37	8 (16%)	59,90,113	1.09	4 (6%)
24	DD6	U	212	-	40,45,45	1.23	7 (17%)	51,67,67	1.85	13 (25%)
18	CLA	Q	213	30	61,65,73	1.24	6 (9%)	72,103,113	1.01	5 (6%)
18	CLA	Q	205	13	64,68,73	1.19	7 (10%)	76,107,113	0.99	4 (5%)
18	CLA	T	202	-	46,50,73	1.44	7 (15%)	53,85,113	1.19	4 (7%)
24	DD6	A	854	-	40,45,45	1.29	6 (15%)	51,67,67	1.54	10 (19%)
18	CLA	B	820	-	67,71,73	1.21	8 (11%)	79,110,113	1.01	6 (7%)
18	CLA	A	833	1	69,73,73	1.25	7 (10%)	82,113,113	1.02	6 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
24	DD6	S	202	-	40,45,45	1.30	9 (22%)	51,67,67	1.49	9 (17%)
18	CLA	L	203	9	69,73,73	1.20	7 (10%)	82,113,113	0.90	4 (4%)
18	CLA	B	810	2	58,62,73	1.43	9 (15%)	71,100,113	0.90	3 (4%)
18	CLA	B	831	2	69,73,73	1.22	7 (10%)	82,113,113	0.88	3 (3%)
18	CLA	A	824	1	66,70,73	1.21	7 (10%)	78,109,113	1.06	5 (6%)
28	KC1	P	305	12	49,53,53	1.95	10 (20%)	61,89,89	1.12	6 (9%)
19	PQN	B	836	-	34,34,34	0.43	0	43,45,45	0.50	1 (2%)
18	CLA	P	313	12	49,53,73	1.42	6 (12%)	58,89,113	1.13	5 (8%)
21	BCR	A	842	-	41,41,41	1.08	2 (4%)	56,56,56	1.14	4 (7%)
18	CLA	B	804	2	49,53,73	1.46	8 (16%)	58,89,113	1.02	3 (5%)
18	CLA	A	835	1	69,73,73	1.24	7 (10%)	82,113,113	0.94	4 (4%)
22	CL0	A	848	1	58,73,73	0.84	2 (3%)	60,113,113	1.73	8 (13%)
21	BCR	J	104	-	41,41,41	1.06	2 (4%)	56,56,56	1.14	4 (7%)
24	DD6	U	214	-	24,26,45	1.50	5 (20%)	29,35,67	1.57	6 (20%)
28	KC1	P	302	12	49,53,53	1.69	10 (20%)	61,89,89	1.10	6 (9%)
18	CLA	S	201	30	69,73,73	1.13	6 (8%)	82,113,113	1.07	7 (8%)
27	LMG	J	102	-	39,39,55	0.86	0	47,47,63	1.22	3 (6%)
18	CLA	Q	216	30	69,73,73	1.15	8 (11%)	82,113,113	0.94	4 (4%)
18	CLA	Q	203	13	52,56,73	1.33	6 (11%)	61,92,113	1.08	5 (8%)
24	DD6	Q	202	-	40,45,45	1.28	7 (17%)	51,67,67	1.67	12 (23%)
28	KC1	S	208	15	49,53,53	1.61	9 (18%)	61,89,89	1.05	3 (4%)
29	A86	U	202	-	47,50,50	1.40	5 (10%)	51,76,76	1.32	7 (13%)
18	CLA	O	204	11	69,73,73	1.21	9 (13%)	82,113,113	0.93	5 (6%)
21	BCR	A	841	-	41,41,41	1.04	2 (4%)	56,56,56	1.26	4 (7%)
18	CLA	P	306	12	69,73,73	1.17	7 (10%)	82,113,113	0.99	6 (7%)
18	CLA	A	802	1	59,63,73	1.32	7 (11%)	70,101,113	0.97	3 (4%)
18	CLA	A	823	1	69,73,73	1.24	8 (11%)	82,113,113	0.87	3 (3%)
21	BCR	I	101	-	41,41,41	1.10	2 (4%)	56,56,56	1.36	8 (14%)
18	CLA	B	830	2	62,66,73	1.29	8 (12%)	73,104,113	1.01	6 (8%)
18	CLA	B	809	2	69,73,73	1.19	8 (11%)	82,113,113	0.94	5 (6%)
18	CLA	F	803	6	50,54,73	1.38	7 (14%)	59,90,113	1.10	4 (6%)
24	DD6	O	201	-	40,45,45	1.42	8 (20%)	51,67,67	1.57	9 (17%)
18	CLA	A	847	1	69,73,73	1.21	9 (13%)	82,113,113	1.03	5 (6%)
18	CLA	A	822	30	69,73,73	1.14	5 (7%)	82,113,113	0.94	4 (4%)
18	CLA	B	844	2	69,73,73	1.27	7 (10%)	82,113,113	0.90	5 (6%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
24	DD6	S	203	-	40,45,45	1.36	8 (20%)	51,67,67	1.86	11 (21%)
18	CLA	O	208	11	64,68,73	1.17	8 (12%)	76,107,113	1.02	5 (6%)
18	CLA	A	817	1	49,53,73	1.42	7 (14%)	58,89,113	1.10	4 (6%)
18	CLA	S	207	15	56,60,73	1.26	7 (12%)	65,97,113	1.12	5 (7%)
18	CLA	L	202	9	53,57,73	1.36	7 (13%)	61,93,113	1.39	6 (9%)
18	CLA	P	308	-	56,60,73	1.28	7 (12%)	65,97,113	1.02	5 (7%)
18	CLA	A	812	1	49,53,73	1.47	7 (14%)	58,89,113	1.08	4 (6%)
21	BCR	B	837	-	41,41,41	1.10	2 (4%)	56,56,56	1.14	4 (7%)
18	CLA	A	852	1	69,73,73	1.17	8 (11%)	82,113,113	0.93	5 (6%)
18	CLA	A	836	30	69,73,73	1.18	9 (13%)	82,113,113	1.00	7 (8%)
28	KC1	O	210	11	49,53,53	1.61	10 (20%)	61,89,89	1.11	4 (6%)
18	CLA	A	849	1	69,73,73	1.23	6 (8%)	82,113,113	0.89	3 (3%)
18	CLA	A	810	1	58,62,73	1.27	6 (10%)	68,99,113	1.02	5 (7%)
18	CLA	A	809	1	66,70,73	1.23	7 (10%)	78,109,113	0.87	3 (3%)
18	CLA	B	812	2	58,62,73	1.34	8 (13%)	68,99,113	0.99	4 (5%)
18	CLA	B	825	2	69,73,73	1.27	8 (11%)	82,113,113	0.92	5 (6%)
23	SF4	A	850	2,1	0,12,12	-	-	-	-	-
18	CLA	T	212	-	51,55,73	1.34	6 (11%)	60,91,113	1.14	5 (8%)
24	DD6	J	101	-	40,45,45	1.24	7 (17%)	51,67,67	1.59	9 (17%)
20	LHG	A	839	-	47,47,48	0.66	0	50,53,54	1.21	5 (10%)
18	CLA	B	806	2	69,73,73	1.25	8 (11%)	82,113,113	0.89	4 (4%)
18	CLA	F	802	30	52,56,73	1.40	7 (13%)	61,92,113	1.16	5 (8%)
18	CLA	A	815	1	69,73,73	1.22	7 (10%)	82,113,113	0.95	4 (4%)
18	CLA	B	802	2	69,73,73	1.27	7 (10%)	82,113,113	0.80	3 (3%)
18	CLA	B	834	2	69,73,73	1.26	9 (13%)	82,113,113	0.98	4 (4%)
18	CLA	B	826	2	54,58,73	1.42	8 (14%)	64,95,113	1.02	4 (6%)
18	CLA	Q	209	13	69,73,73	1.14	6 (8%)	82,113,113	0.95	3 (3%)
24	DD6	P	304	-	40,45,45	1.31	6 (15%)	51,67,67	1.42	10 (19%)
18	CLA	A	825	1	69,73,73	1.25	8 (11%)	82,113,113	0.87	3 (3%)
24	DD6	S	209	-	40,45,45	1.27	5 (12%)	51,67,67	1.44	8 (15%)
28	KC1	T	209	17	49,53,53	1.67	9 (18%)	61,89,89	1.00	2 (3%)
21	BCR	M	101	-	41,41,41	1.07	2 (4%)	56,56,56	1.31	7 (12%)
18	CLA	A	818	30	69,73,73	1.18	7 (10%)	82,113,113	0.90	3 (3%)
18	CLA	T	211	17	50,54,73	1.40	7 (14%)	59,90,113	1.21	7 (11%)
18	CLA	A	804	1	69,73,73	1.16	8 (11%)	82,113,113	0.99	6 (7%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
24	DD6	S	213	-	40,45,45	1.72	9 (22%)	51,67,67	1.63	11 (21%)
29	A86	P	303	-	47,50,50	1.39	6 (12%)	51,76,76	1.46	7 (13%)
18	CLA	U	205	16	69,73,73	1.17	7 (10%)	82,113,113	0.91	3 (3%)

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
18	CLA	A	846	1	-	9/33/109/115	-
21	BCR	F	804	-	-	9/29/63/63	0/2/2/2
18	CLA	A	826	1	1/1/15/20	3/39/115/115	-
18	CLA	A	845	30	1/1/15/20	6/39/115/115	-
18	CLA	O	209	11	-	12/39/115/115	-
18	CLA	A	853	30	1/1/15/20	2/39/115/115	-
24	DD6	Q	215	-	-	8/26/80/80	0/3/3/3
18	CLA	T	210	17	-	0/10/86/115	-
23	SF4	C	102	3	-	-	0/6/5/5
18	CLA	B	807	2	-	2/39/115/115	-
18	CLA	B	805	2	1/1/15/20	8/39/115/115	-
18	CLA	O	205	11	1/1/15/20	6/39/115/115	-
18	CLA	U	206	16	1/1/11/20	2/15/91/115	-
18	CLA	U	207	-	-	6/39/115/115	-
18	CLA	A	838	20	1/1/12/20	4/24/100/115	-
18	CLA	P	312	12	1/1/10/20	2/10/86/115	-
18	CLA	A	851	1	1/1/15/20	11/39/115/115	-
18	CLA	U	204	-	1/1/14/20	9/35/111/115	-
18	CLA	S	215	15	1/1/15/20	4/39/115/115	-
18	CLA	O	203	-	1/1/11/20	1/15/91/115	-
18	CLA	U	209	16	-	0/12/88/115	-
18	CLA	P	315	12	1/1/11/20	1/18/94/115	-
18	CLA	B	828	2	-	2/31/107/115	-
18	CLA	U	210	16	-	9/39/115/115	-
18	CLA	B	833	-	1/1/15/20	3/39/115/115	-
24	DD6	P	317	-	-	10/15/46/80	0/2/2/3
27	LMG	S	211	-	-	25/44/64/70	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
27	LMG	B	849	-	-	8/19/39/70	0/1/1/1
24	DD6	O	215	-	-	12/26/80/80	0/3/3/3
18	CLA	B	811	2	-	0/27/103/115	-
18	CLA	B	803	-	1/1/15/20	6/39/115/115	-
26	SQD	M	102	-	-	15/41/61/69	0/1/1/1
18	CLA	O	206	-	1/1/15/20	0/39/115/115	-
18	CLA	A	803	1	1/1/15/20	4/39/115/115	-
18	CLA	T	203	17	1/1/10/20	1/10/86/115	-
18	CLA	O	202	11	-	3/13/89/115	-
18	CLA	A	828	1	1/1/15/20	4/39/115/115	-
18	CLA	A	832	1	1/1/12/20	0/23/99/115	-
18	CLA	S	214	30	1/1/15/20	1/39/115/115	-
18	CLA	B	848	2	1/1/12/20	6/21/97/115	-
18	CLA	A	827	1	-	0/21/97/115	-
18	CLA	A	807	1	-	6/39/115/115	-
24	DD6	S	212	-	-	16/26/80/80	0/3/3/3
18	CLA	A	830	1	-	0/21/97/115	-
24	DD6	O	213	-	-	5/26/80/80	0/3/3/3
18	CLA	R	205	14	1/1/15/20	11/39/115/115	-
18	CLA	A	808	1	1/1/13/20	5/29/105/115	-
18	CLA	Q	211	13	1/1/10/20	4/10/86/115	-
24	DD6	O	214	-	-	8/26/80/80	0/3/3/3
18	CLA	A	821	-	1/1/15/20	8/39/115/115	-
18	CLA	T	207	17	1/1/10/20	0/12/88/115	-
28	KC1	P	311	12	-	4/12/68/71	-
20	LHG	R	201	-	-	30/53/53/53	-
29	A86	R	204	-	-	1/30/84/90	0/3/3/3
21	BCR	R	203	-	-	10/27/61/63	0/2/2/2
24	DD6	U	203	-	-	16/26/80/80	0/3/3/3
23	SF4	C	101	3	-	-	0/6/5/5
18	CLA	A	829	1	1/1/15/20	4/39/115/115	-
18	CLA	J	103	8	1/1/10/20	2/12/88/115	-
18	CLA	U	208	16	1/1/11/20	2/17/93/115	-
19	PQN	A	837	-	-	7/23/43/43	0/2/2/2
24	DD6	P	316	-	-	14/26/80/80	0/3/3/3
18	CLA	P	310	12	-	4/21/97/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
18	CLA	B	814	2	-	3/27/103/115	-
18	CLA	B	816	2	1/1/14/20	2/33/109/115	-
21	BCR	F	801	-	-	5/29/63/63	0/2/2/2
18	CLA	B	829	30	1/1/15/20	1/39/115/115	-
18	CLA	P	307	12	1/1/13/20	11/29/105/115	-
18	CLA	S	204	15	1/1/11/20	0/17/93/115	-
18	CLA	B	815	2	-	7/32/108/115	-
18	CLA	B	818	2	-	1/17/93/115	-
24	DD6	T	213	-	-	12/26/80/80	0/3/3/3
21	BCR	A	844	-	-	9/29/63/63	0/2/2/2
18	CLA	B	808	2	1/1/15/20	5/39/115/115	-
18	CLA	R	202	30	1/1/11/20	0/15/91/115	-
24	DD6	O	212	-	-	11/26/80/80	0/3/3/3
28	KC1	U	213	16	-	2/15/71/71	-
27	LMG	P	301	-	-	12/29/49/70	0/1/1/1
29	A86	T	201	-	-	18/34/90/90	0/3/3/3
18	CLA	A	814	30	1/1/11/20	3/15/91/115	-
18	CLA	A	801	-	-	4/39/115/115	-
18	CLA	B	822	2	1/1/15/20	0/39/115/115	-
18	CLA	T	205	-	1/1/13/20	10/30/106/115	-
29	A86	Q	201	-	-	23/34/90/90	0/3/3/3
18	CLA	B	817	-	1/1/15/20	7/39/115/115	-
18	CLA	B	821	-	1/1/14/20	4/38/114/115	-
21	BCR	B	840	-	-	5/29/63/63	0/2/2/2
18	CLA	B	845	2	1/1/15/20	2/39/115/115	-
21	BCR	L	205	-	-	8/29/63/63	0/2/2/2
28	KC1	S	210	15	-	5/15/71/71	-
18	CLA	L	204	-	1/1/12/20	8/21/97/115	-
18	CLA	P	309	30	-	2/18/94/115	-
18	CLA	S	205	15	1/1/11/20	5/17/93/115	-
27	LMG	Q	217	-	-	23/50/70/70	0/1/1/1
18	CLA	A	811	1	1/1/15/20	4/39/115/115	-
18	CLA	Q	208	13	1/1/12/20	1/21/97/115	-
18	CLA	S	206	15	1/1/11/20	2/15/91/115	-
18	CLA	B	846	2	1/1/15/20	7/39/115/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
18	CLA	B	813	2	-	5/20/96/115	-
29	A86	Q	214	-	-	8/34/90/90	0/3/3/3
18	CLA	Q	212	13	-	15/39/115/115	-
18	CLA	O	211	11	-	0/10/86/115	-
18	CLA	A	816	1	1/1/15/20	0/39/115/115	-
18	CLA	Q	207	13	1/1/11/20	4/17/93/115	-
27	LMG	U	201	-	-	18/27/47/70	0/1/1/1
18	CLA	Q	204	-	1/1/14/20	8/35/111/115	-
18	CLA	B	823	2	1/1/15/20	6/39/115/115	-
18	CLA	B	801	-	1/1/15/20	1/39/115/115	-
18	CLA	B	835	30	1/1/15/20	1/39/115/115	-
18	CLA	A	820	1	1/1/12/20	3/23/99/115	-
21	BCR	B	841	-	-	3/29/63/63	0/2/2/2
24	DD6	P	314	-	-	13/26/80/80	0/3/3/3
25	DGD	B	842	-	-	20/49/89/95	0/2/2/2
29	A86	R	206	-	-	4/34/90/90	0/3/3/3
18	CLA	B	832	2	1/1/11/20	2/18/94/115	-
18	CLA	A	819	1	-	0/13/89/115	-
21	BCR	B	838	-	-	7/29/63/63	0/2/2/2
21	BCR	B	839	-	-	2/29/63/63	0/2/2/2
21	BCR	I	102	-	-	9/29/63/63	0/2/2/2
18	CLA	B	843	2	1/1/15/20	5/39/115/115	-
18	CLA	A	834	1	-	7/39/115/115	-
24	DD6	T	214	-	-	12/26/80/80	0/3/3/3
18	CLA	B	819	2	-	5/25/101/115	-
18	CLA	Q	206	13	1/1/12/20	7/22/98/115	-
21	BCR	A	843	-	-	4/29/63/63	0/2/2/2
18	CLA	T	204	17	1/1/11/20	3/17/93/115	-
18	CLA	B	827	2	1/1/11/20	2/20/96/115	-
18	CLA	A	806	1	1/1/15/20	5/39/115/115	-
18	CLA	A	831	1	1/1/11/20	0/15/91/115	-
28	KC1	Q	210	13	-	2/15/71/71	-
18	CLA	U	211	16	1/1/12/20	3/24/100/115	-
20	LHG	A	840	18	-	7/31/31/53	-
26	SQD	B	847	-	-	23/45/65/69	0/1/1/1

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
18	CLA	T	208	17	-	12/39/115/115	-
21	BCR	L	201	-	-	10/29/63/63	0/2/2/2
18	CLA	A	813	1	-	1/21/97/115	-
18	CLA	T	206	-	1/1/11/20	5/17/93/115	-
18	CLA	A	805	1	-	2/20/96/115	-
18	CLA	B	824	2	-	9/39/115/115	-
18	CLA	O	207	11	1/1/11/20	1/17/93/115	-
24	DD6	U	212	-	-	11/26/80/80	0/3/3/3
18	CLA	Q	213	30	1/1/13/20	6/30/106/115	-
18	CLA	Q	205	13	1/1/14/20	5/33/109/115	-
18	CLA	T	202	-	1/1/10/20	1/12/88/115	-
24	DD6	A	854	-	-	13/26/80/80	0/3/3/3
18	CLA	B	820	-	1/1/14/20	5/37/113/115	-
18	CLA	A	833	1	1/1/15/20	7/39/115/115	-
24	DD6	S	202	-	-	8/26/80/80	0/3/3/3
18	CLA	L	203	9	-	0/39/115/115	-
18	CLA	B	810	2	-	2/25/101/115	-
18	CLA	B	831	2	1/1/15/20	4/39/115/115	-
18	CLA	A	824	1	1/1/14/20	5/36/112/115	-
28	KC1	P	305	12	-	3/15/71/71	-
19	PQN	B	836	-	-	1/23/43/43	0/2/2/2
18	CLA	P	313	12	1/1/11/20	0/15/91/115	-
21	BCR	A	842	-	-	9/29/63/63	0/2/2/2
18	CLA	B	804	2	1/1/11/20	5/15/91/115	-
18	CLA	A	835	1	1/1/15/20	4/39/115/115	-
22	CL0	A	848	1	-	3/37/135/135	-
21	BCR	J	104	-	-	11/29/63/63	0/2/2/2
24	DD6	U	214	-	-	7/14/37/80	0/1/1/3
28	KC1	P	302	12	-	2/15/71/71	-
18	CLA	S	201	30	-	9/39/115/115	-
27	LMG	J	102	-	-	26/34/54/70	0/1/1/1
18	CLA	Q	216	30	-	11/39/115/115	-
18	CLA	Q	203	13	-	5/19/95/115	-
24	DD6	Q	202	-	-	6/26/80/80	0/3/3/3
28	KC1	S	208	15	-	1/15/71/71	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
29	A86	U	202	-	-	13/34/90/90	0/3/3/3
18	CLA	O	204	11	1/1/15/20	3/39/115/115	-
21	BCR	A	841	-	-	7/29/63/63	0/2/2/2
18	CLA	P	306	12	1/1/15/20	7/39/115/115	-
18	CLA	A	802	1	1/1/13/20	5/27/103/115	-
18	CLA	A	823	1	1/1/15/20	8/39/115/115	-
21	BCR	I	101	-	-	6/29/63/63	0/2/2/2
18	CLA	B	830	2	1/1/13/20	2/31/107/115	-
18	CLA	B	809	2	1/1/15/20	7/39/115/115	-
18	CLA	F	803	6	1/1/11/20	4/17/93/115	-
24	DD6	O	201	-	-	3/26/80/80	0/3/3/3
18	CLA	A	847	1	-	7/39/115/115	-
18	CLA	A	822	30	1/1/15/20	1/39/115/115	-
18	CLA	B	844	2	1/1/15/20	2/39/115/115	-
24	DD6	S	203	-	-	8/26/80/80	0/3/3/3
18	CLA	O	208	11	1/1/14/20	3/33/109/115	-
18	CLA	A	817	1	1/1/11/20	0/15/91/115	-
18	CLA	S	207	15	-	3/24/100/115	-
18	CLA	L	202	9	-	7/20/96/115	-
18	CLA	P	308	-	1/1/12/20	3/24/100/115	-
18	CLA	A	812	1	1/1/11/20	0/15/91/115	-
21	BCR	B	837	-	-	9/29/63/63	0/2/2/2
18	CLA	A	852	1	1/1/15/20	4/39/115/115	-
18	CLA	A	836	30	1/1/15/20	3/39/115/115	-
28	KC1	O	210	11	-	0/15/71/71	-
18	CLA	A	849	1	-	4/39/115/115	-
18	CLA	A	810	1	1/1/12/20	0/26/102/115	-
18	CLA	A	809	1	-	2/36/112/115	-
18	CLA	B	812	2	-	1/26/102/115	-
18	CLA	B	825	2	-	7/39/115/115	-
24	DD6	J	101	-	-	7/26/80/80	0/3/3/3
18	CLA	T	212	-	1/1/11/20	4/18/94/115	-
23	SF4	A	850	2,1	-	-	0/6/5/5
20	LHG	A	839	-	-	19/52/52/53	-
18	CLA	B	806	2	1/1/15/20	8/39/115/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
18	CLA	F	802	30	1/1/11/20	5/19/95/115	-
18	CLA	A	815	1	1/1/15/20	12/39/115/115	-
18	CLA	B	802	2	1/1/15/20	1/39/115/115	-
18	CLA	B	834	2	-	7/39/115/115	-
18	CLA	B	826	2	-	0/21/97/115	-
18	CLA	Q	209	13	-	8/39/115/115	-
24	DD6	P	304	-	-	8/26/80/80	0/3/3/3
18	CLA	A	825	1	-	5/39/115/115	-
24	DD6	S	209	-	-	6/26/80/80	0/3/3/3
28	KC1	T	209	17	-	2/15/71/71	-
21	BCR	M	101	-	-	4/29/63/63	0/2/2/2
18	CLA	A	818	30	1/1/15/20	3/39/115/115	-
18	CLA	T	211	17	-	4/17/93/115	-
18	CLA	A	804	1	1/1/15/20	11/39/115/115	-
24	DD6	S	213	-	-	10/26/80/80	0/3/3/3
29	A86	P	303	-	-	14/34/90/90	0/3/3/3
18	CLA	U	205	16	-	7/39/115/115	-

The worst 5 of 1399 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
29	Q	214	A86	C13-C11	-6.85	1.36	1.49
24	S	213	DD6	C28-C27	-6.73	1.46	1.50
29	T	201	A86	C13-C11	-6.70	1.37	1.49
29	U	202	A86	C13-C11	-6.59	1.37	1.49
29	R	206	A86	C13-C11	-6.57	1.37	1.49

The worst 5 of 1171 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
22	A	848	CL0	C1B-CHB-C4A	8.99	127.11	121.32
18	B	824	CLA	C1-O2A-CGA	8.56	137.38	116.65
24	U	212	DD6	C3-C4-C5	7.16	138.16	123.52
24	P	317	DD6	C3-C4-C5	5.92	131.04	123.29
24	S	203	DD6	C3-C4-C5	5.83	135.44	123.52

5 of 98 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
18	A	802	CLA	ND
18	A	803	CLA	ND
18	A	804	CLA	ND
18	A	806	CLA	ND
18	A	808	CLA	ND

5 of 1291 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
18	A	804	CLA	C1A-C2A-CAA-CBA
18	A	804	CLA	C3A-C2A-CAA-CBA
18	A	806	CLA	CHA-CBD-CGD-O1D
18	A	806	CLA	CHA-CBD-CGD-O2D
18	A	808	CLA	CBA-CGA-O2A-C1

There are no ring outliers.

151 monomers are involved in 253 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
18	A	846	CLA	1	0
21	F	804	BCR	3	0
18	A	826	CLA	2	0
18	A	845	CLA	3	0
18	O	209	CLA	2	0
18	A	853	CLA	2	0
18	T	210	CLA	1	0
18	B	807	CLA	4	0
18	O	205	CLA	2	0
18	U	206	CLA	2	0
18	U	207	CLA	3	0
18	A	838	CLA	1	0
18	A	851	CLA	1	0
18	S	215	CLA	2	0
18	O	203	CLA	1	0
18	U	209	CLA	1	0
18	B	828	CLA	1	0
18	P	315	CLA	1	0
18	U	210	CLA	4	0
18	B	833	CLA	2	0
27	S	211	LMG	6	0
27	B	849	LMG	2	0
18	B	803	CLA	3	0
26	M	102	SQD	3	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
18	O	206	CLA	1	0
18	A	803	CLA	2	0
18	O	202	CLA	2	0
18	A	828	CLA	2	0
18	S	214	CLA	4	0
18	B	848	CLA	3	0
18	A	807	CLA	2	0
18	A	830	CLA	1	0
18	R	205	CLA	3	0
18	A	808	CLA	1	0
18	A	821	CLA	1	0
18	T	207	CLA	2	0
20	R	201	LHG	2	0
21	R	203	BCR	3	0
29	R	204	A86	1	0
18	J	103	CLA	2	0
18	U	208	CLA	1	0
19	A	837	PQN	2	0
18	B	816	CLA	2	0
21	F	801	BCR	1	0
18	B	829	CLA	2	0
18	P	307	CLA	1	0
18	B	815	CLA	3	0
18	S	204	CLA	2	0
18	B	818	CLA	1	0
21	A	844	BCR	1	0
18	B	808	CLA	4	0
18	R	202	CLA	1	0
27	P	301	LMG	1	0
18	A	814	CLA	2	0
18	A	801	CLA	4	0
18	B	822	CLA	1	0
18	T	205	CLA	2	0
18	B	817	CLA	2	0
18	B	821	CLA	4	0
21	B	840	BCR	1	0
18	B	845	CLA	1	0
21	L	205	BCR	4	0
27	Q	217	LMG	6	0
18	A	811	CLA	2	0
18	Q	208	CLA	5	0
18	S	206	CLA	2	0

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Mol	Chain	Res	Type	Clashes	Symm-Clashes
18	B	846	CLA	7	0
18	Q	212	CLA	1	0
18	A	816	CLA	2	0
18	Q	207	CLA	1	0
27	U	201	LMG	2	0
18	B	801	CLA	1	0
18	B	835	CLA	8	0
18	A	820	CLA	1	0
21	B	841	BCR	2	0
25	B	842	DGD	5	0
29	R	206	A86	1	0
21	B	838	BCR	1	0
21	B	839	BCR	1	0
21	I	102	BCR	2	0
18	B	843	CLA	3	0
18	A	834	CLA	3	0
18	B	819	CLA	5	0
18	Q	206	CLA	2	0
21	A	843	BCR	2	0
18	B	827	CLA	4	0
18	A	831	CLA	2	0
18	U	211	CLA	3	0
20	A	840	LHG	1	0
26	B	847	SQD	3	0
18	T	208	CLA	3	0
21	L	201	BCR	2	0
18	A	813	CLA	1	0
18	B	824	CLA	3	0
18	O	207	CLA	2	0
18	Q	205	CLA	1	0
18	T	202	CLA	2	0
18	B	820	CLA	3	0
18	A	833	CLA	5	0
18	L	203	CLA	2	0
18	B	810	CLA	1	0
18	B	831	CLA	3	0
18	A	824	CLA	3	0
19	B	836	PQN	1	0
18	P	313	CLA	1	0
21	A	842	BCR	1	0
18	A	835	CLA	3	0
22	A	848	CL0	3	0

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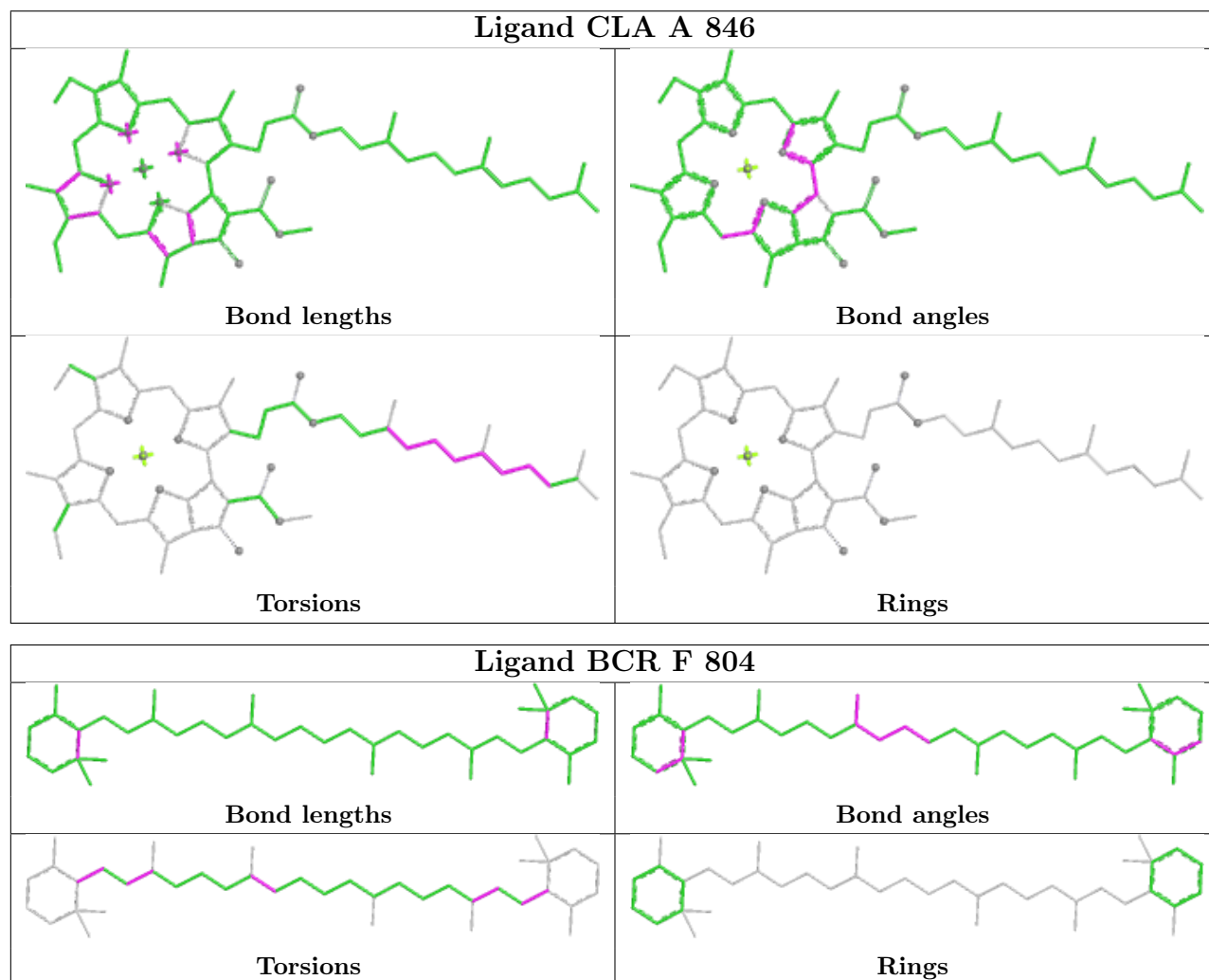
Mol	Chain	Res	Type	Clashes	Symm-Clashes
21	J	104	BCR	1	0
18	S	201	CLA	4	0
27	J	102	LMG	2	0
18	Q	216	CLA	2	0
18	Q	203	CLA	1	0
29	U	202	A86	1	0
18	O	204	CLA	1	0
21	A	841	BCR	1	0
18	P	306	CLA	1	0
18	A	823	CLA	3	0
21	I	101	BCR	2	0
18	B	830	CLA	2	0
18	F	803	CLA	1	0
18	A	847	CLA	3	0
18	A	822	CLA	1	0
18	B	844	CLA	2	0
18	O	208	CLA	2	0
18	A	817	CLA	1	0
18	S	207	CLA	2	0
18	L	202	CLA	2	0
21	B	837	BCR	2	0
18	A	852	CLA	4	0
18	A	836	CLA	2	0
28	O	210	KC1	1	0
18	A	849	CLA	1	0
18	A	810	CLA	1	0
18	B	812	CLA	1	0
18	B	825	CLA	1	0
18	T	212	CLA	1	0
20	A	839	LHG	1	0
18	B	806	CLA	1	0
18	A	815	CLA	2	0
18	B	802	CLA	2	0
18	B	834	CLA	3	0
18	Q	209	CLA	1	0
24	P	304	DD6	1	0
18	A	825	CLA	1	0
28	T	209	KC1	2	0
21	M	101	BCR	4	0
18	A	818	CLA	2	0
18	T	211	CLA	2	0
18	A	804	CLA	1	0

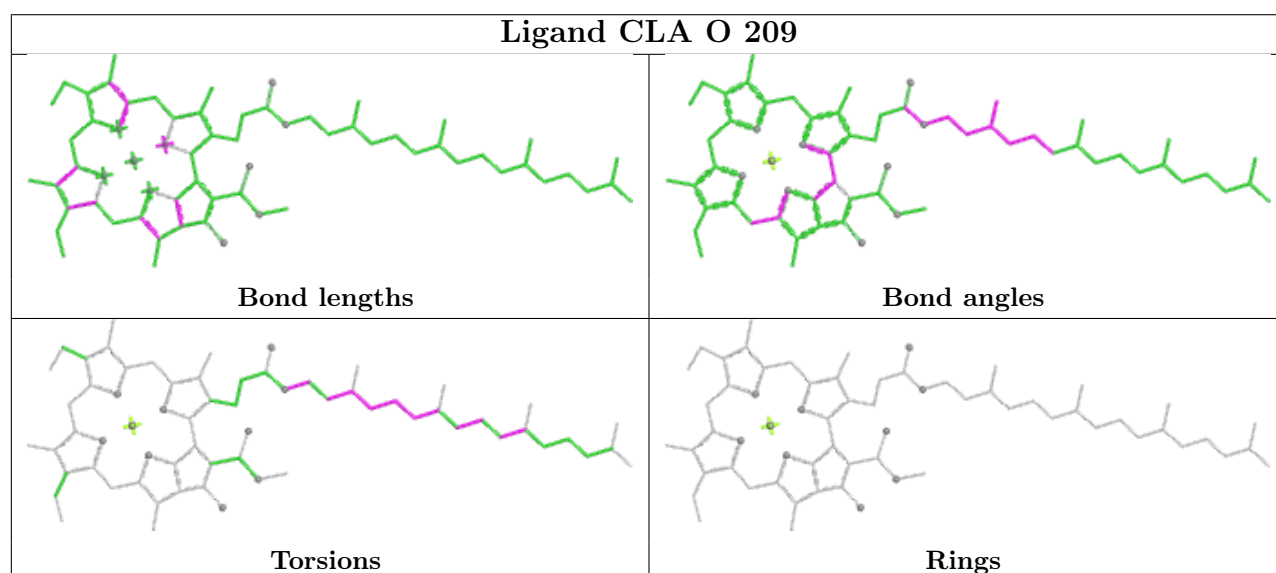
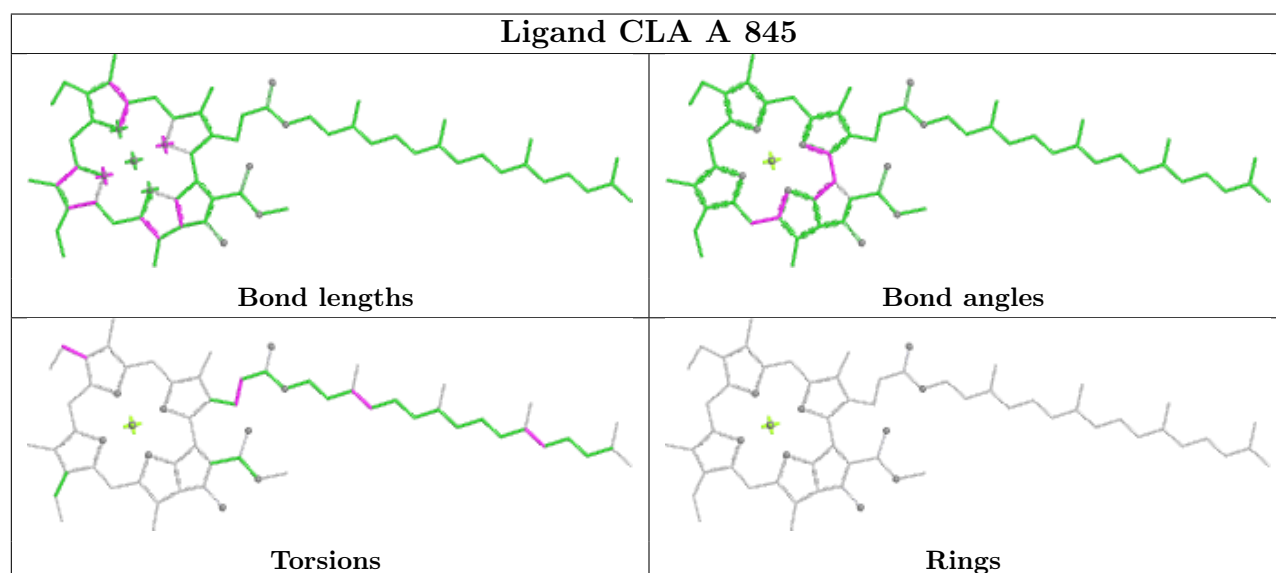
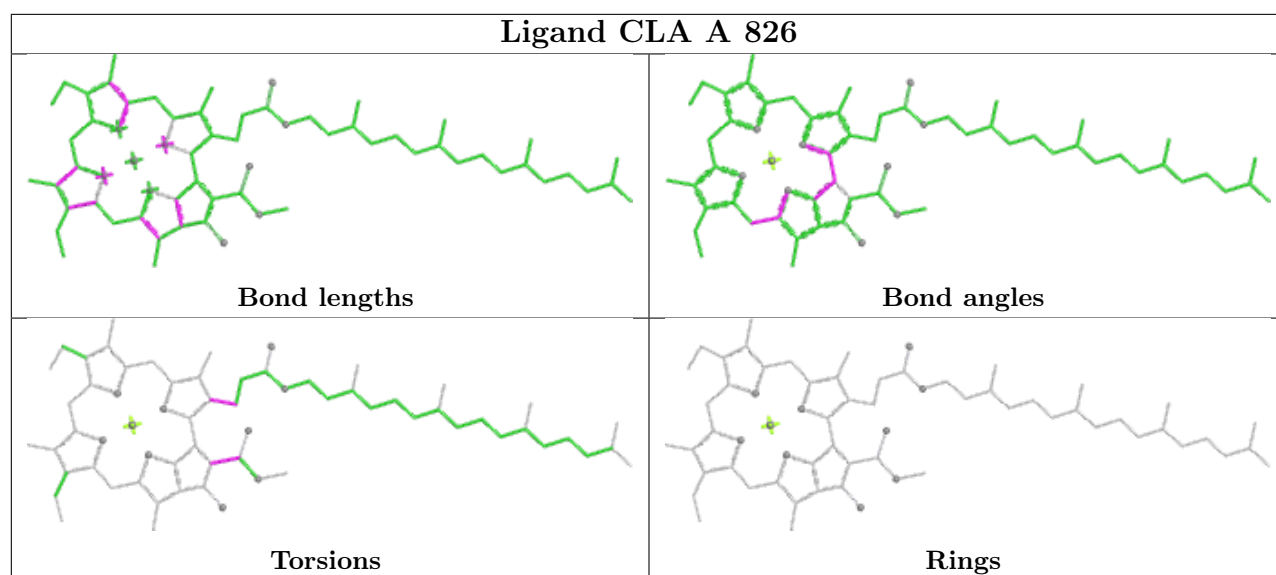
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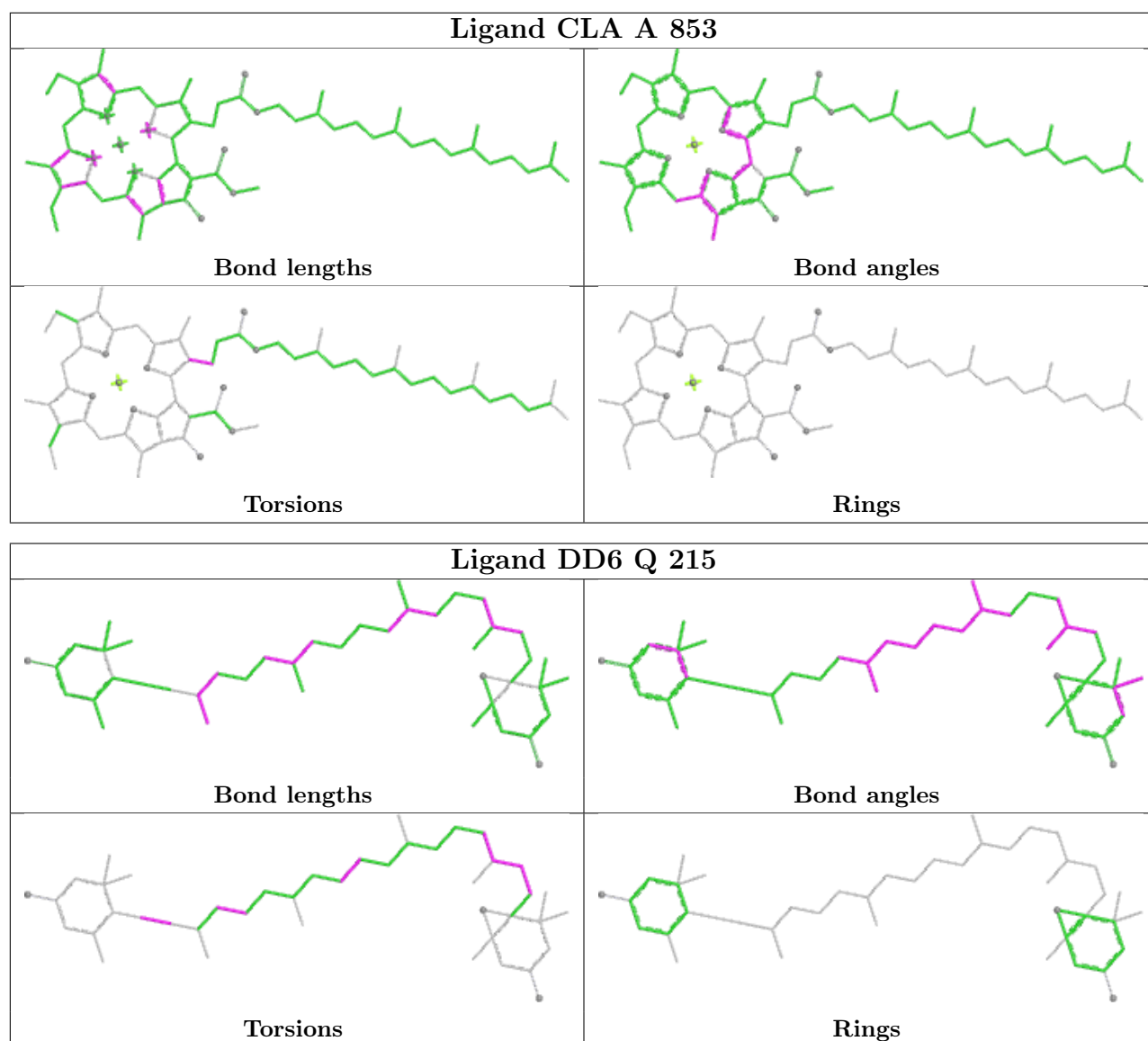
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Mol	Chain	Res	Type	Clashes	Symm-Clashes
18	U	205	CLA	1	0

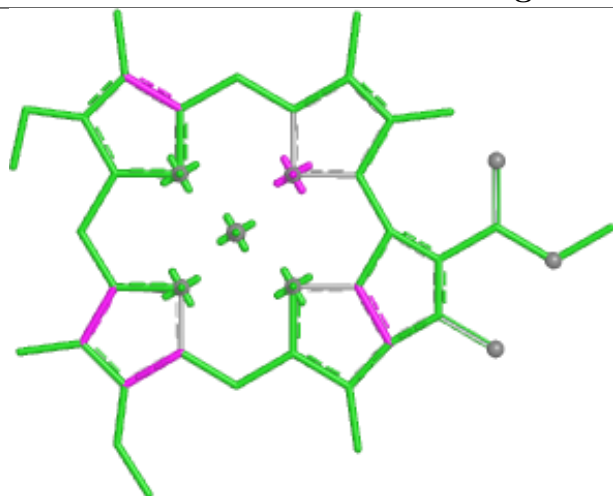
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



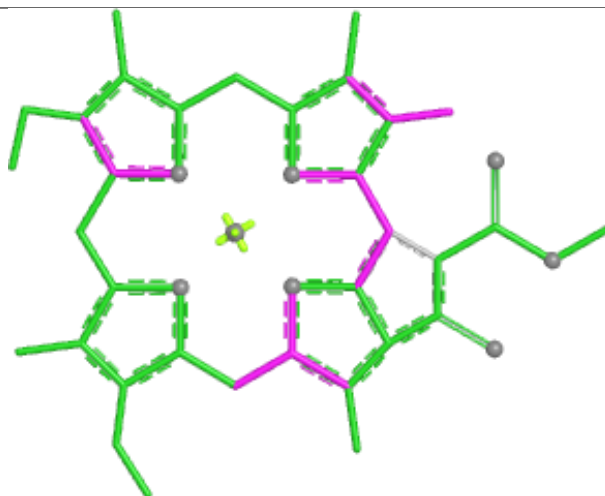




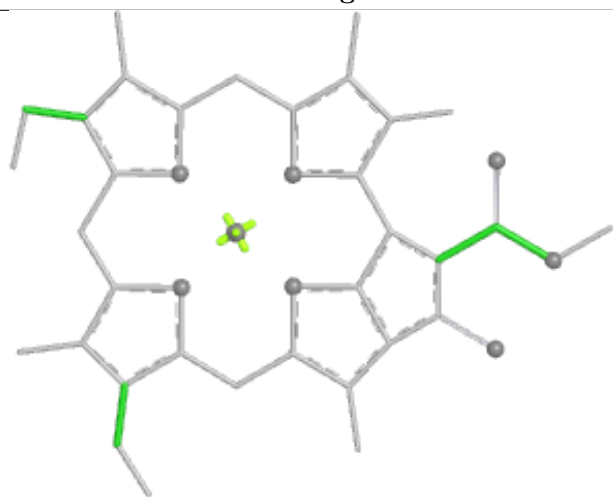
Ligand CLA T 210



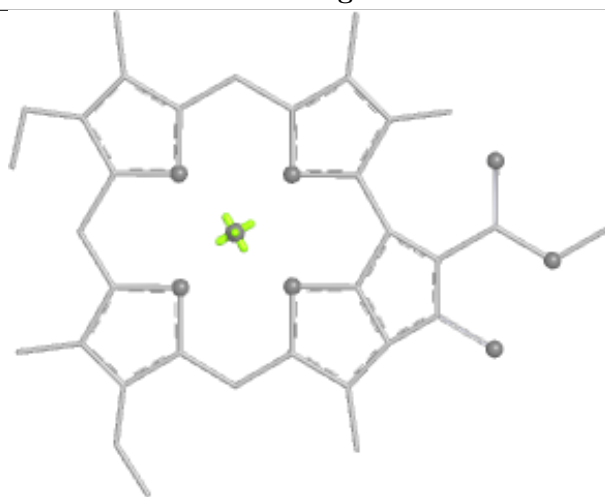
Bond lengths



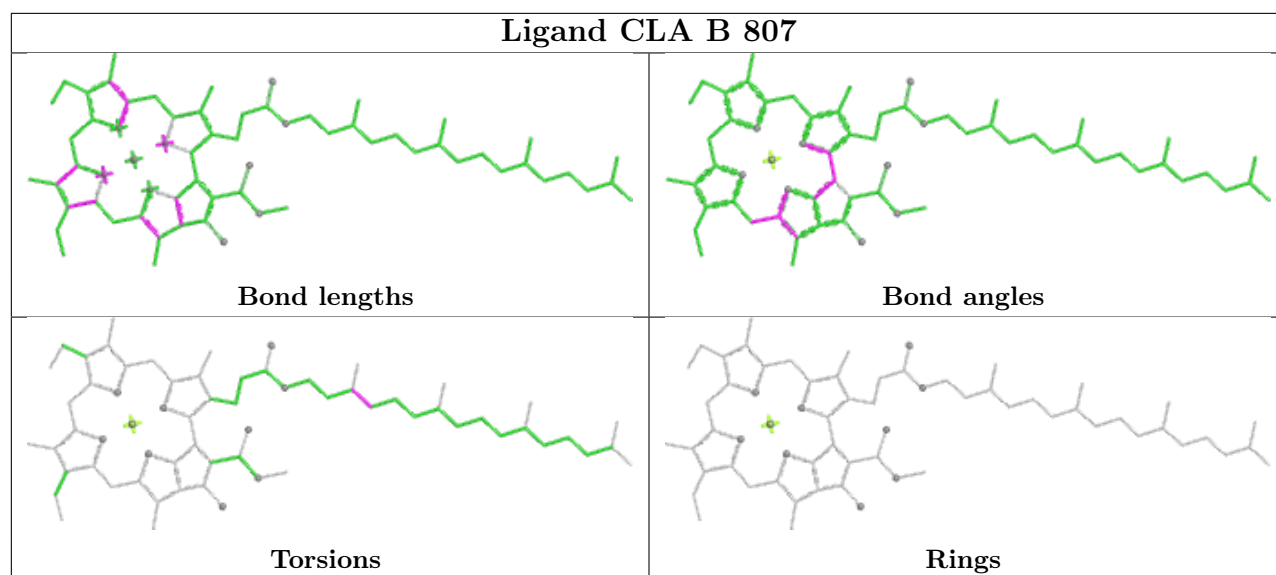
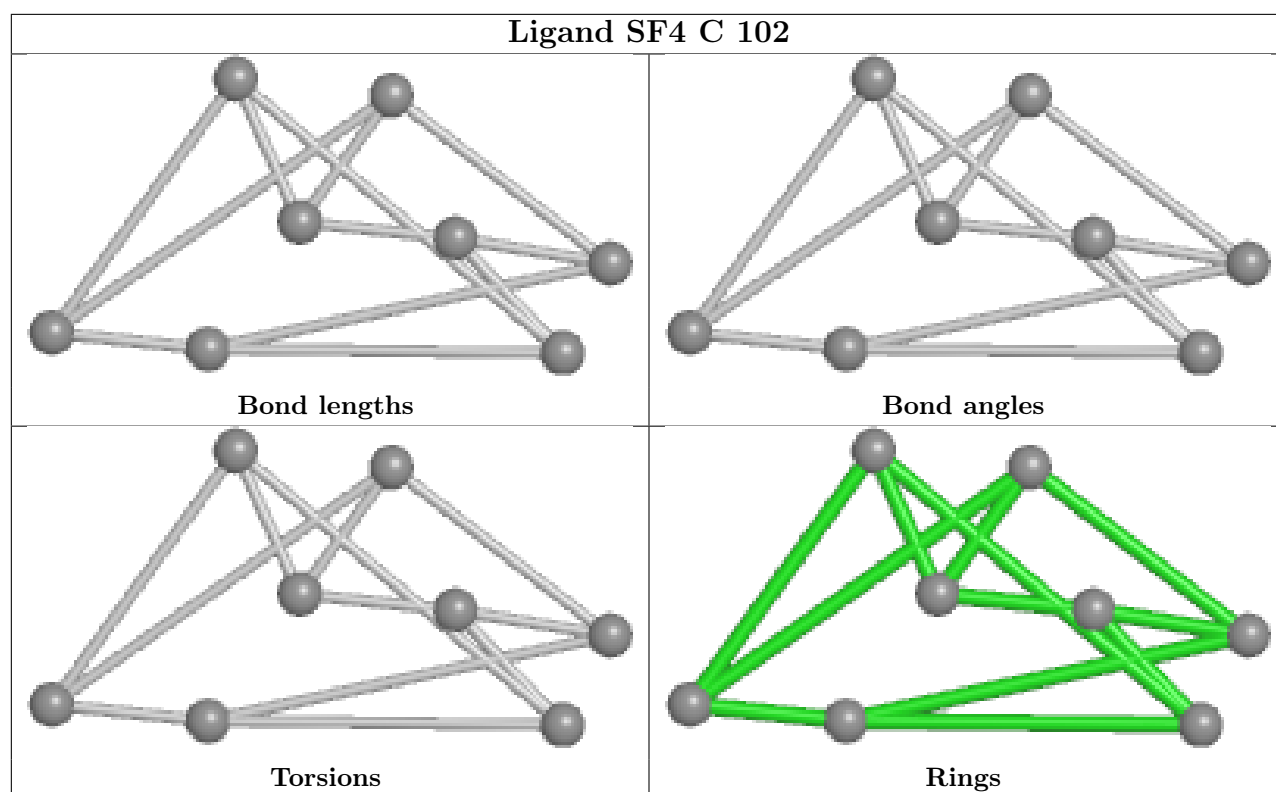
Bond angles

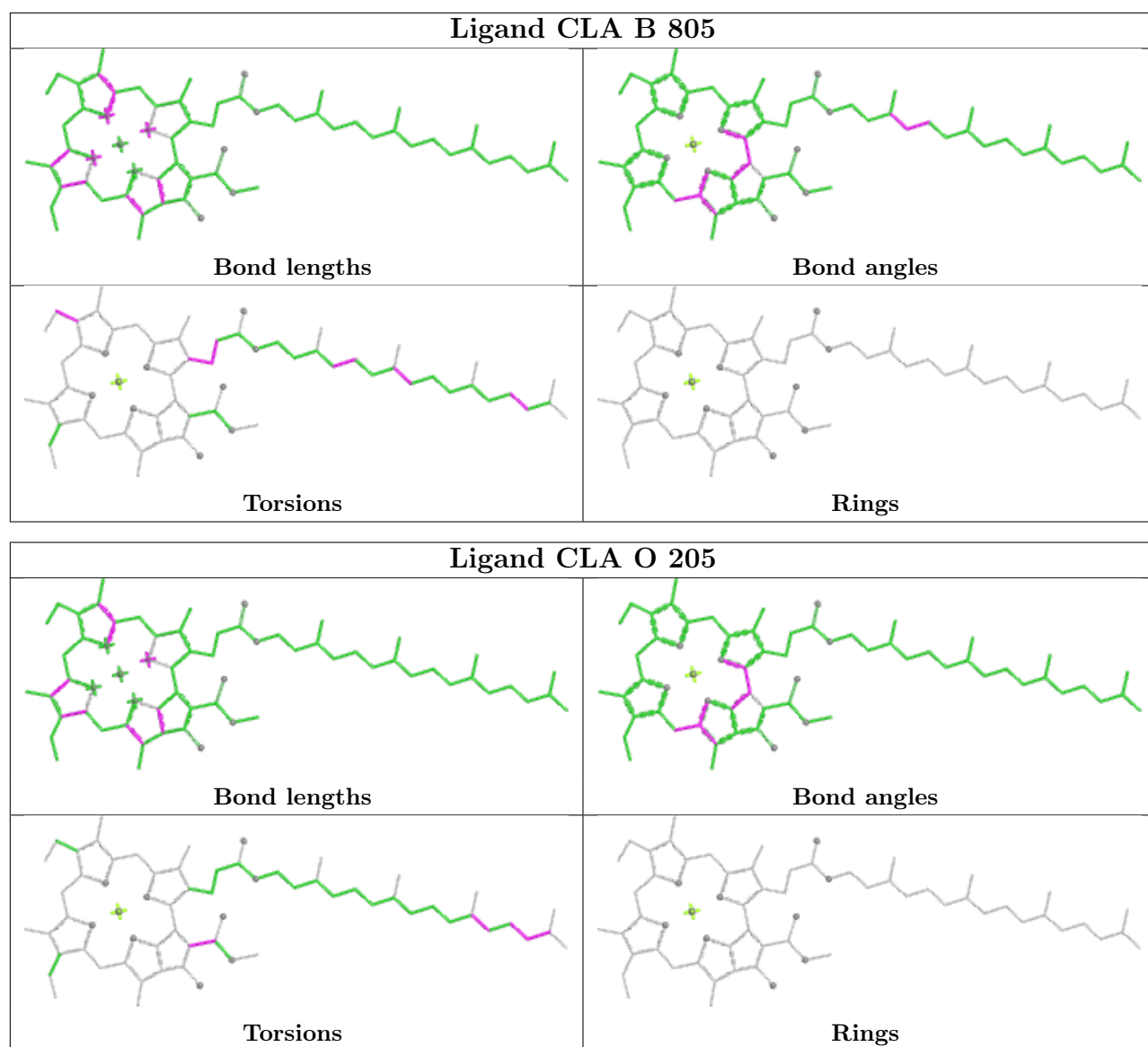


Torsions

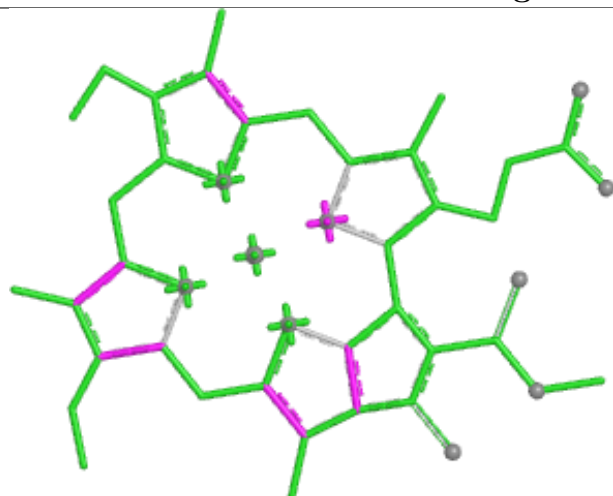


Rings

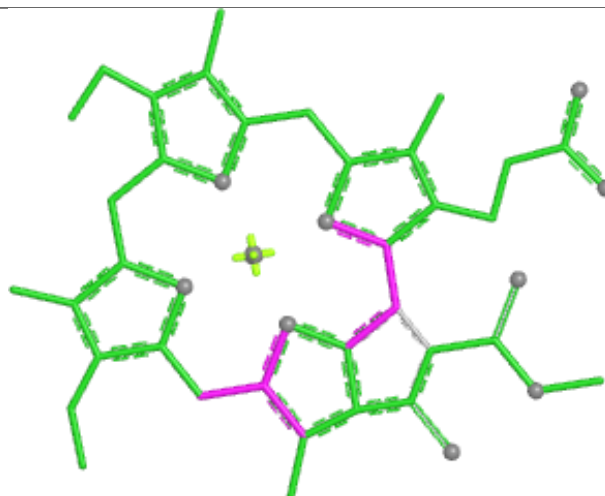




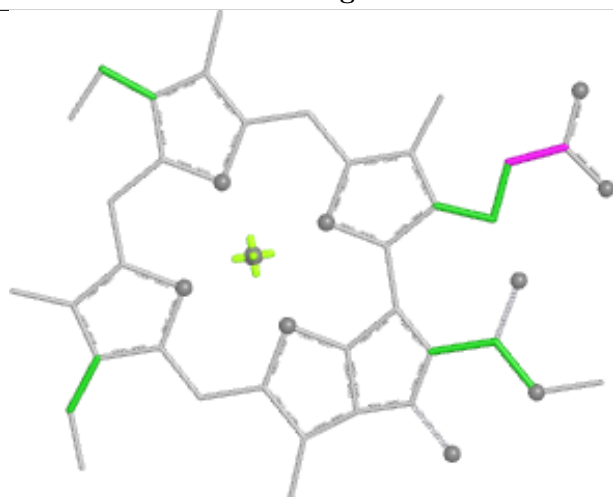
Ligand CLA U 206



Bond lengths



Bond angles

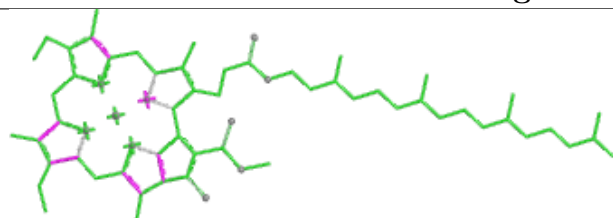


Torsions

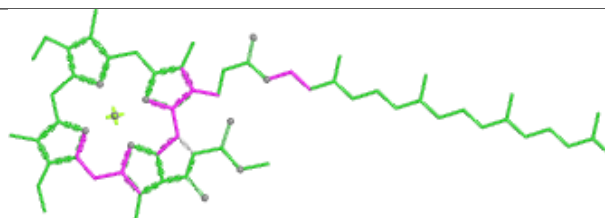


Rings

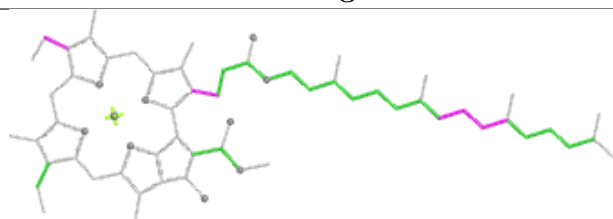
Ligand CLA U 207



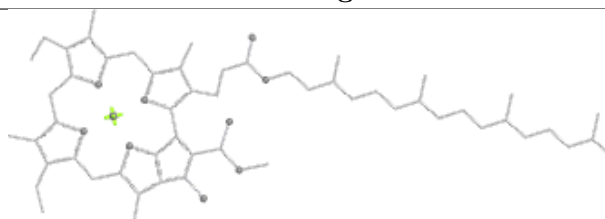
Bond lengths



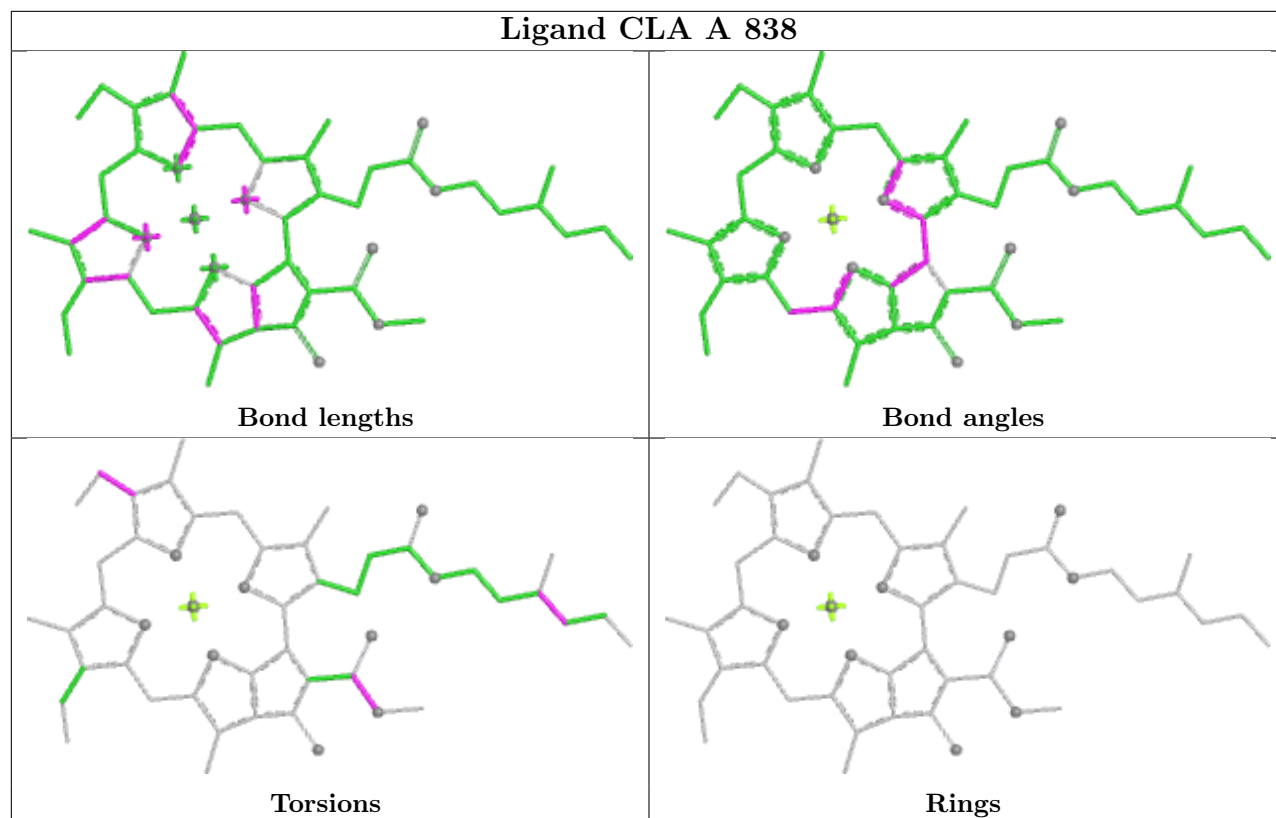
Bond angles



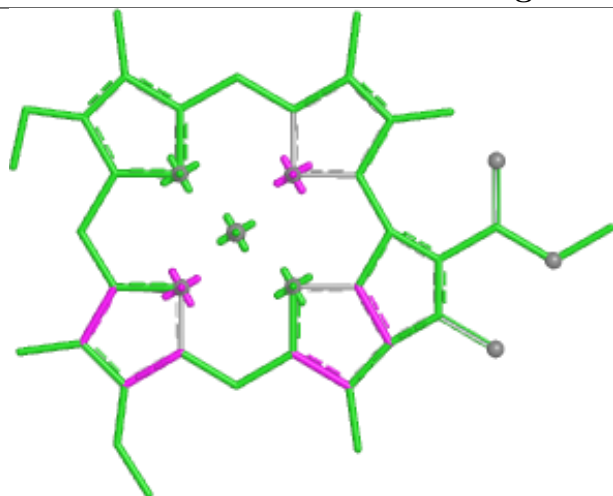
Torsions



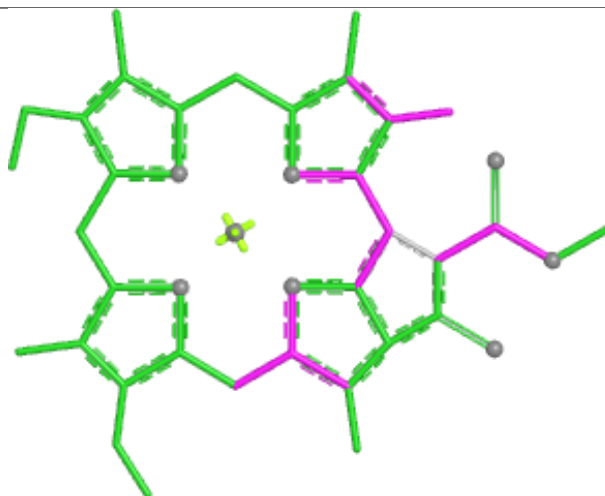
Rings



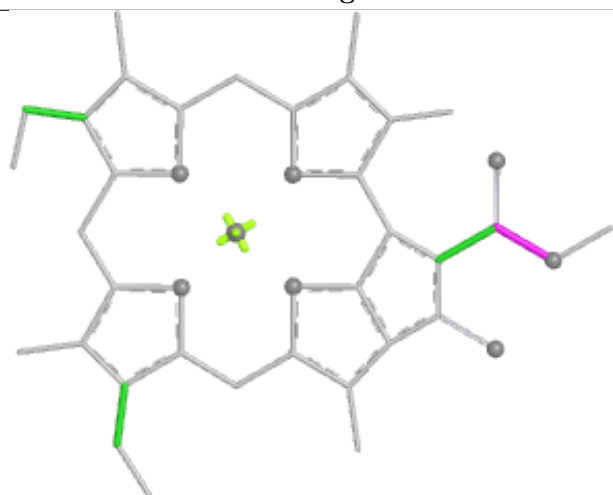
Ligand CLA P 312



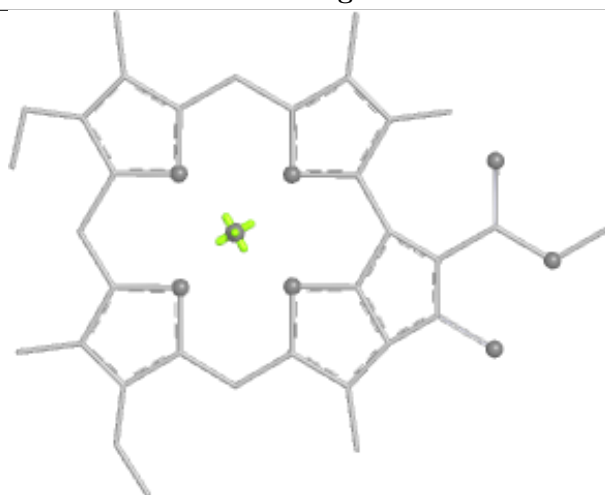
Bond lengths



Bond angles

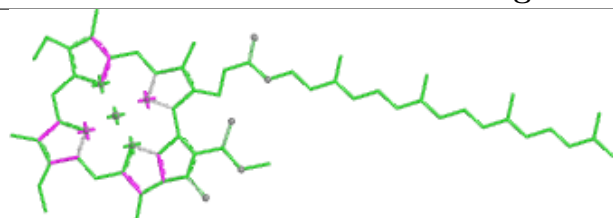


Torsions

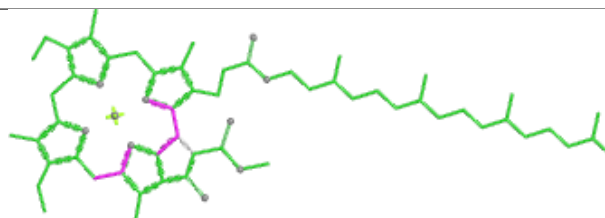


Rings

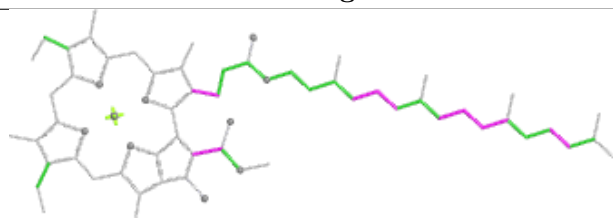
Ligand CLA A 851



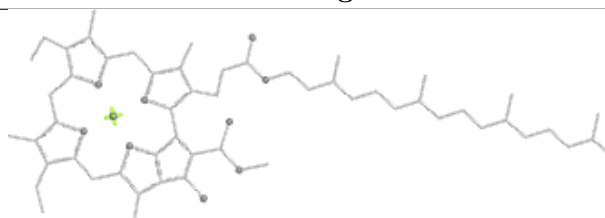
Bond lengths



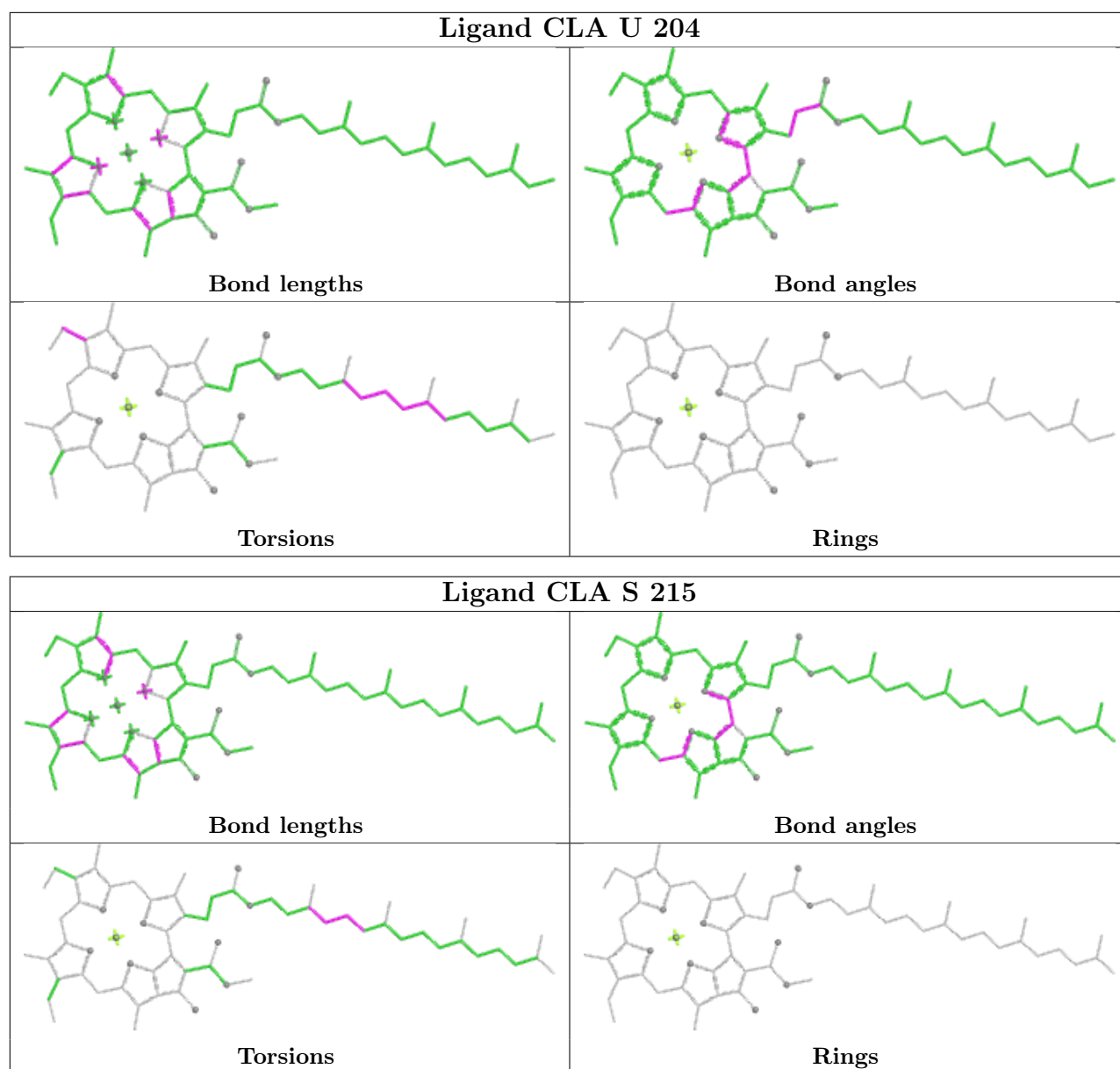
Bond angles

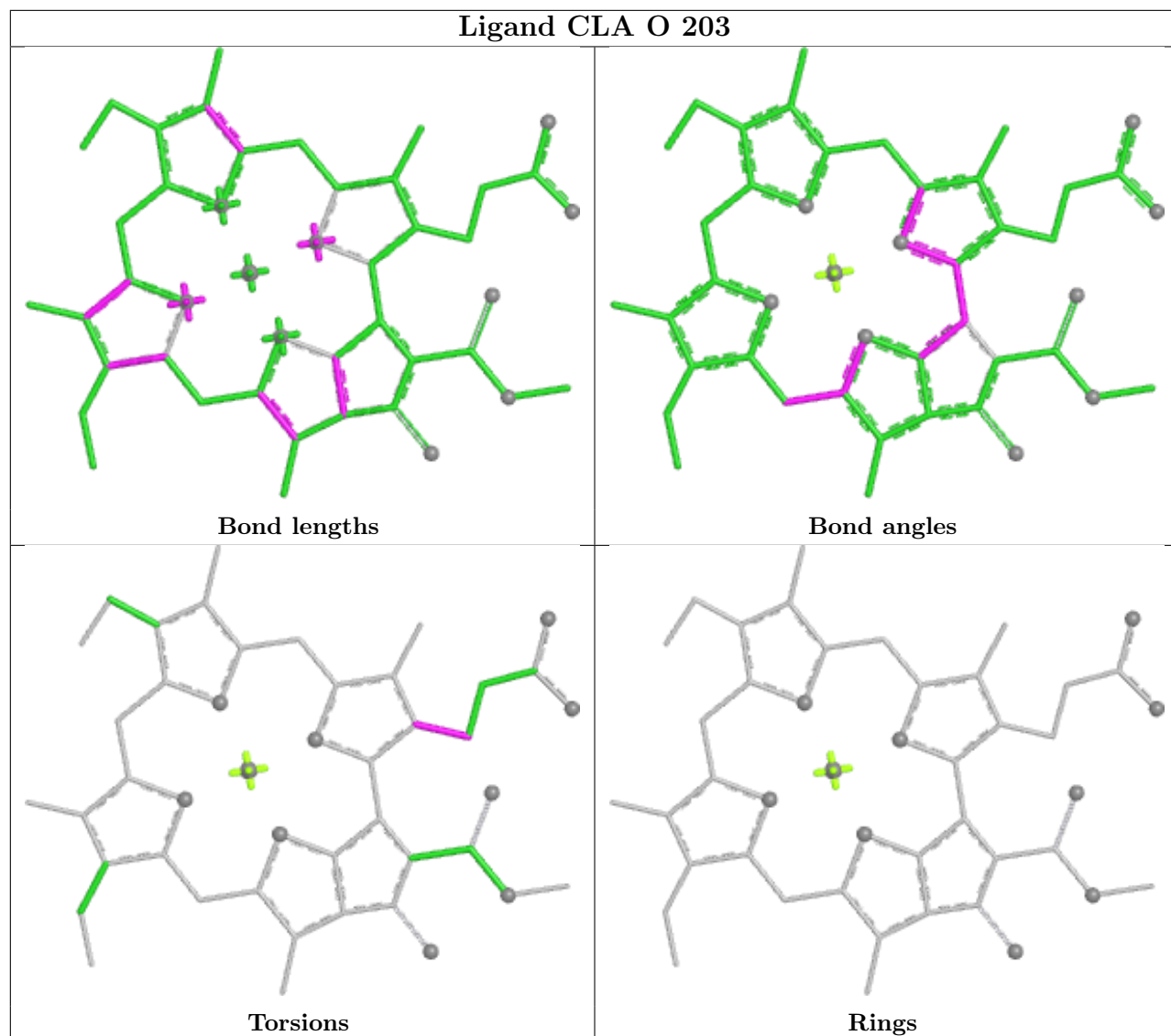


Torsions

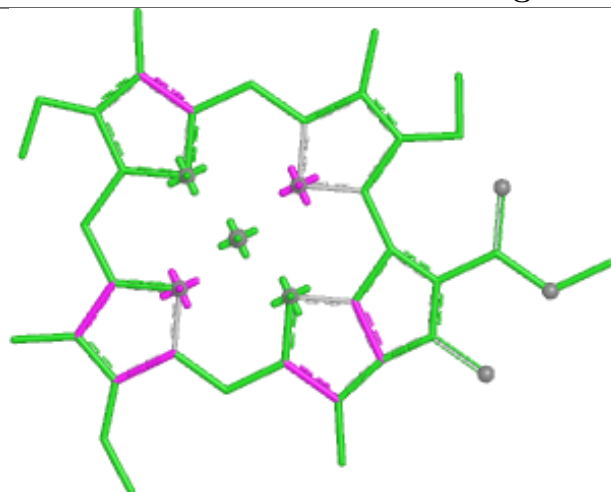


Rings

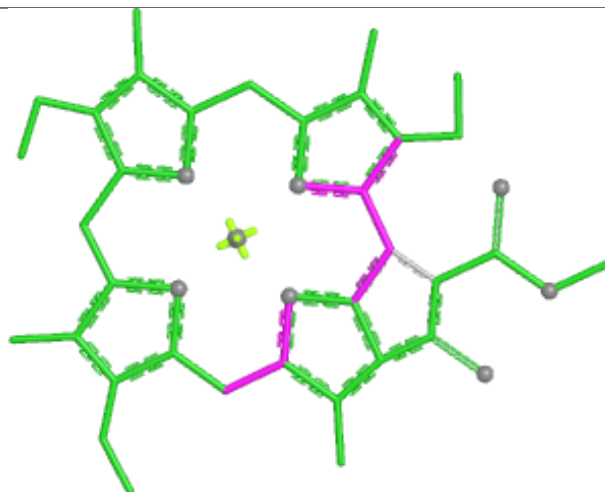




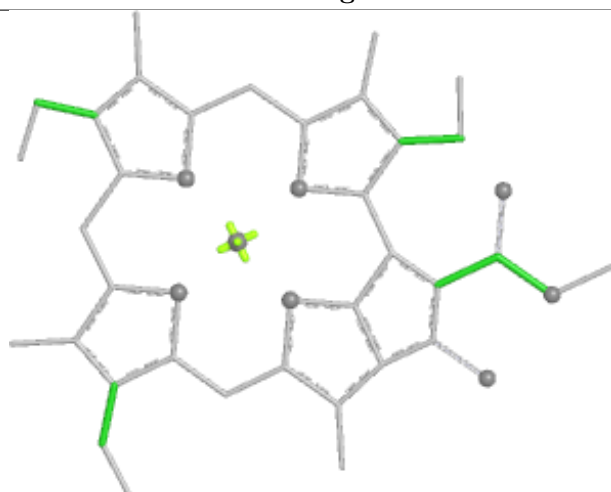
Ligand CLA U 209



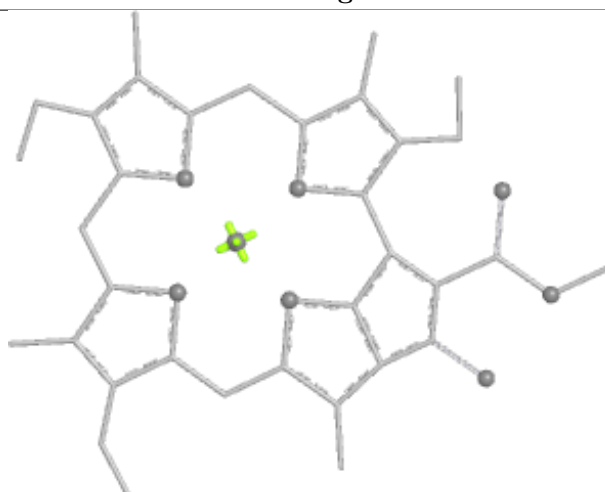
Bond lengths



Bond angles

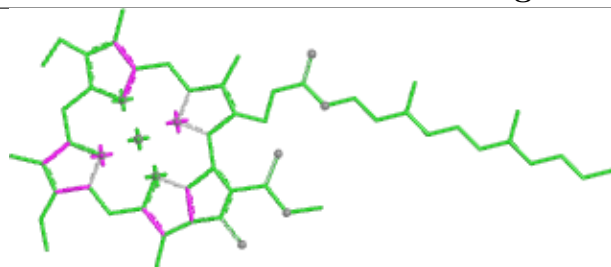


Torsions

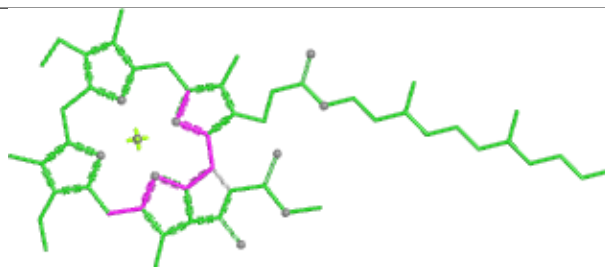


Rings

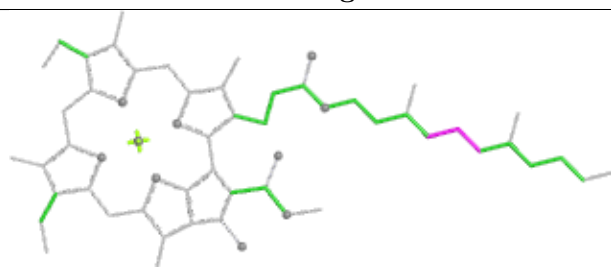
Ligand CLA B 828



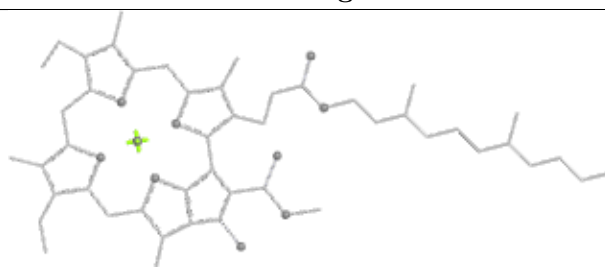
Bond lengths



Bond angles

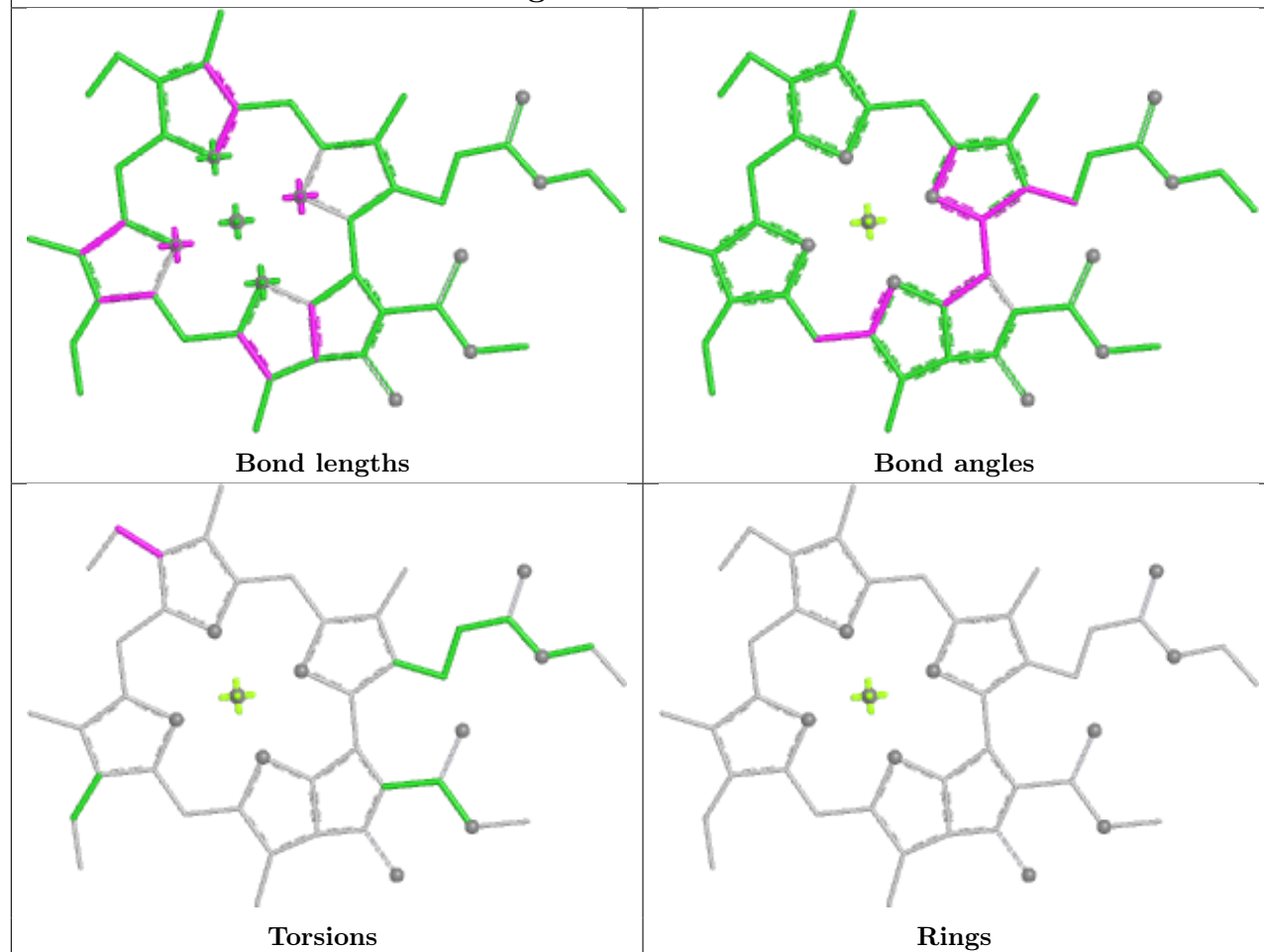


Torsions

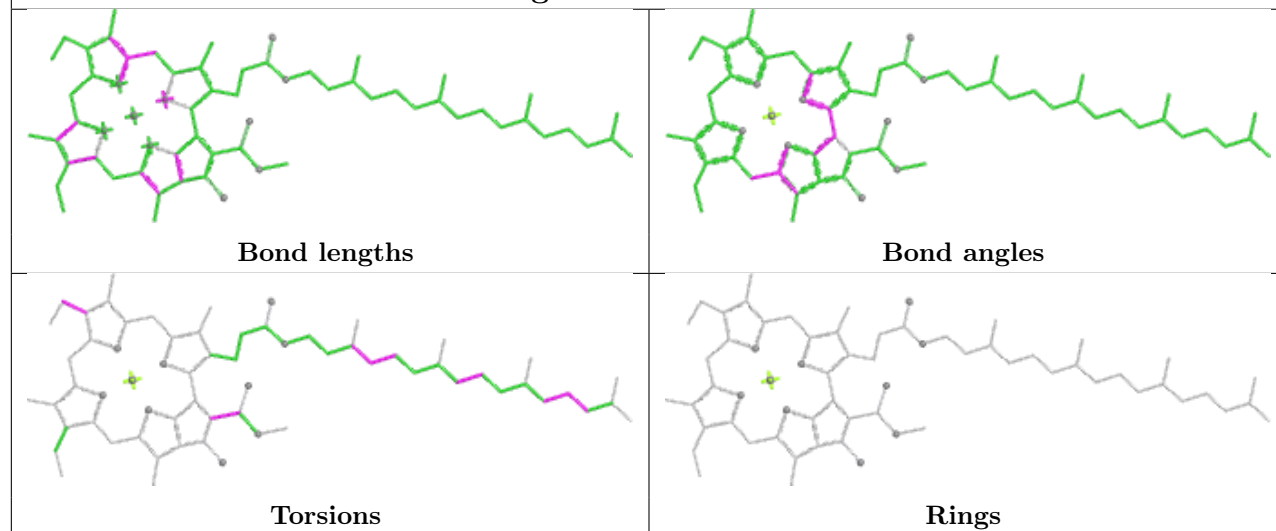


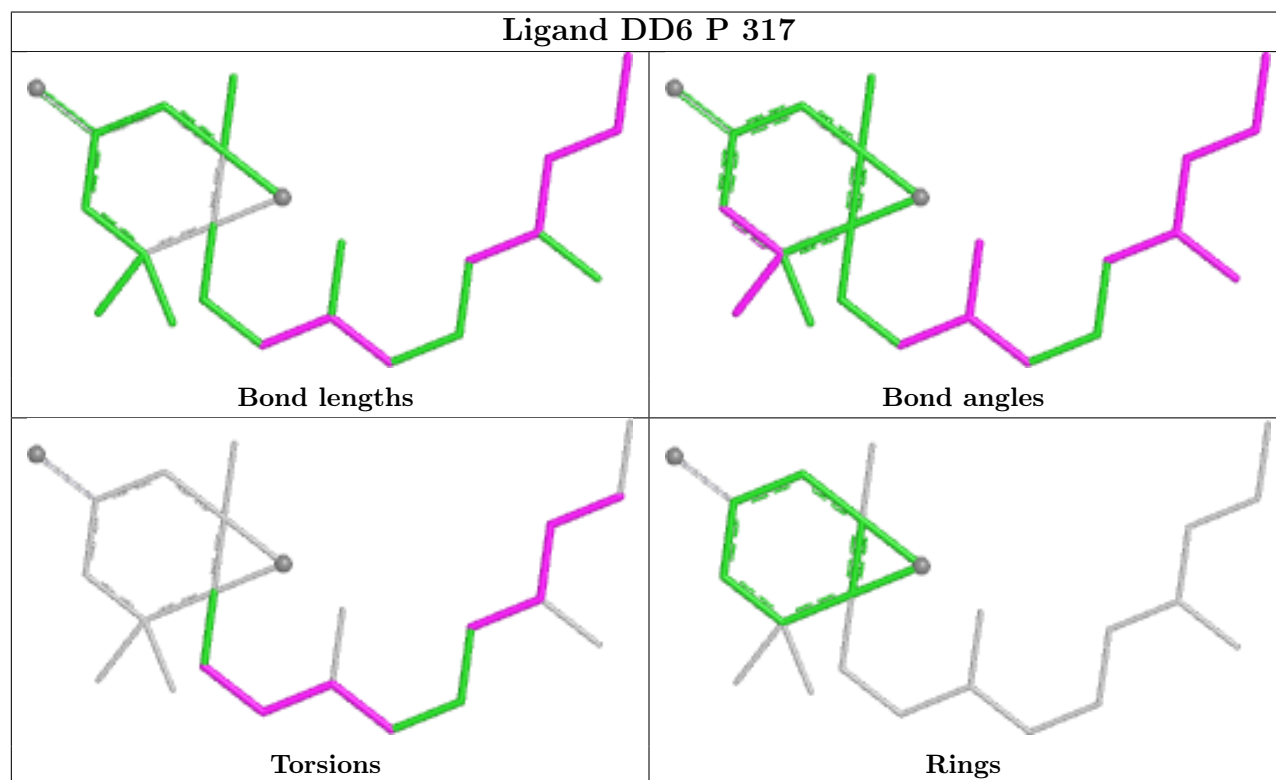
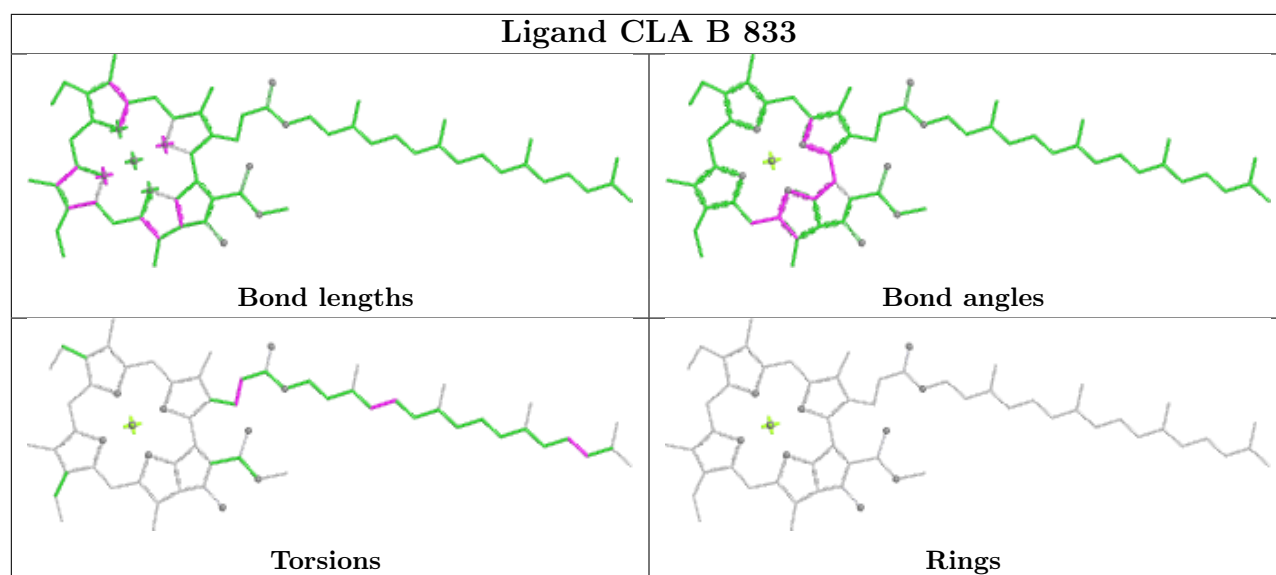
Rings

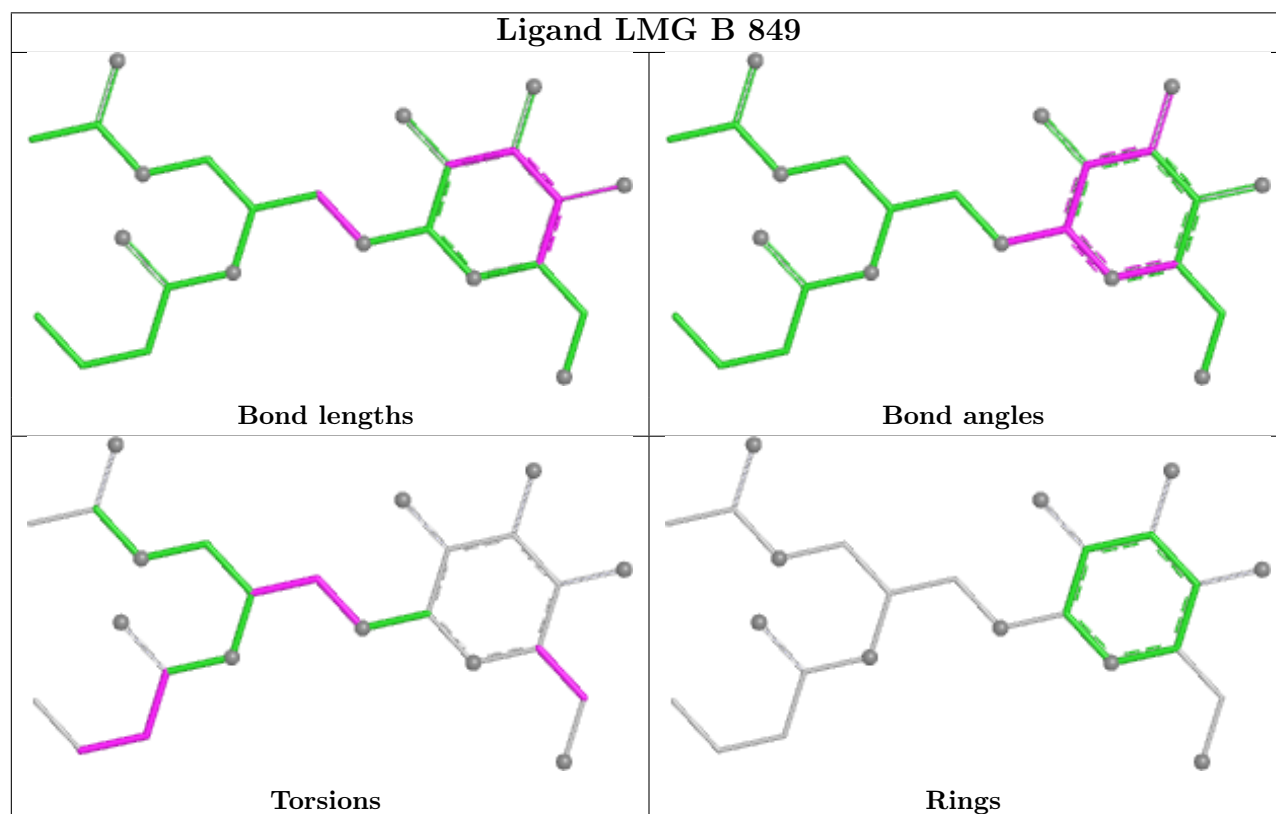
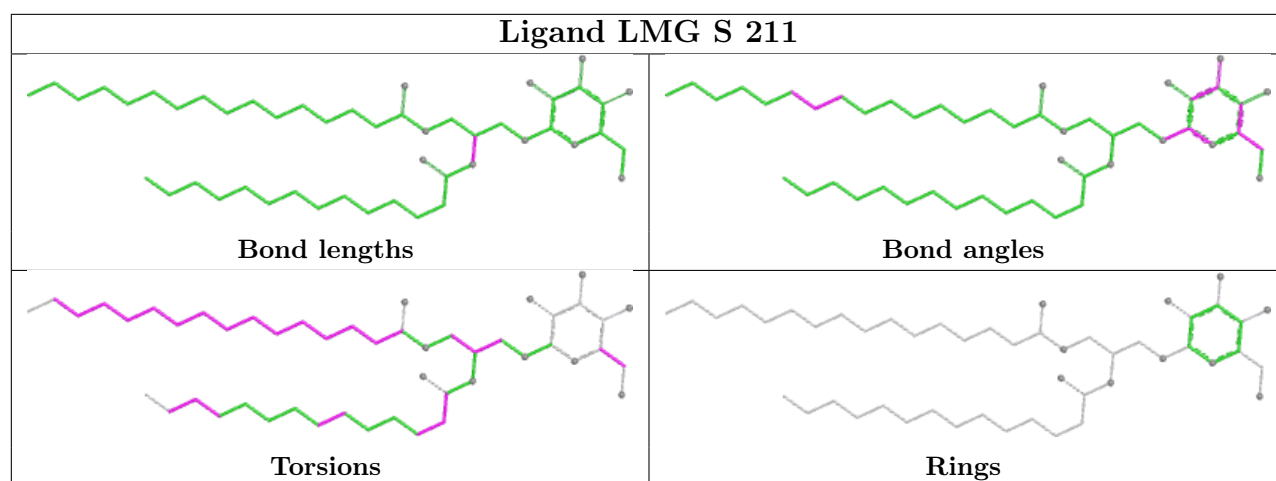
Ligand CLA P 315

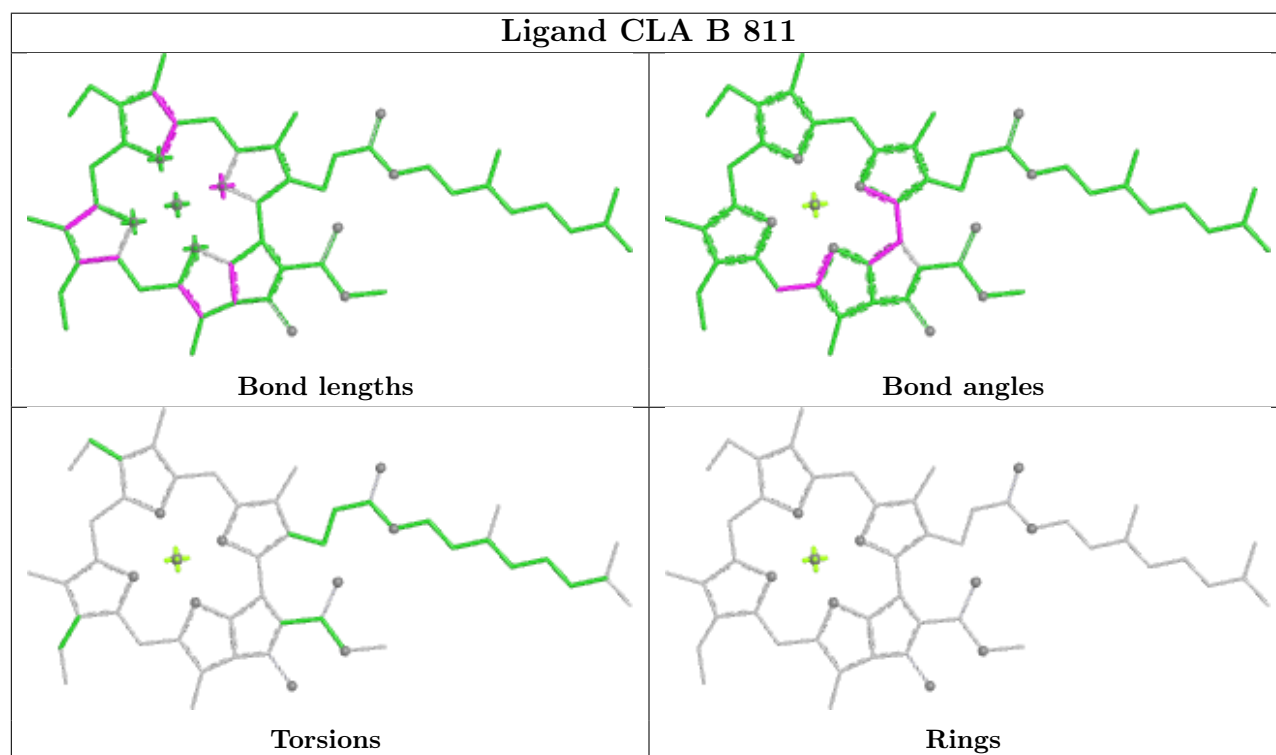
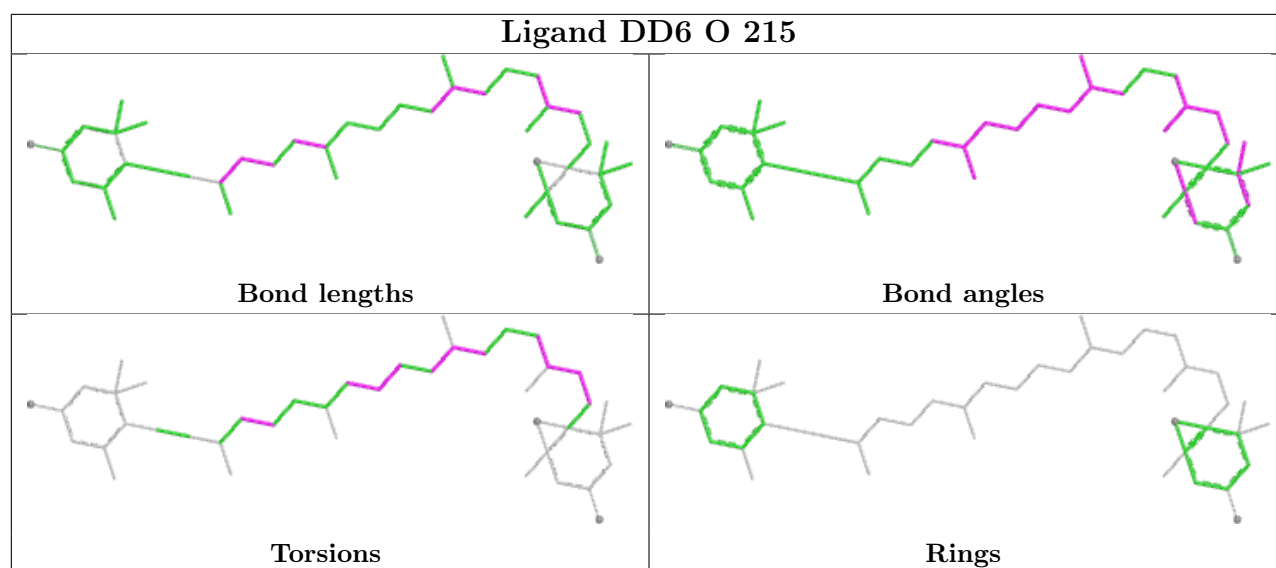


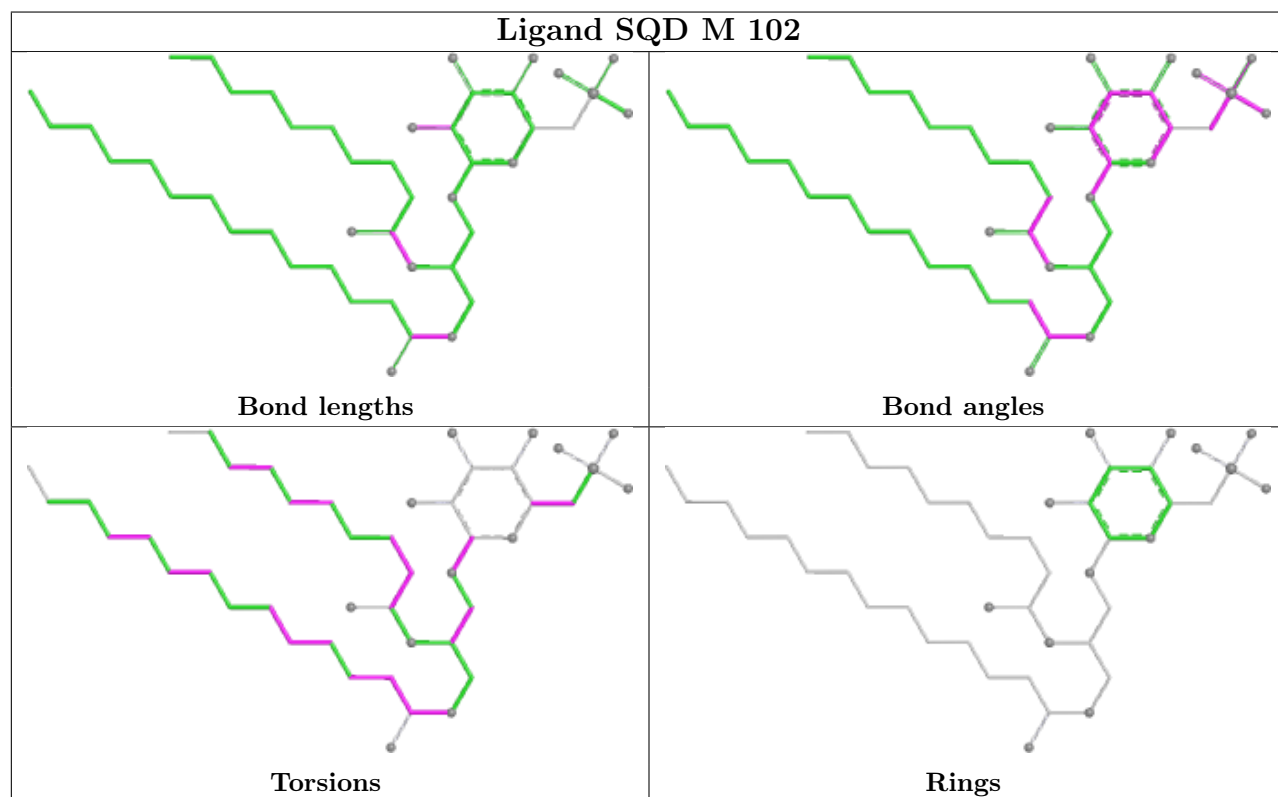
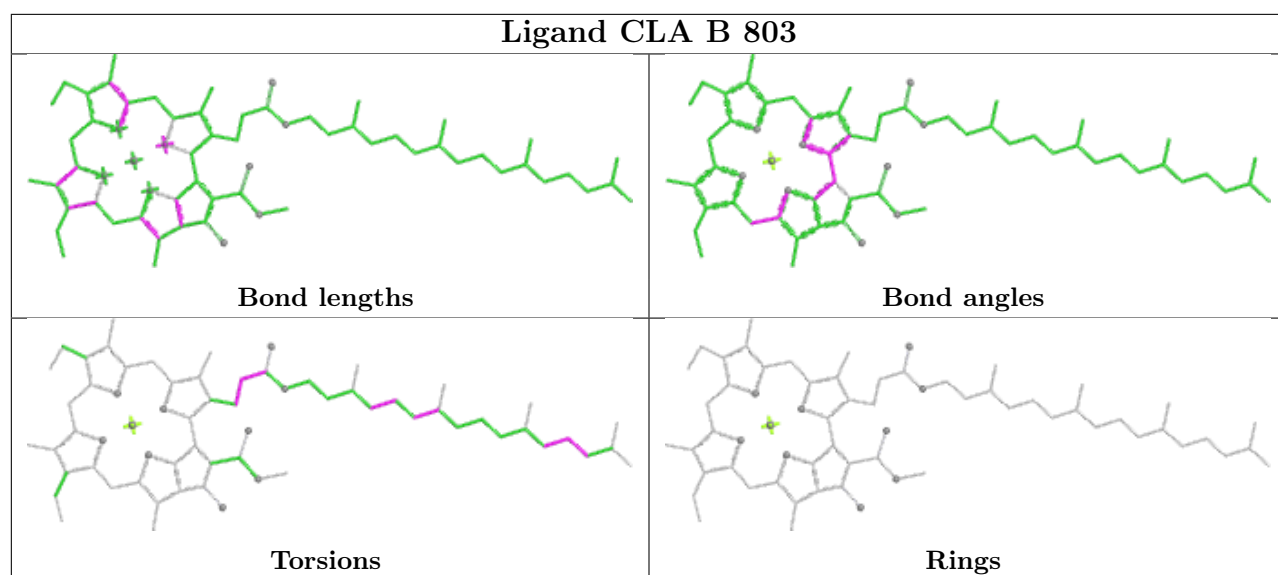
Ligand CLA U 210

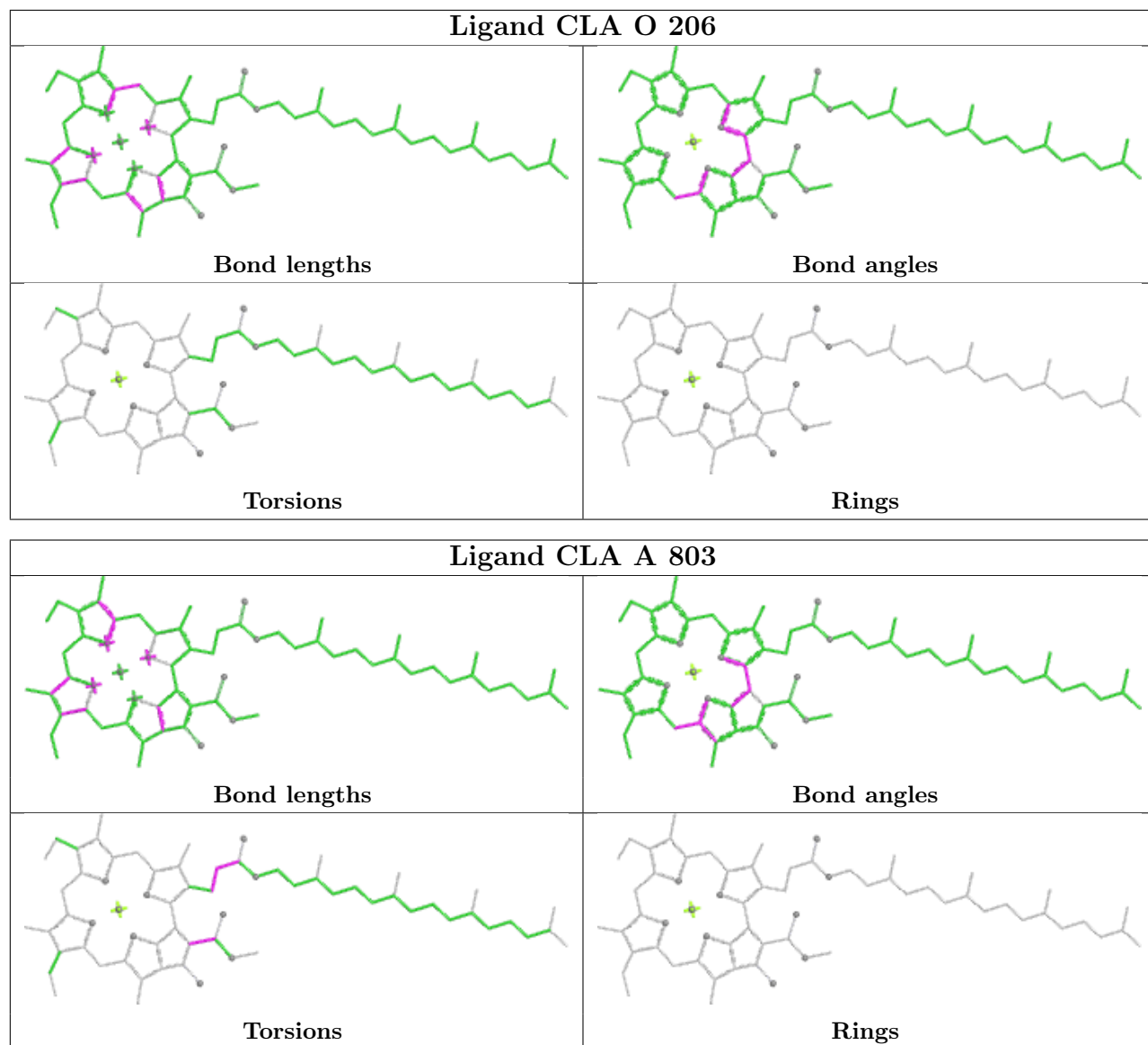




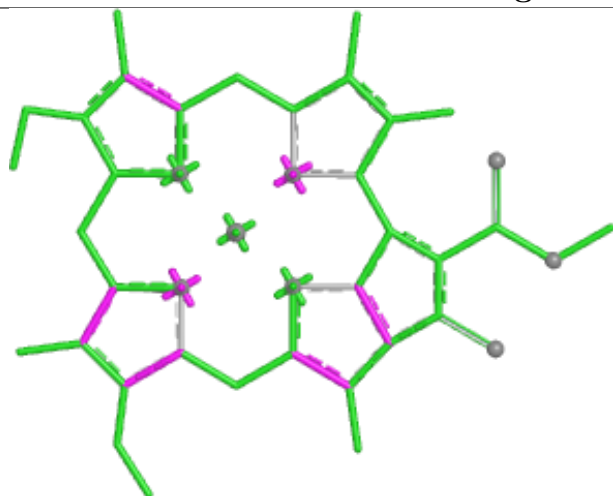




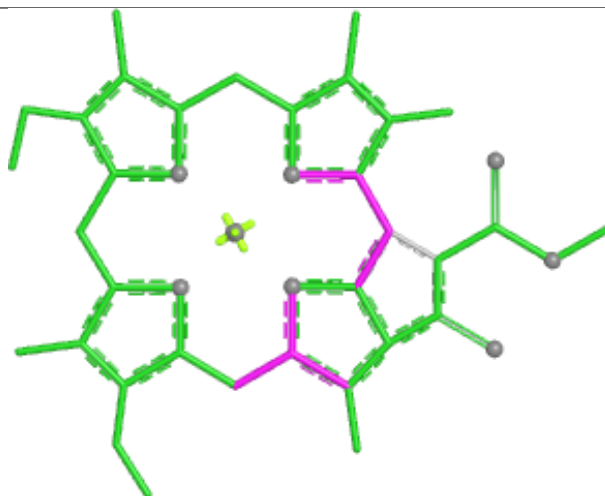




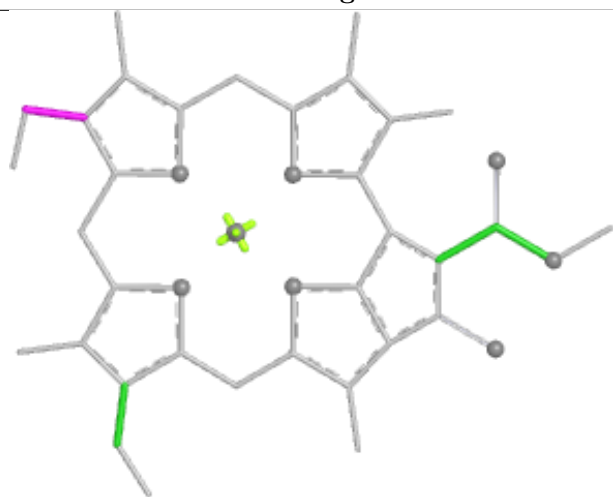
Ligand CLA T 203



Bond lengths



Bond angles

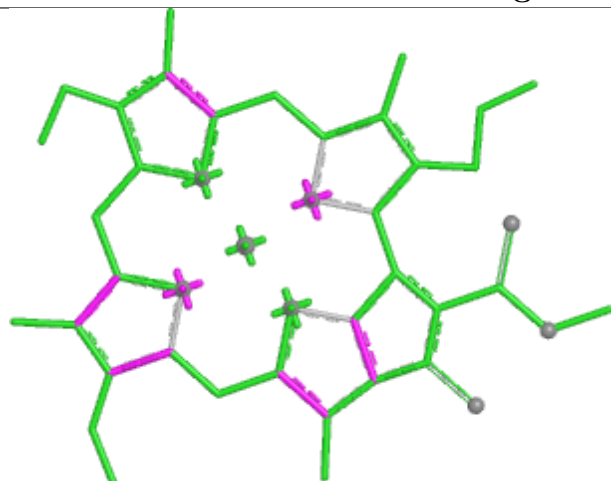


Torsions

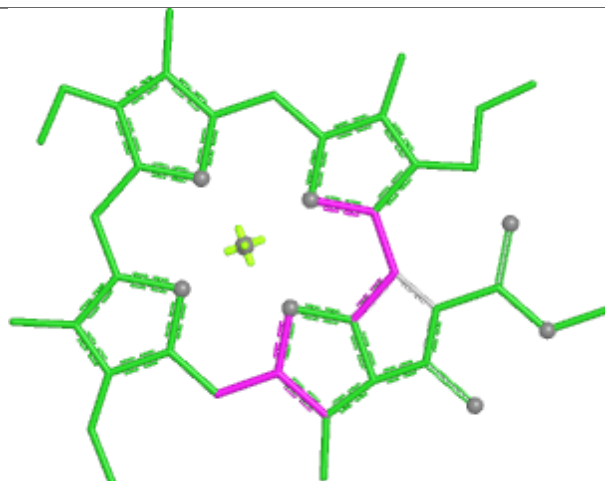


Rings

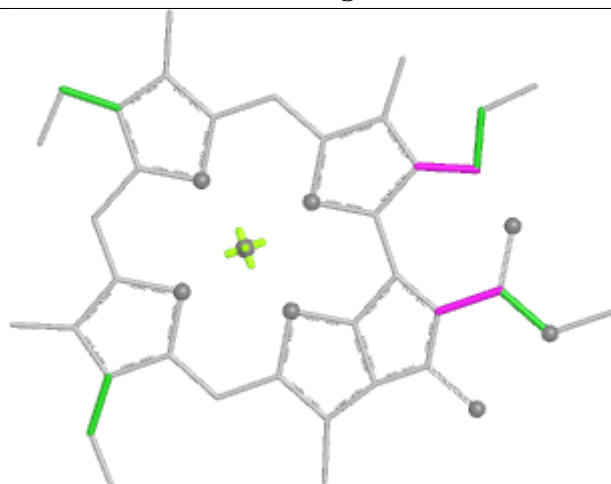
Ligand CLA O 202



Bond lengths



Bond angles

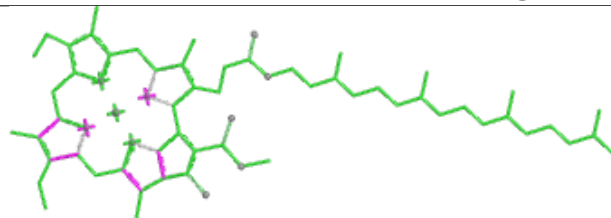


Torsions

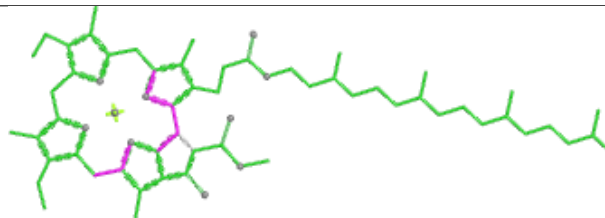


Rings

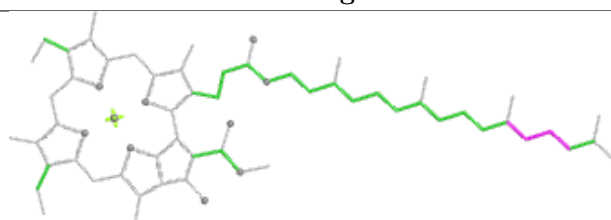
Ligand CLA A 828



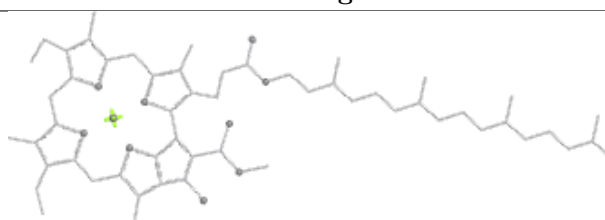
Bond lengths



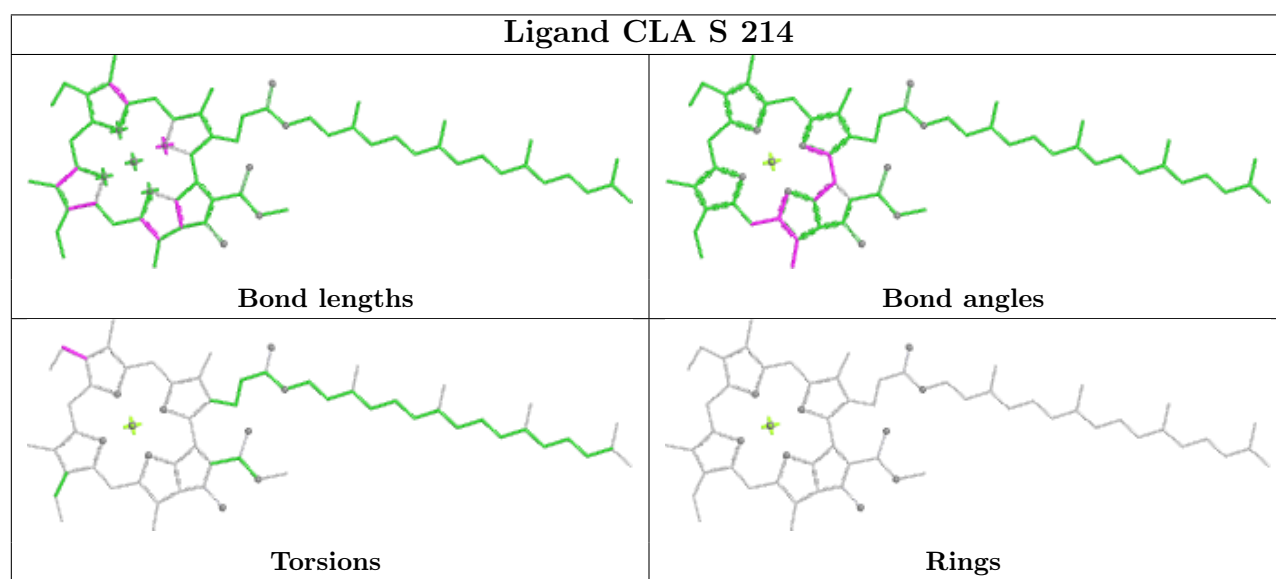
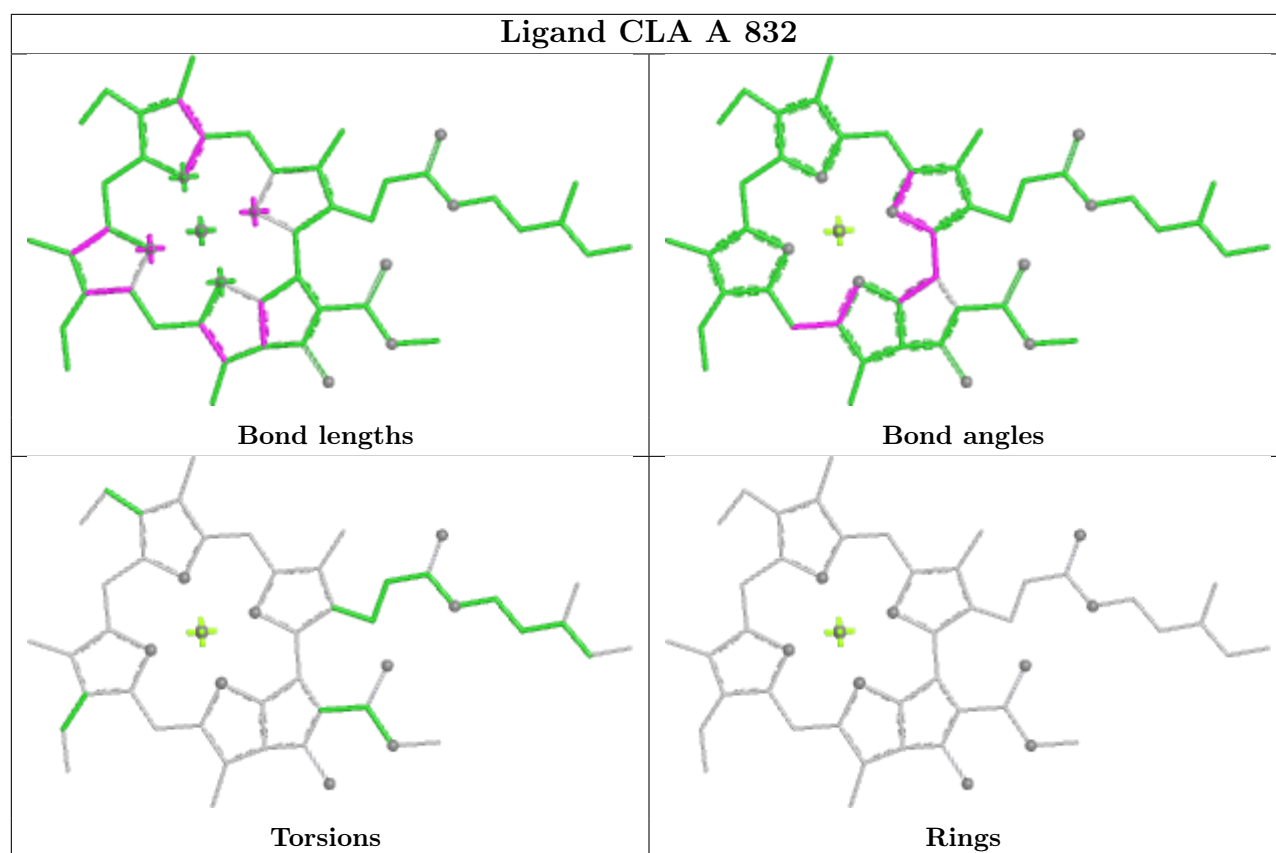
Bond angles

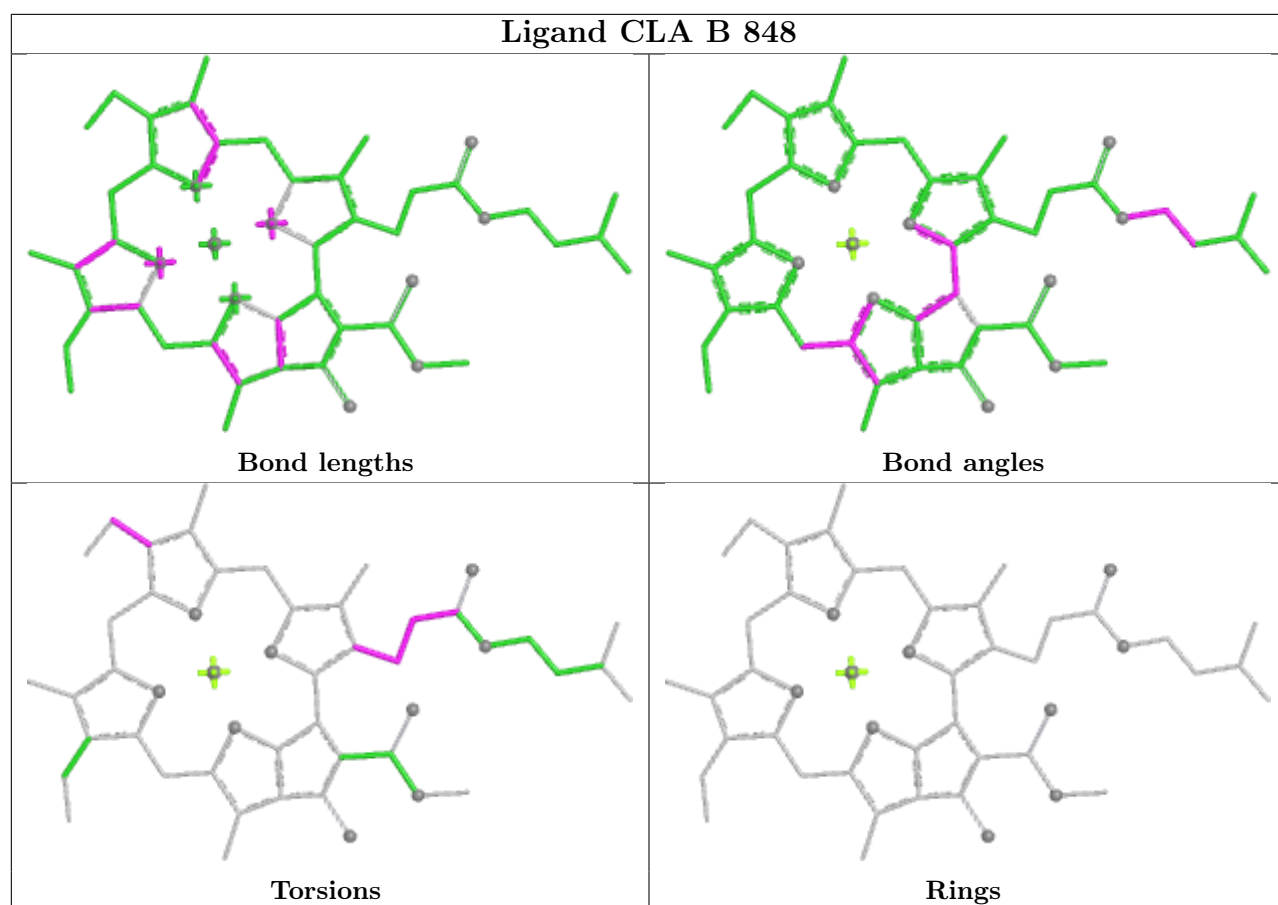


Torsions

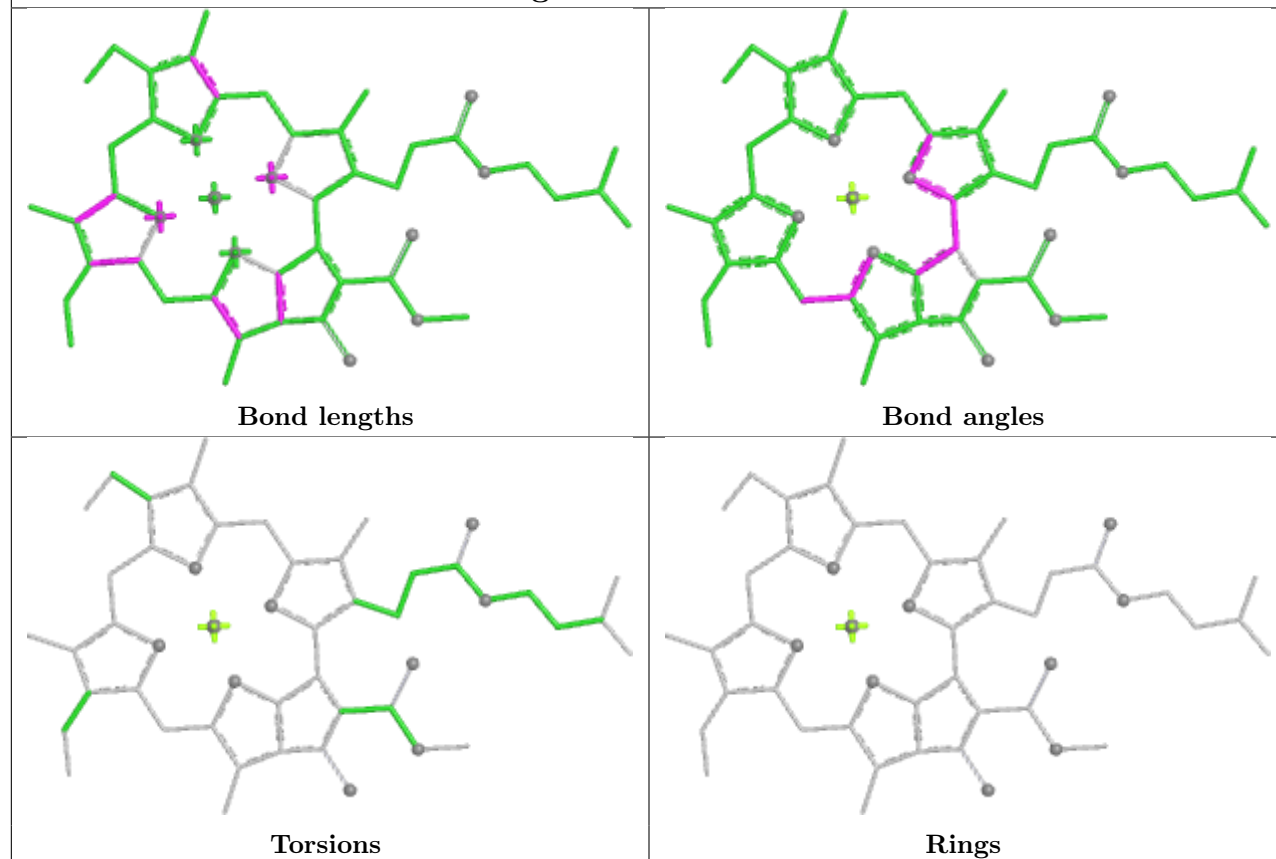


Rings

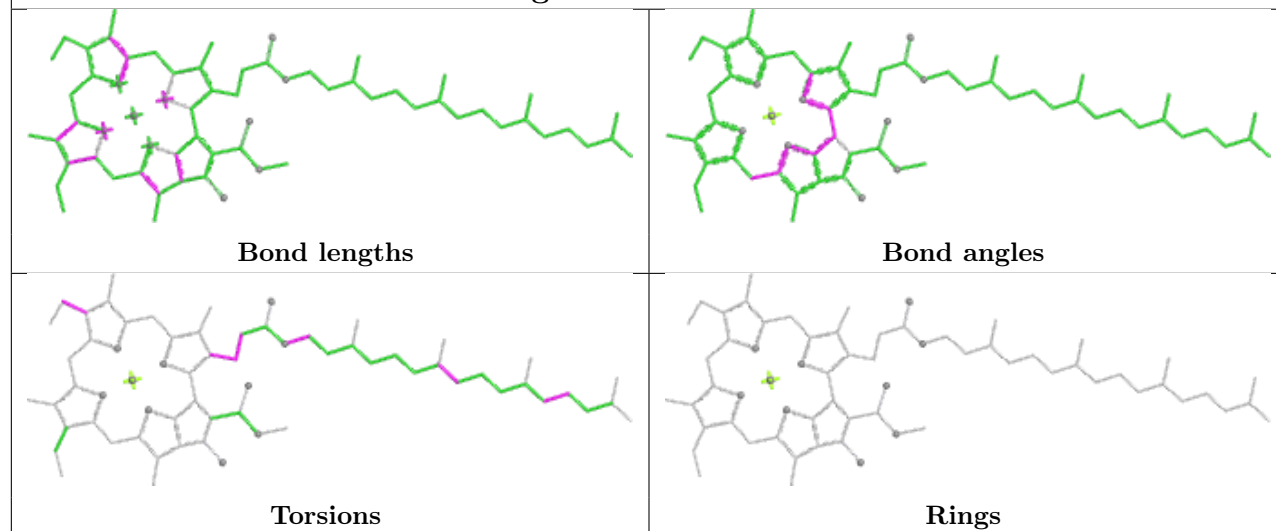




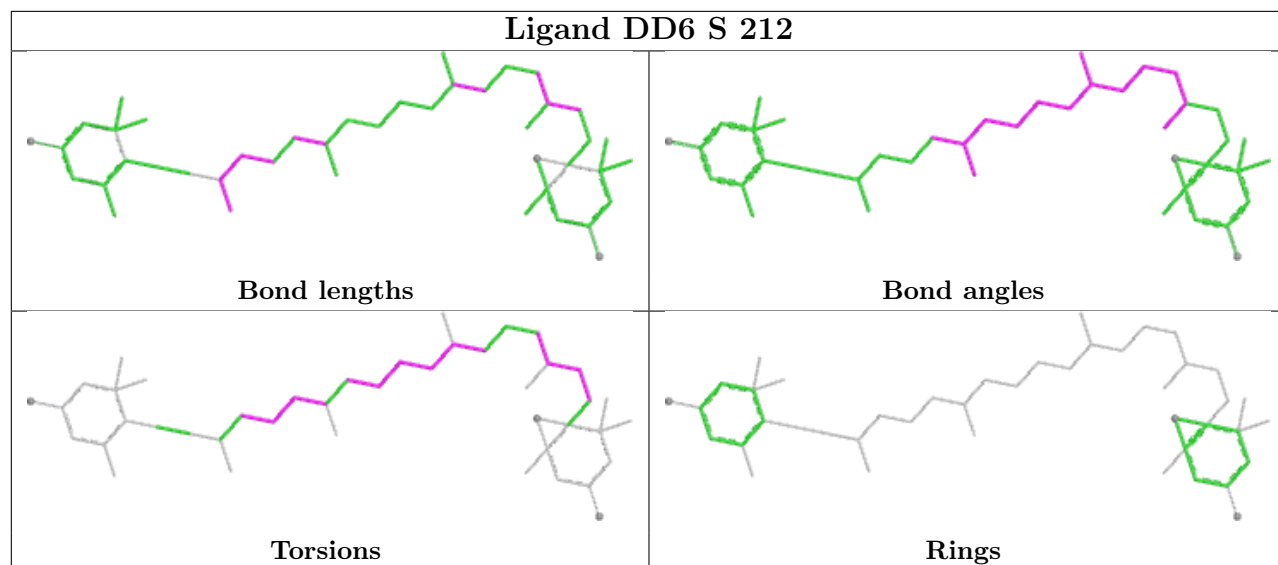
Ligand CLA A 827



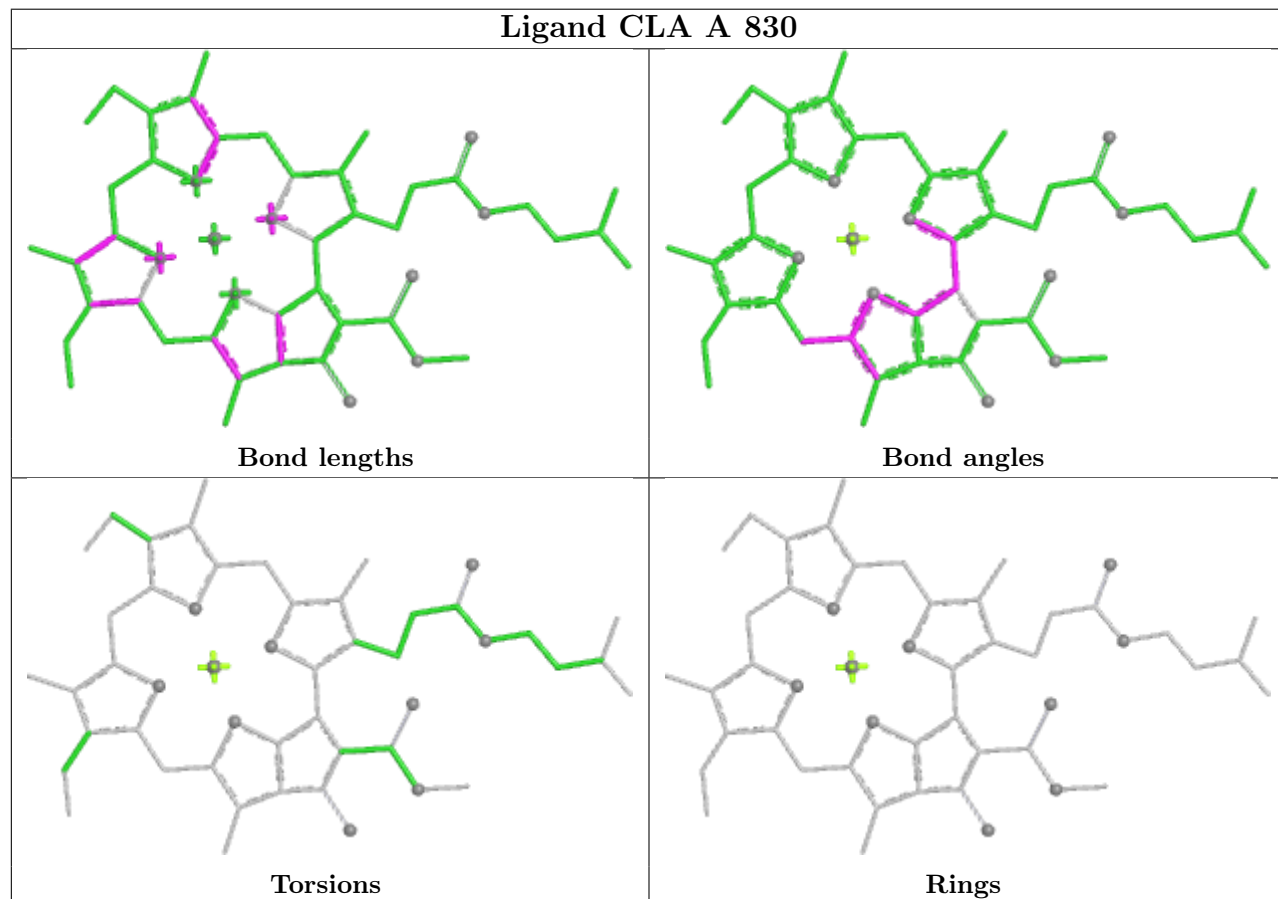
Ligand CLA A 807

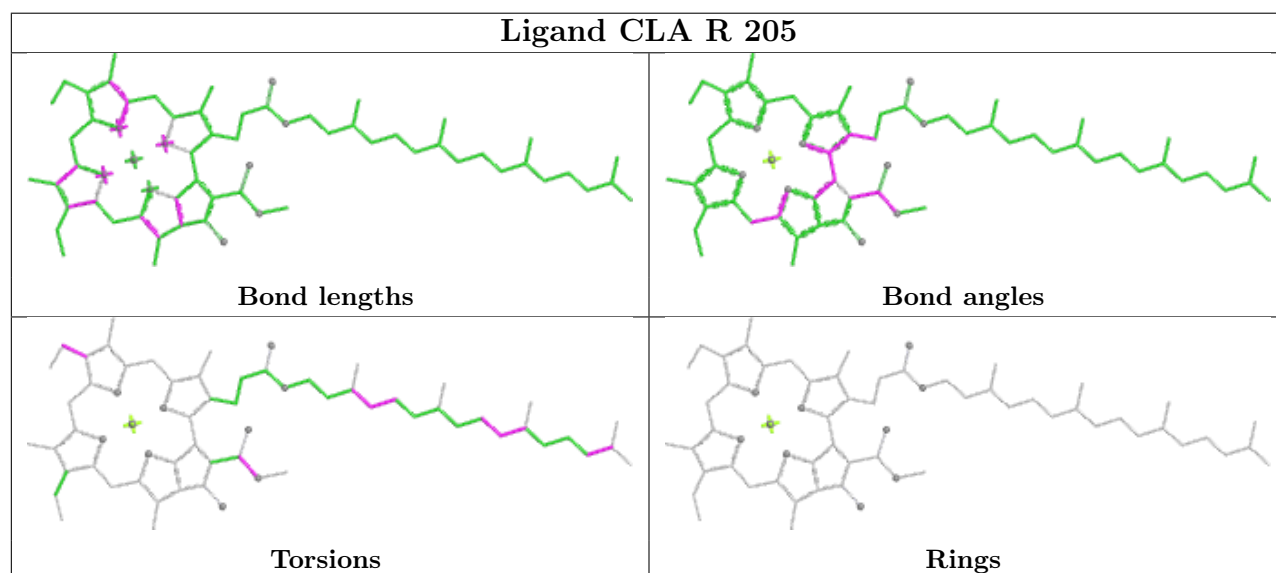
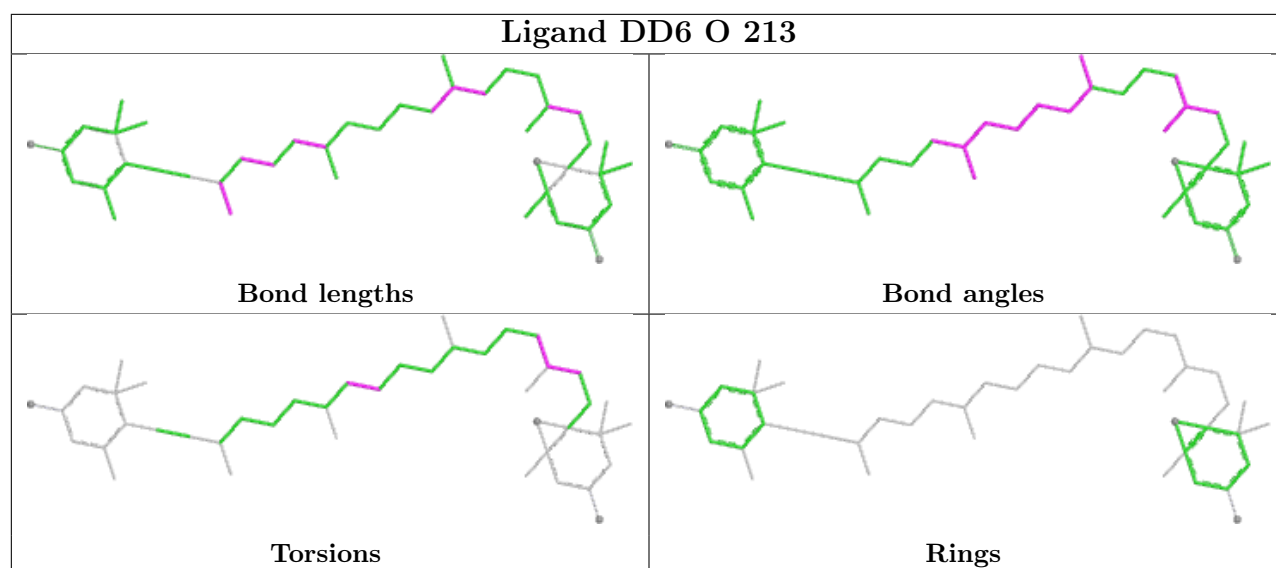


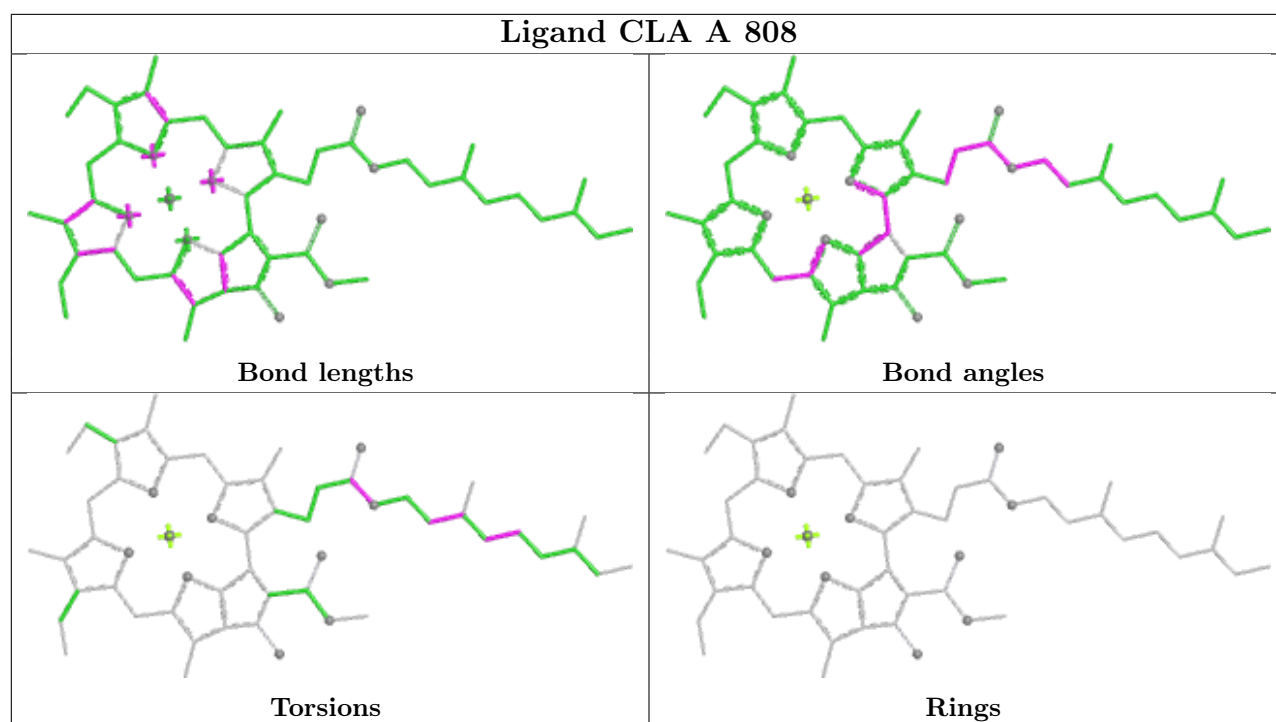
Ligand DD6 S 212



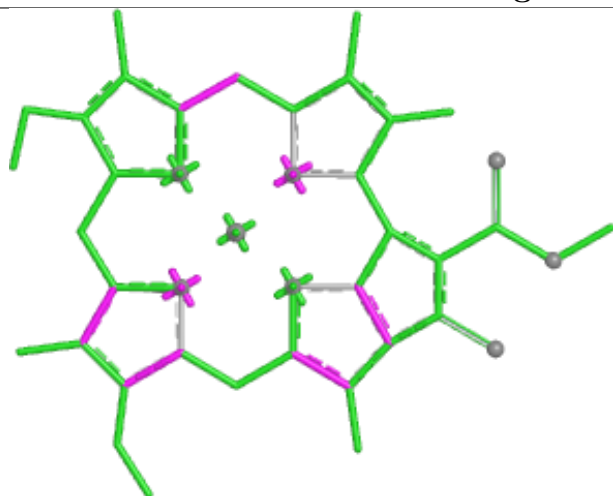
Ligand CLA A 830



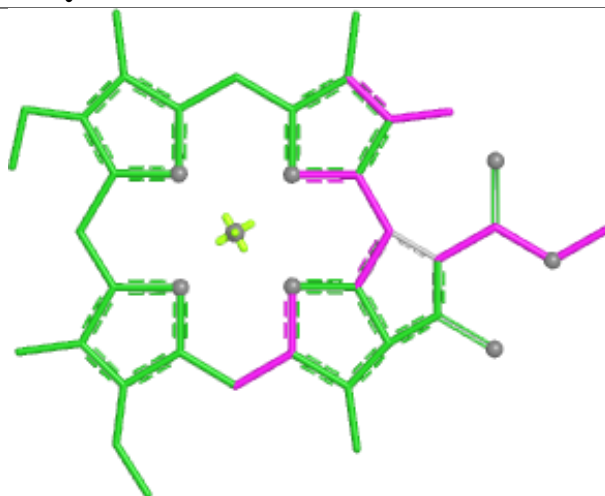




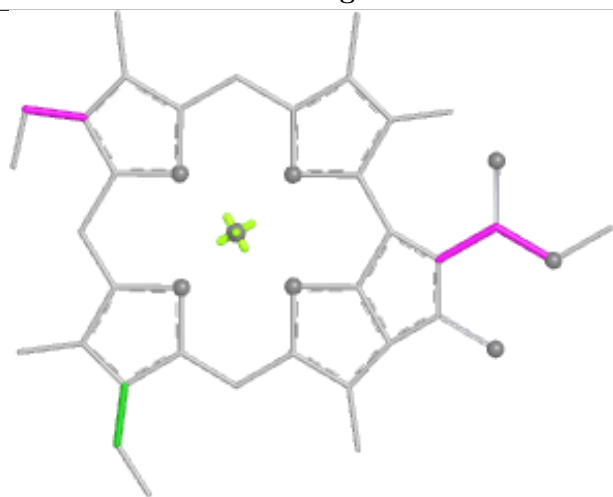
Ligand CLA Q 211



Bond lengths



Bond angles

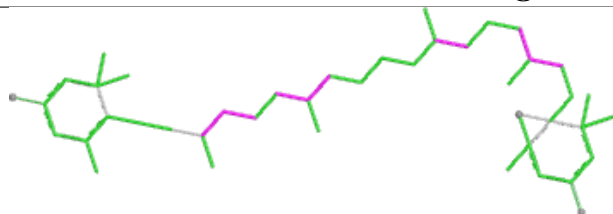


Torsions

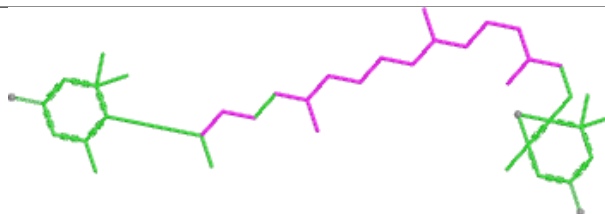


Rings

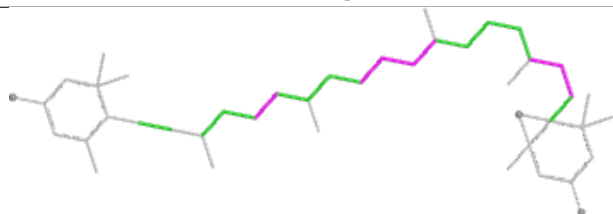
Ligand DD6 O 214



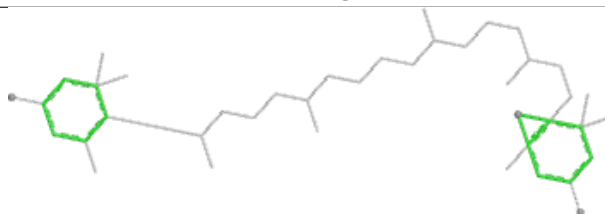
Bond lengths



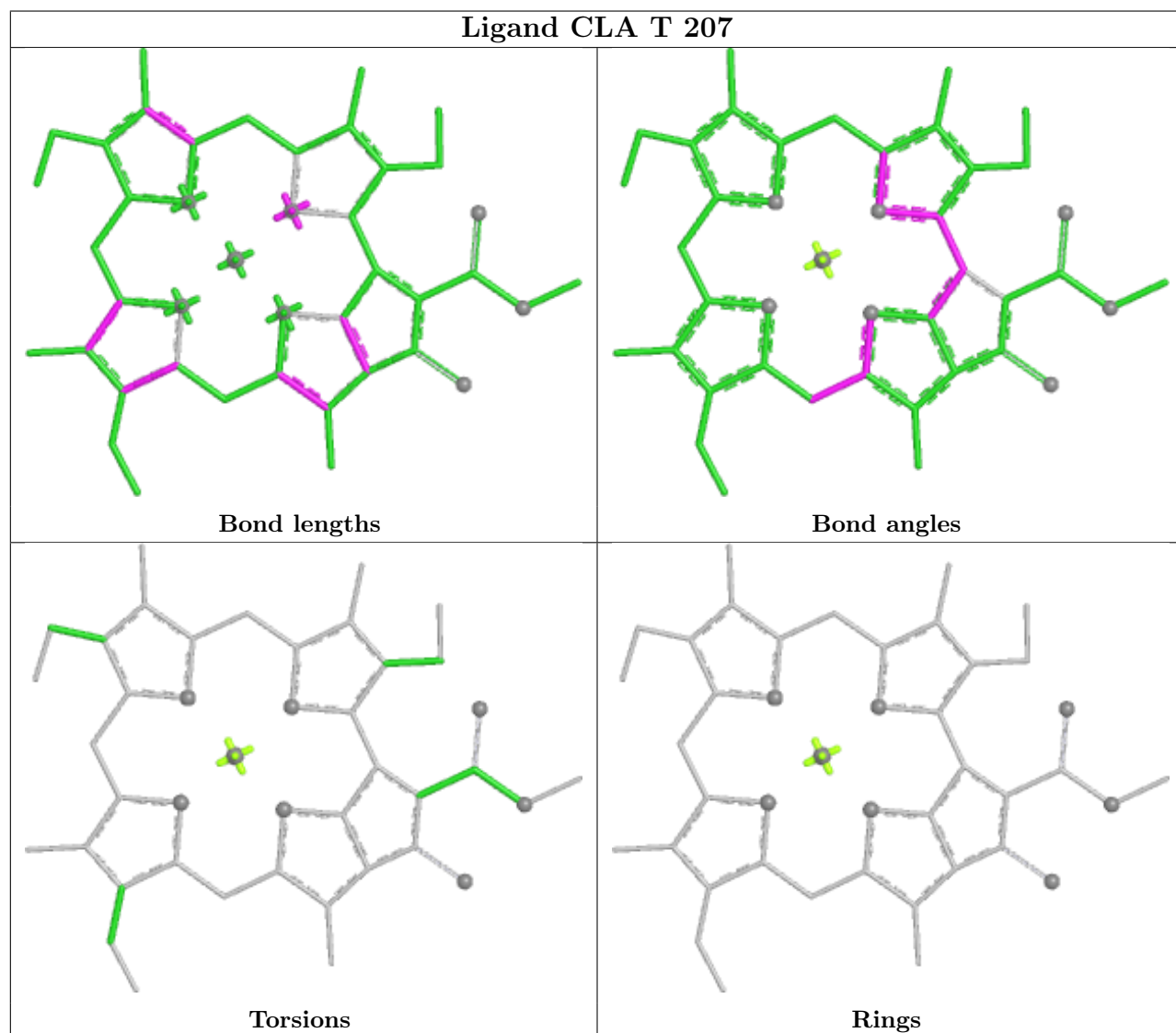
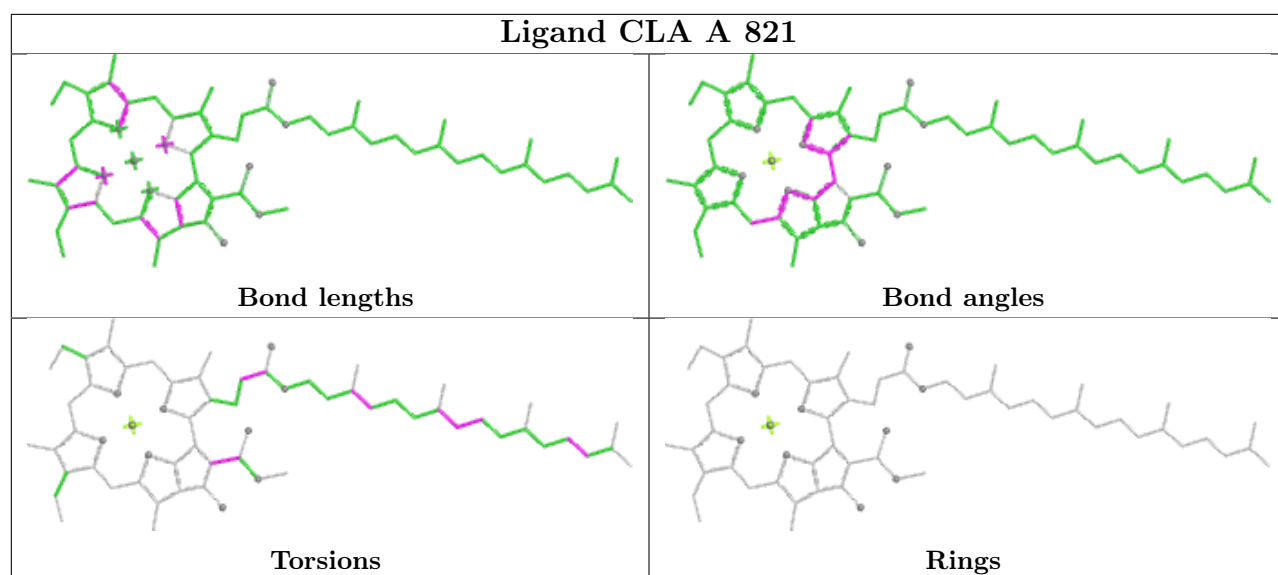
Bond angles



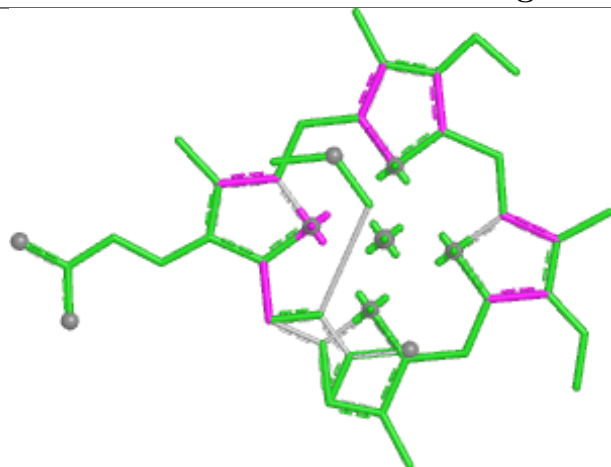
Torsions



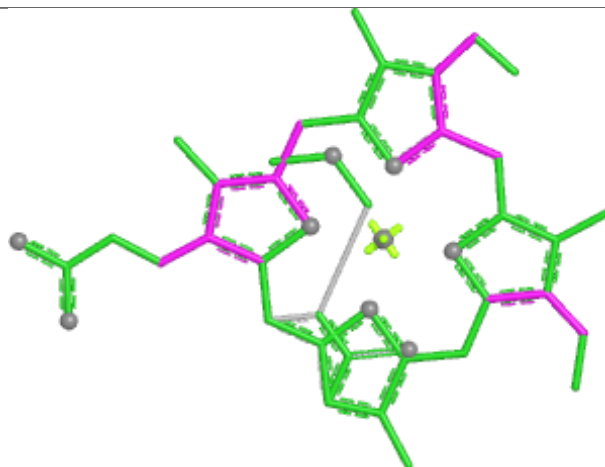
Rings



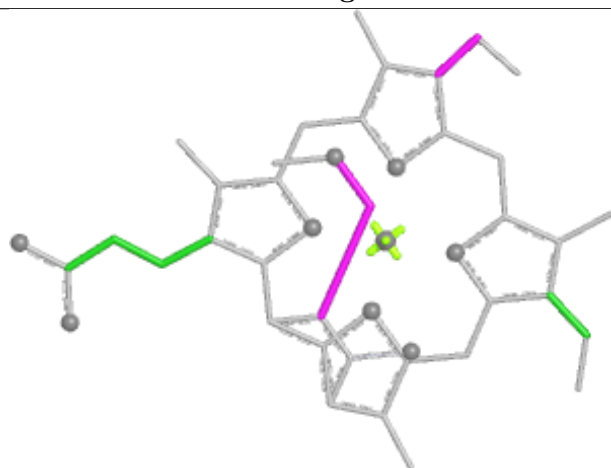
Ligand KC1 P 311



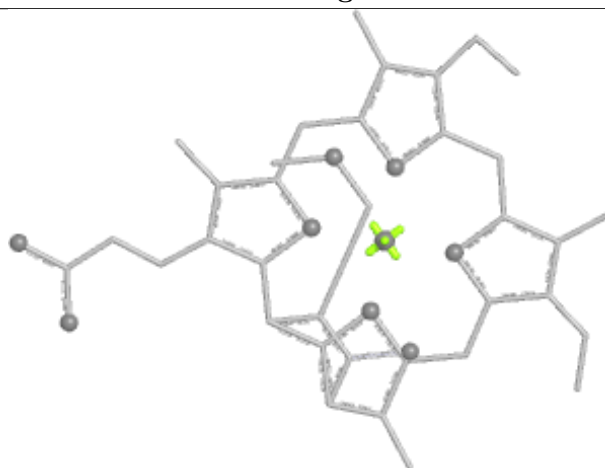
Bond lengths



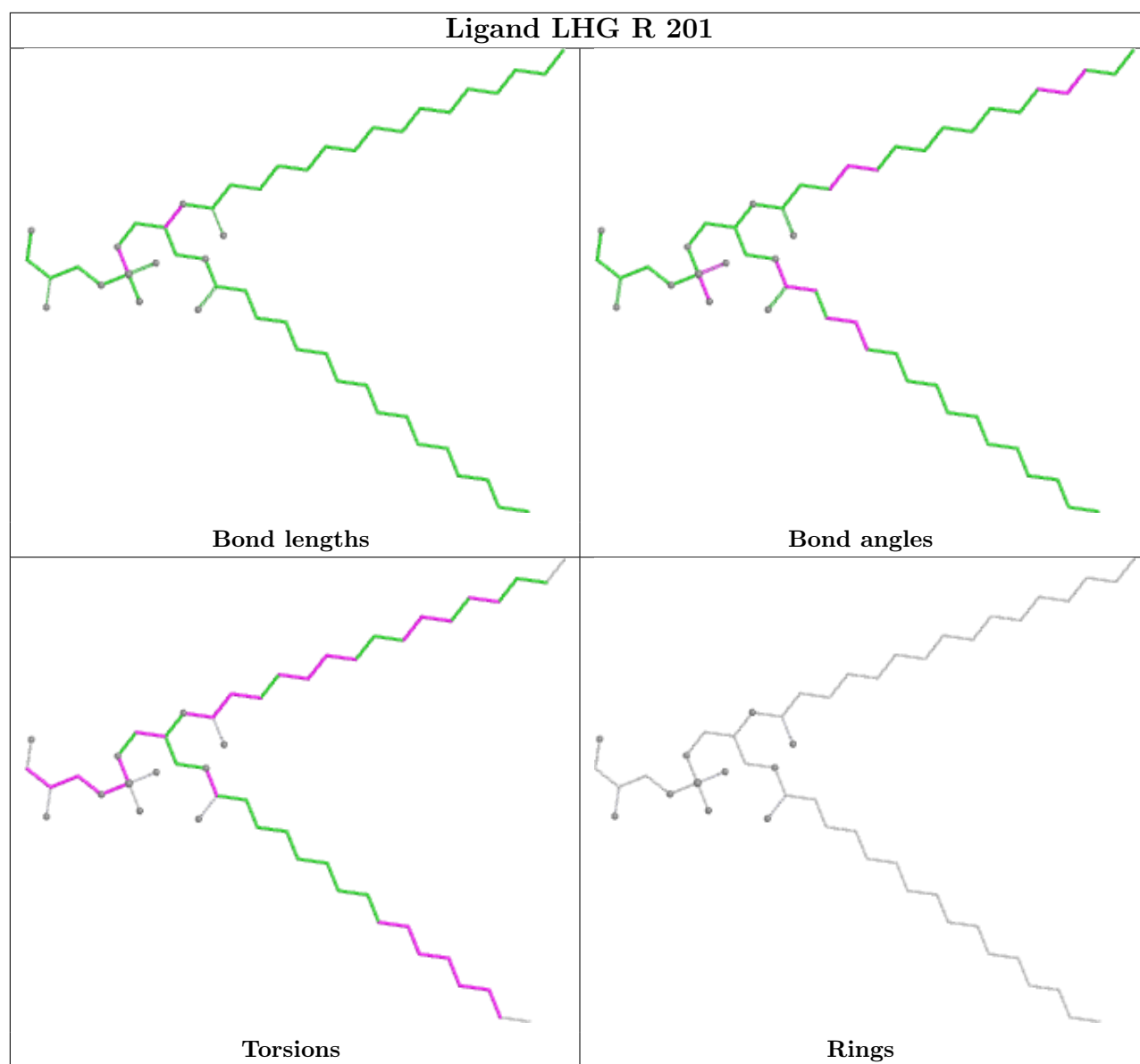
Bond angles

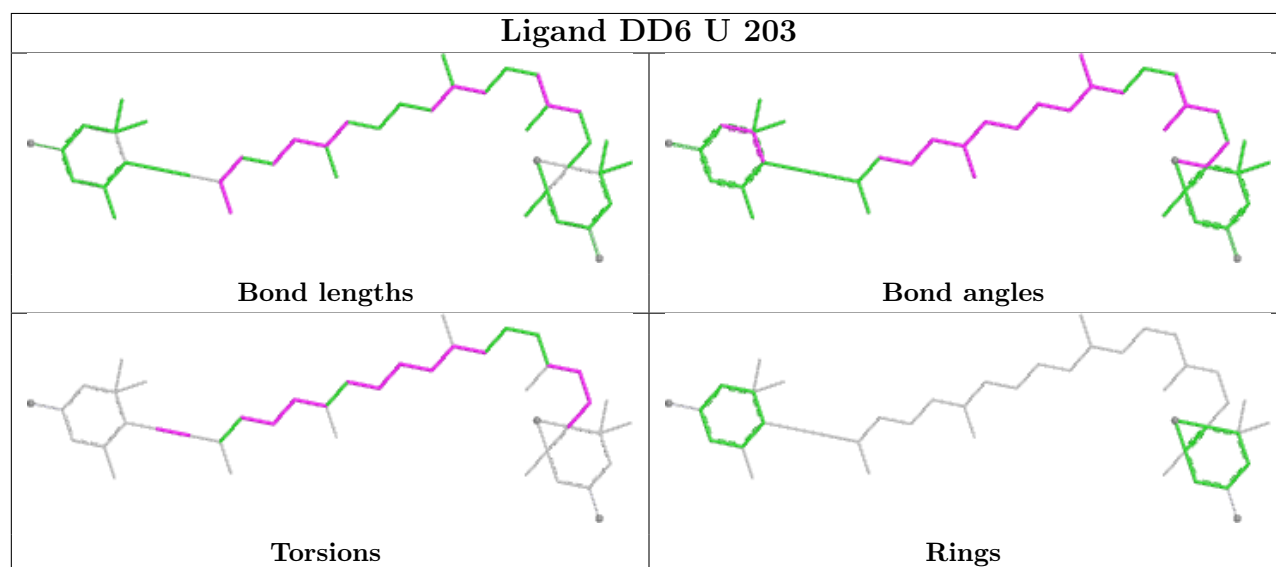
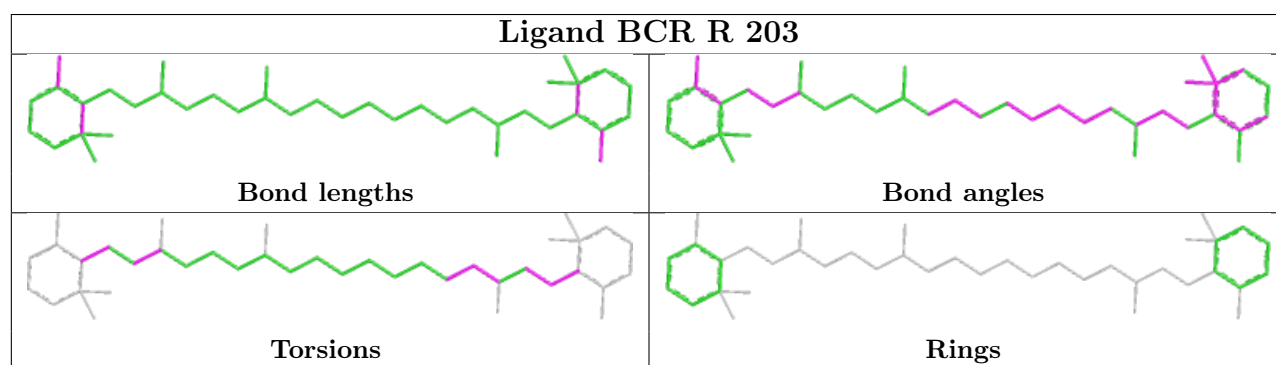
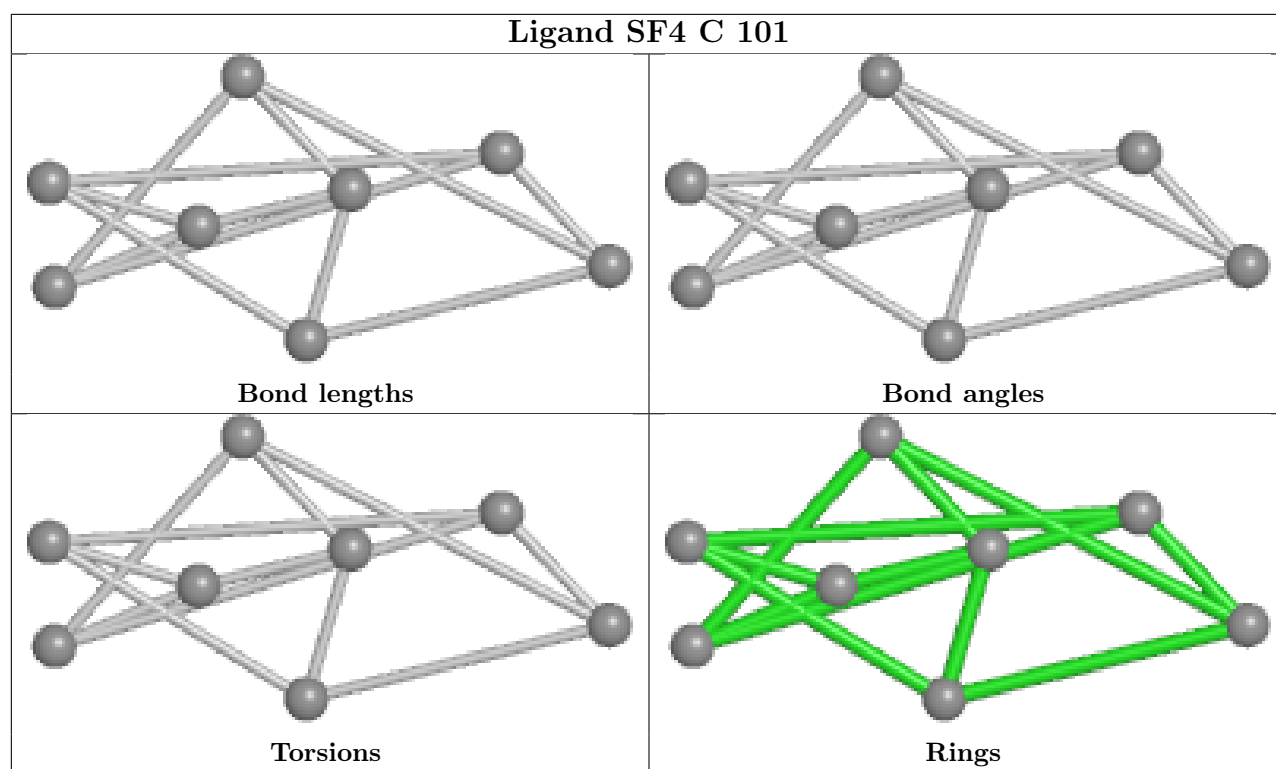


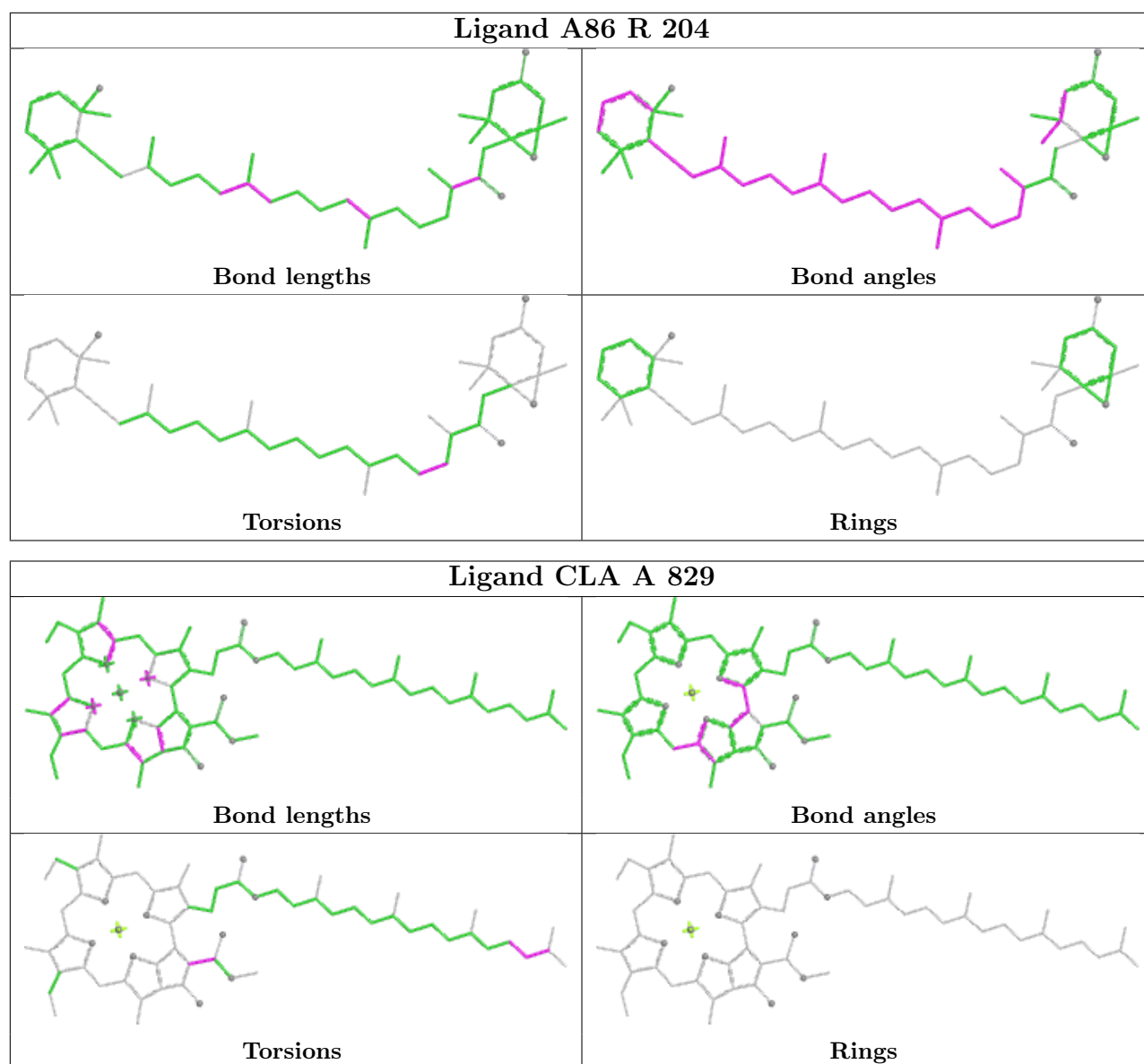
Torsions



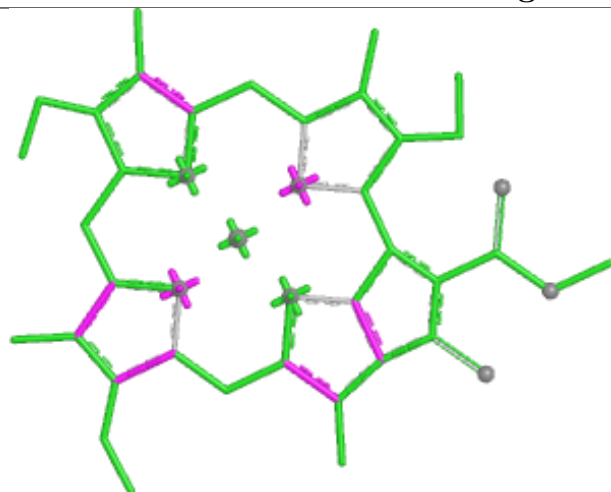
Rings



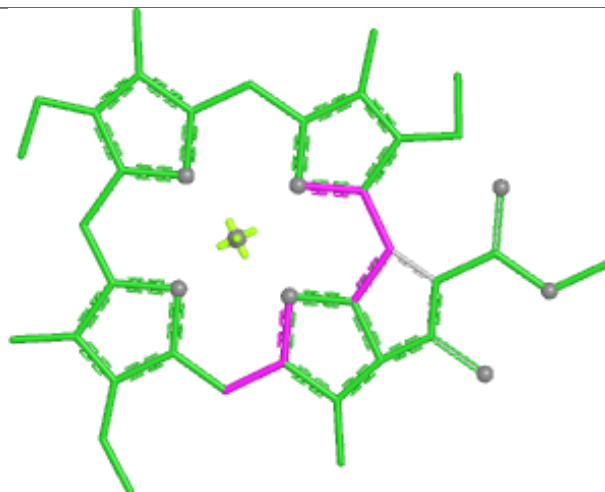




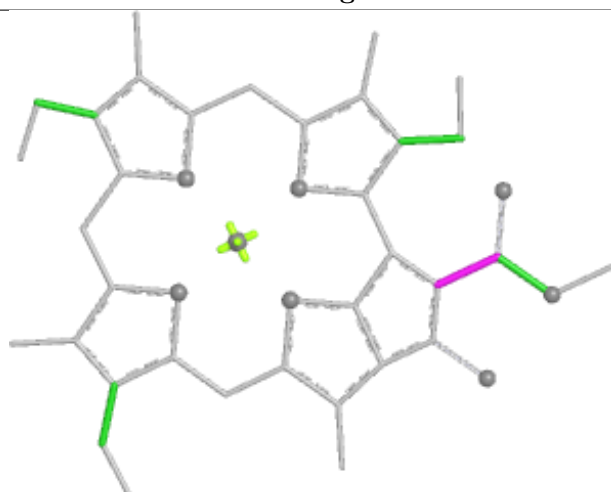
Ligand CLA J 103



Bond lengths



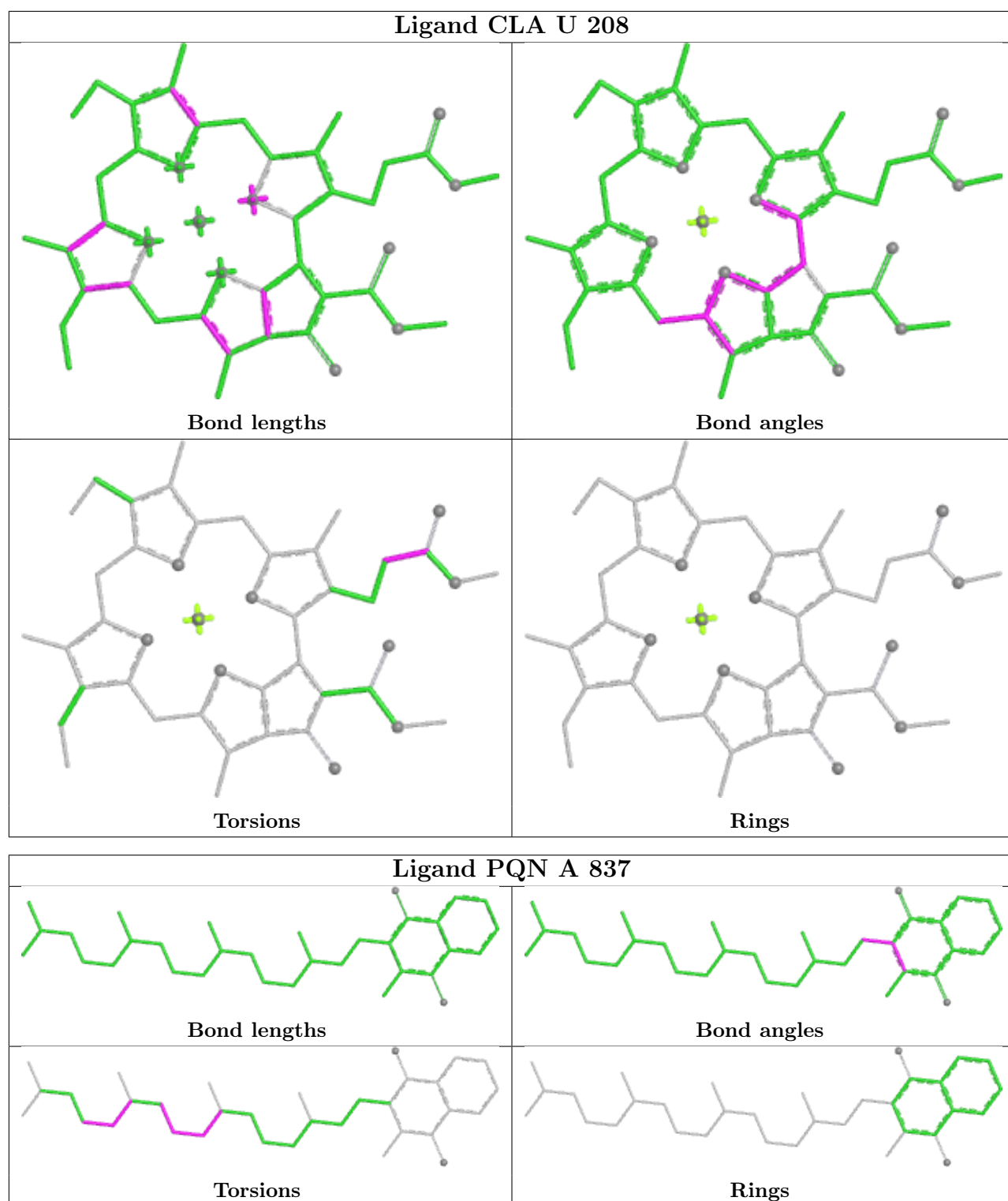
Bond angles



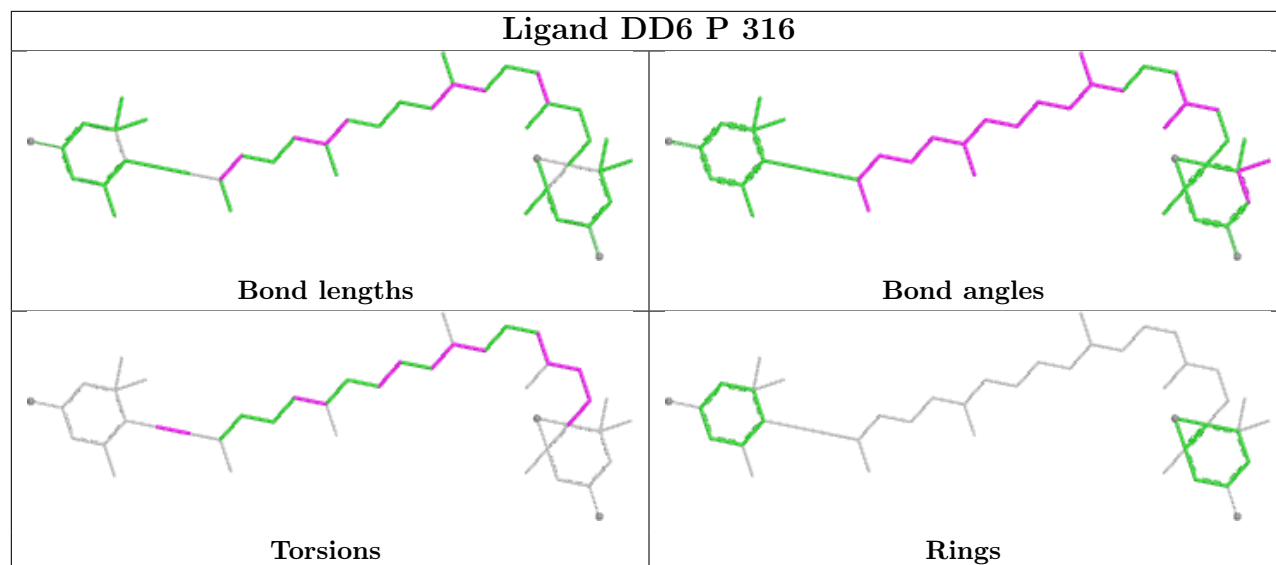
Torsions



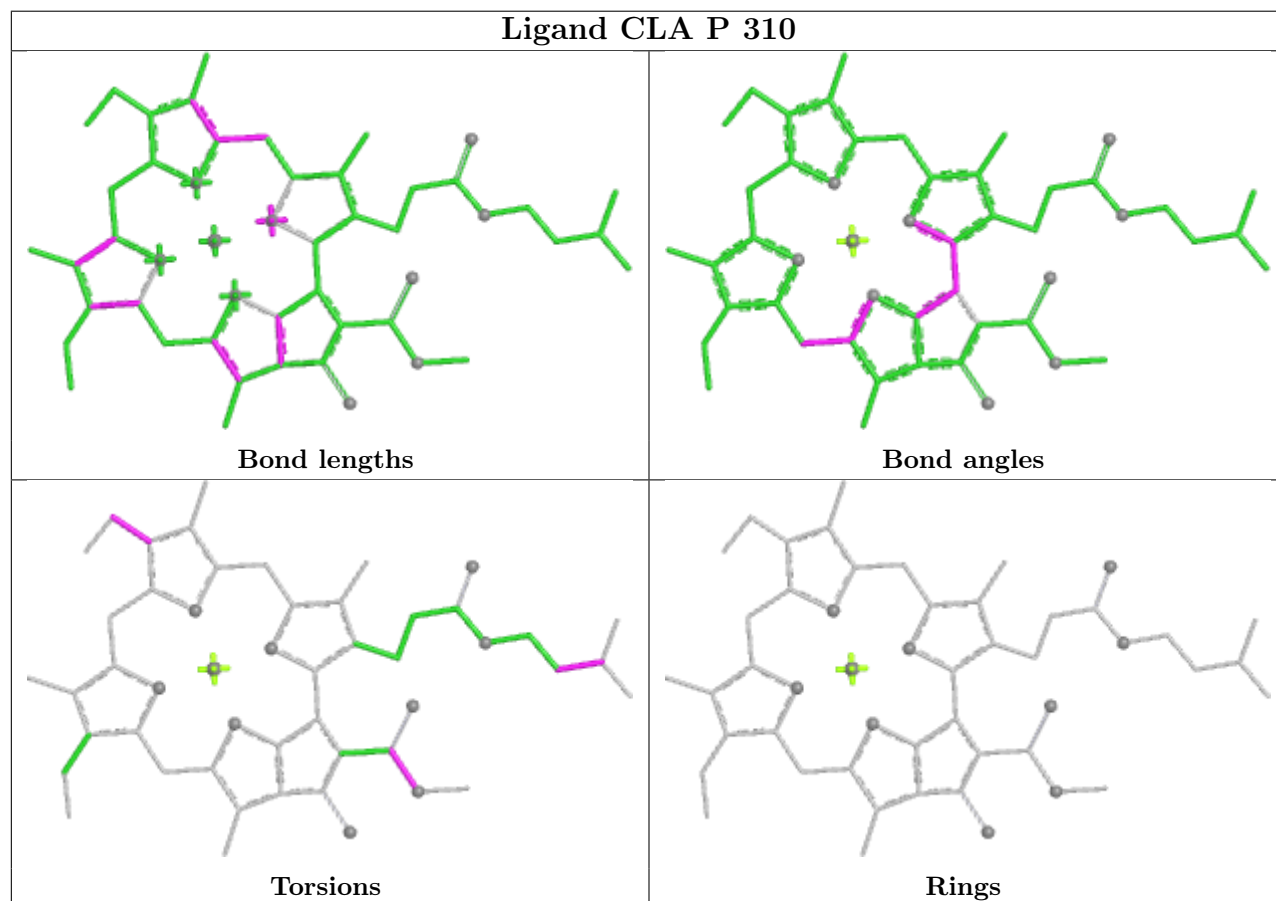
Rings

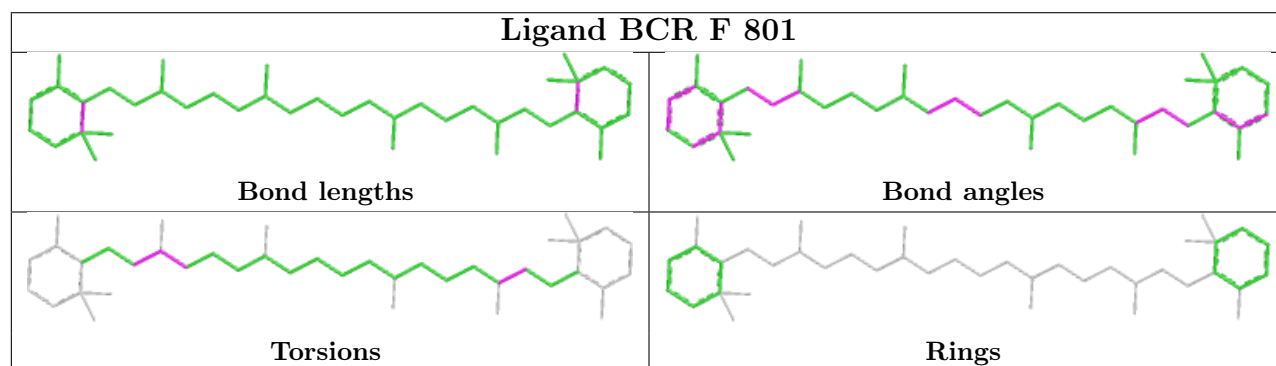
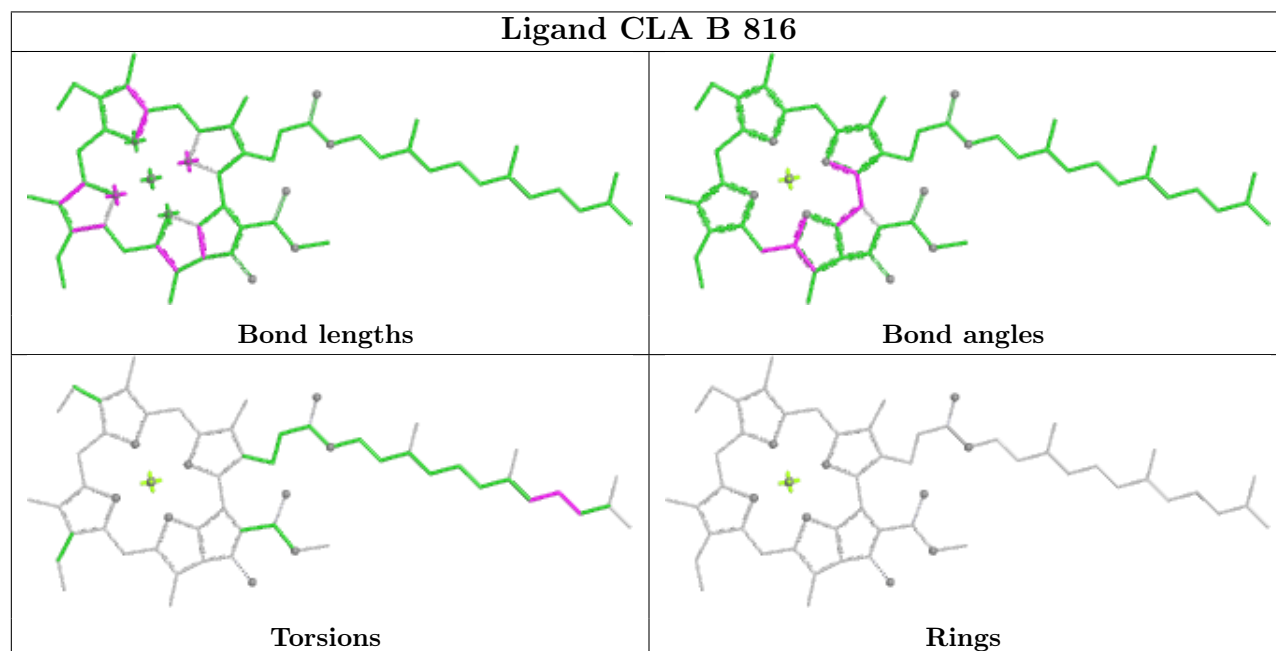
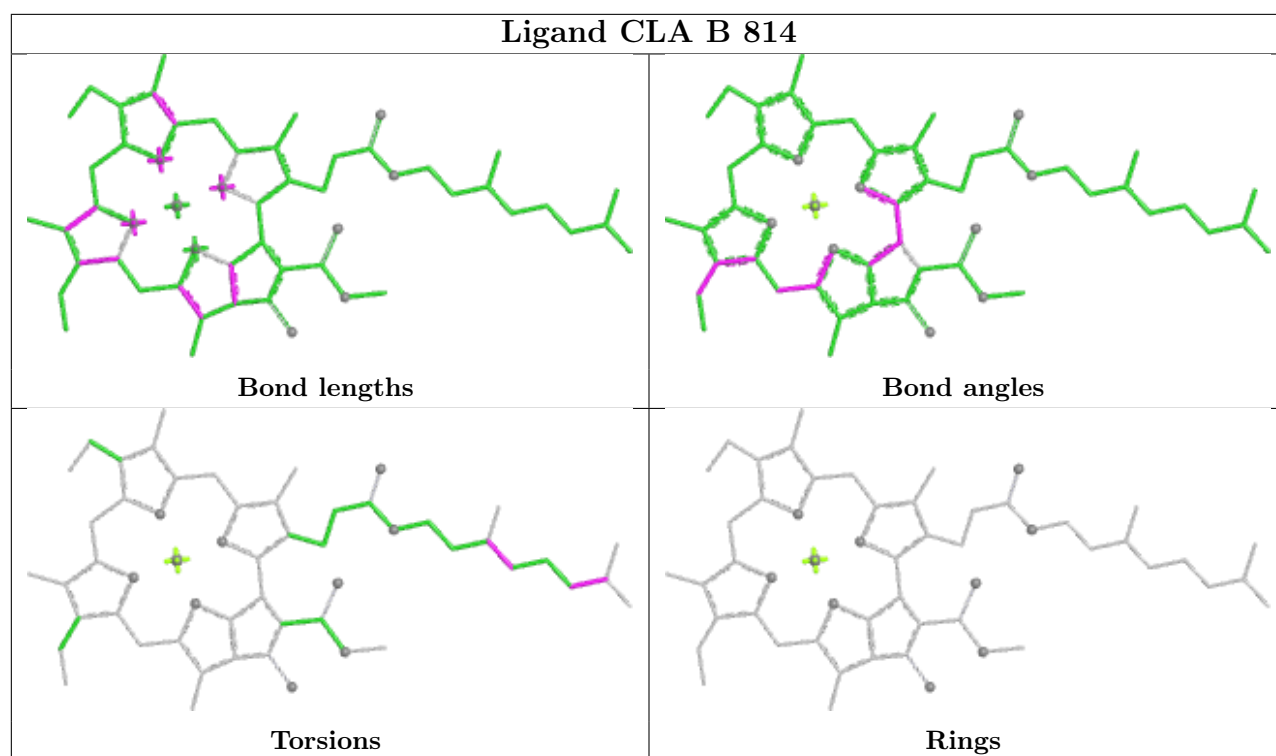


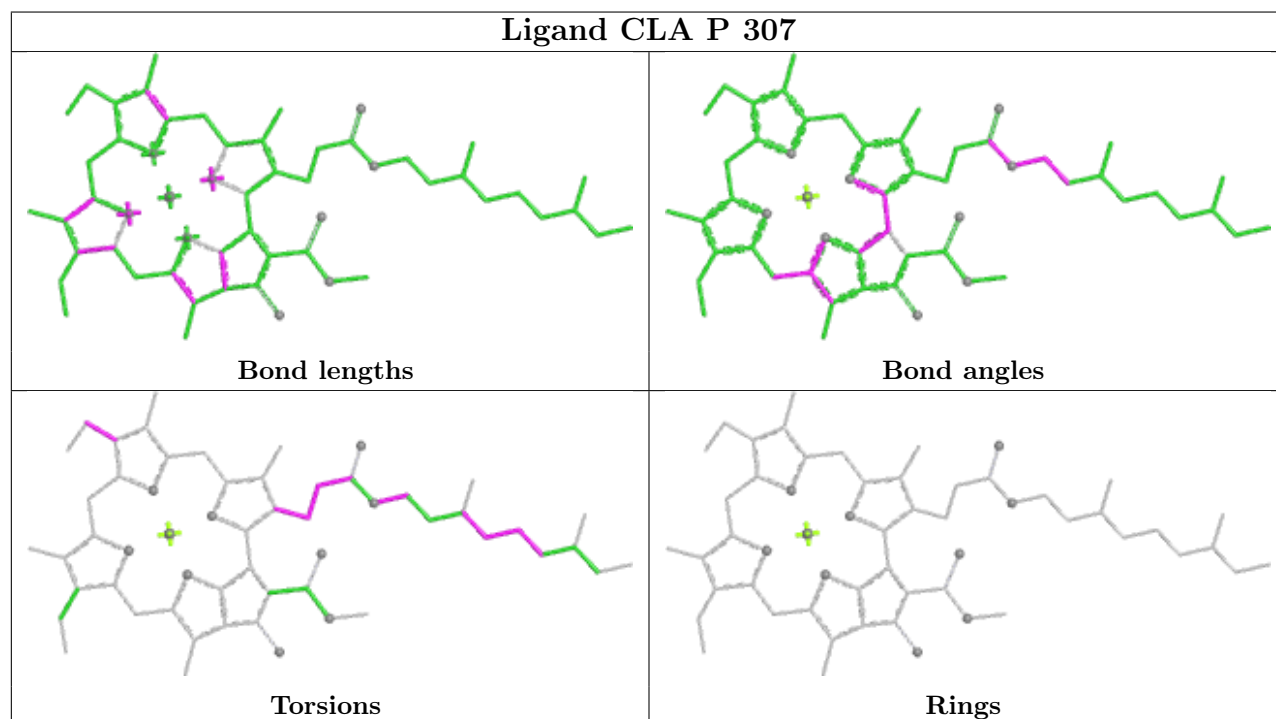
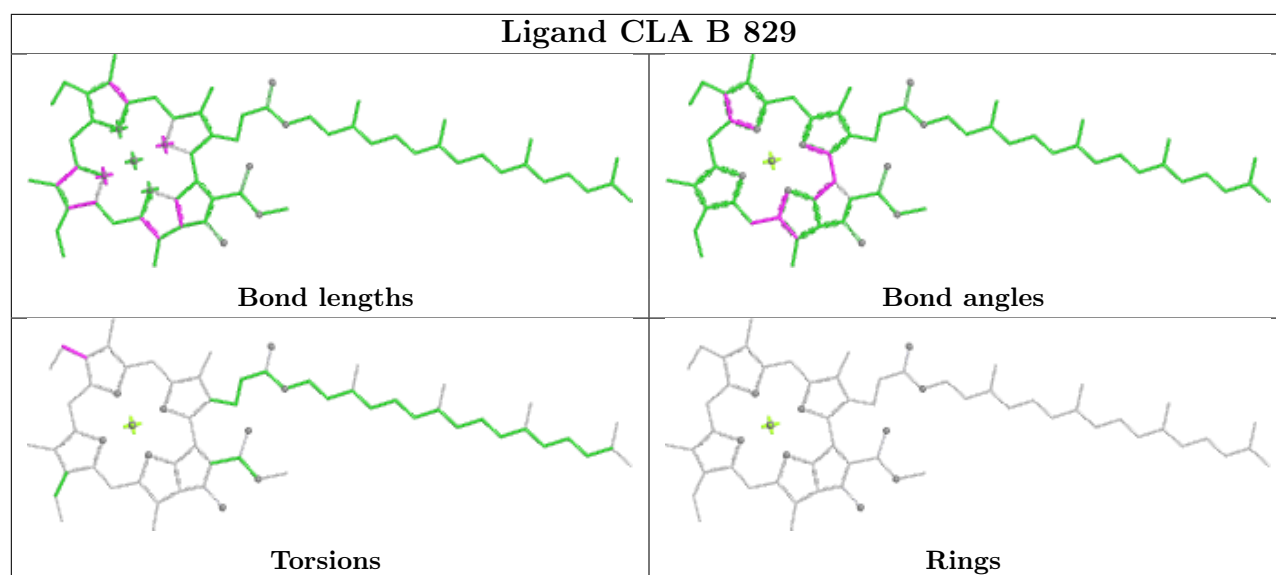
Ligand DD6 P 316



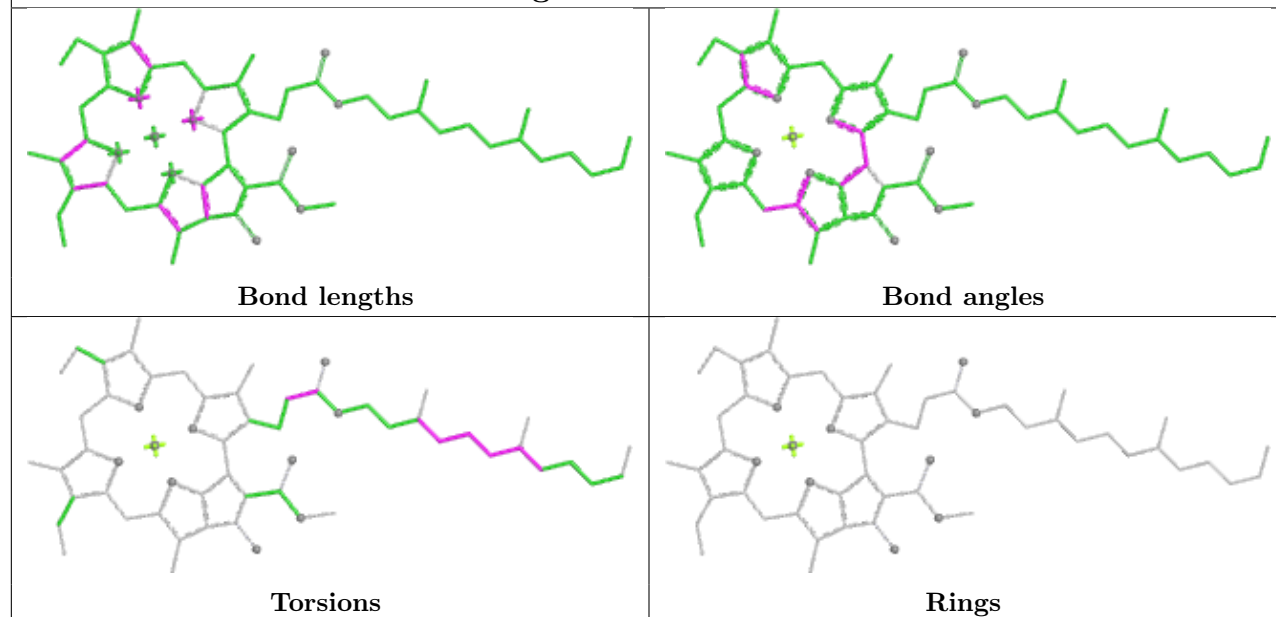
Ligand CLA P 310



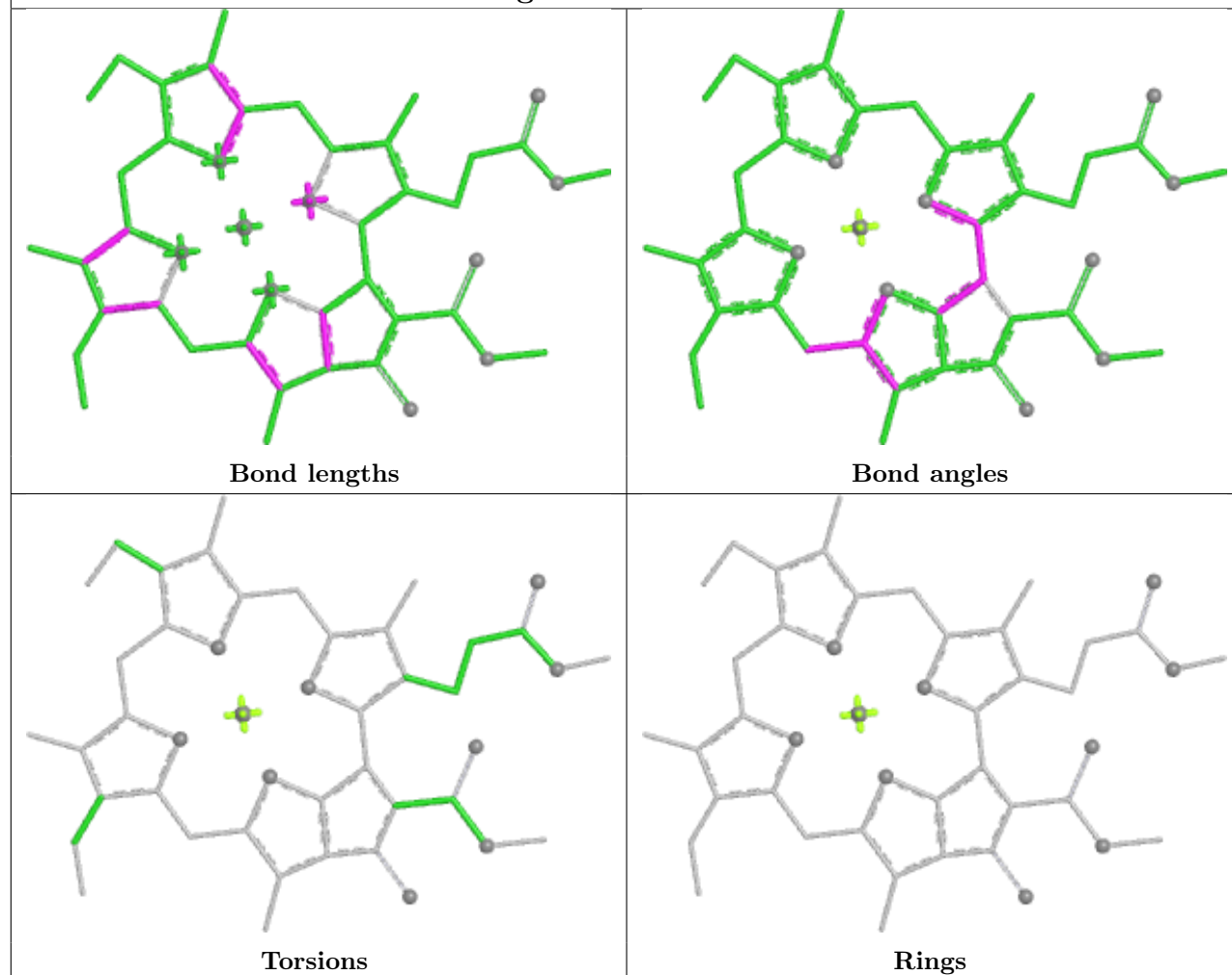




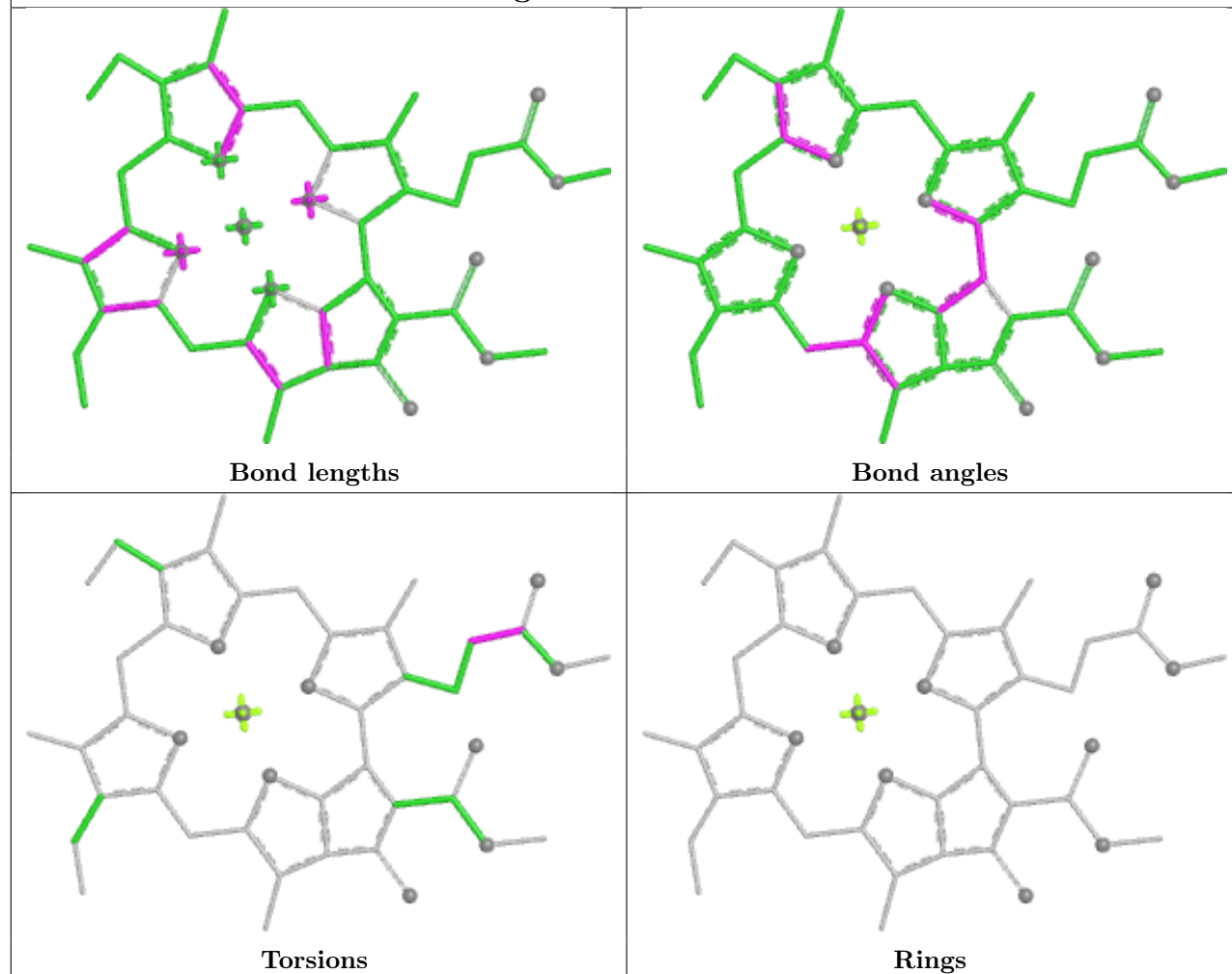
Ligand CLA B 815



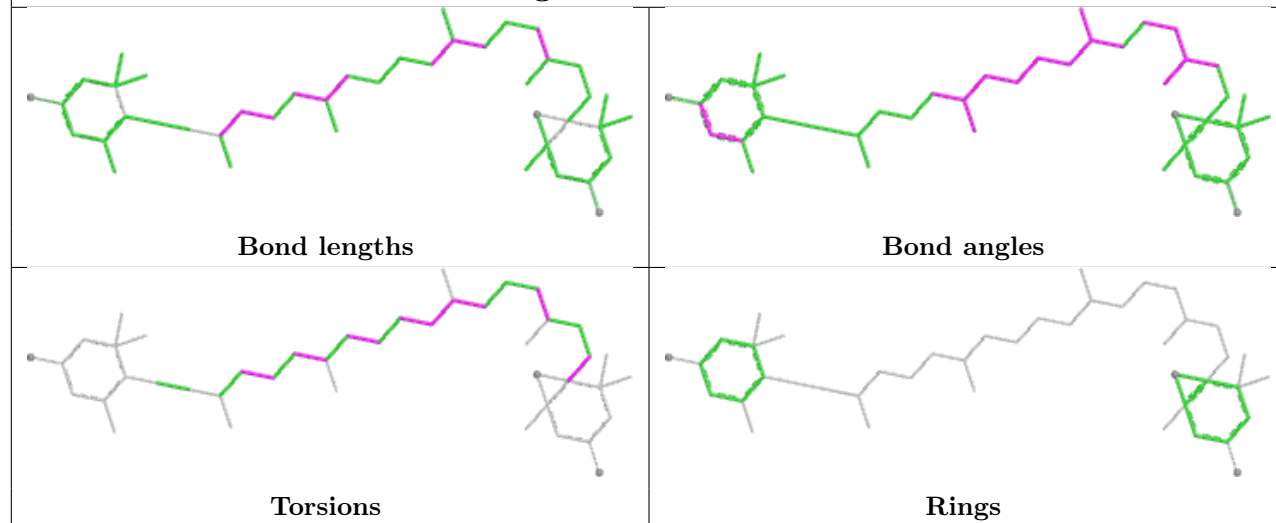
Ligand CLA S 204

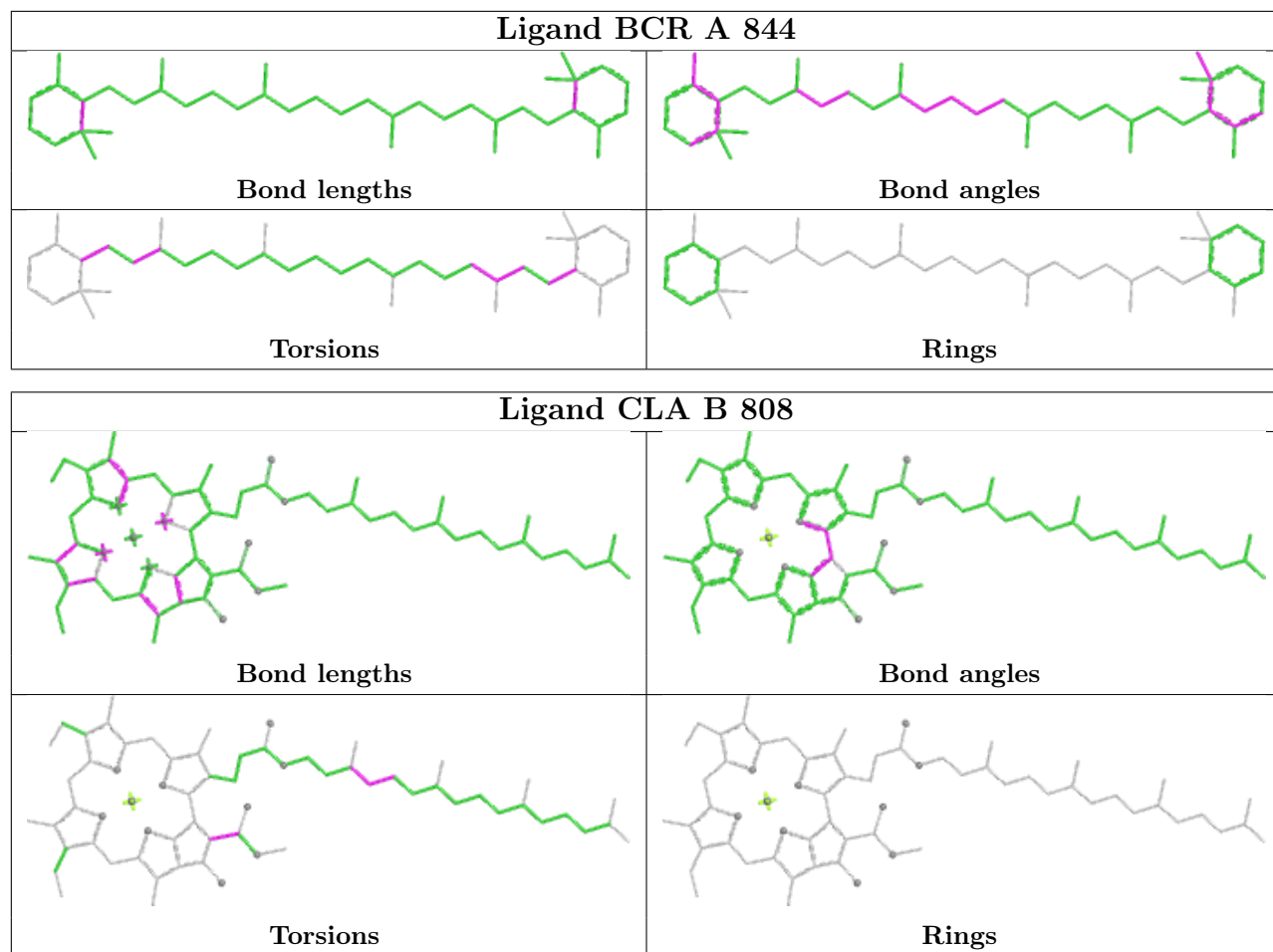


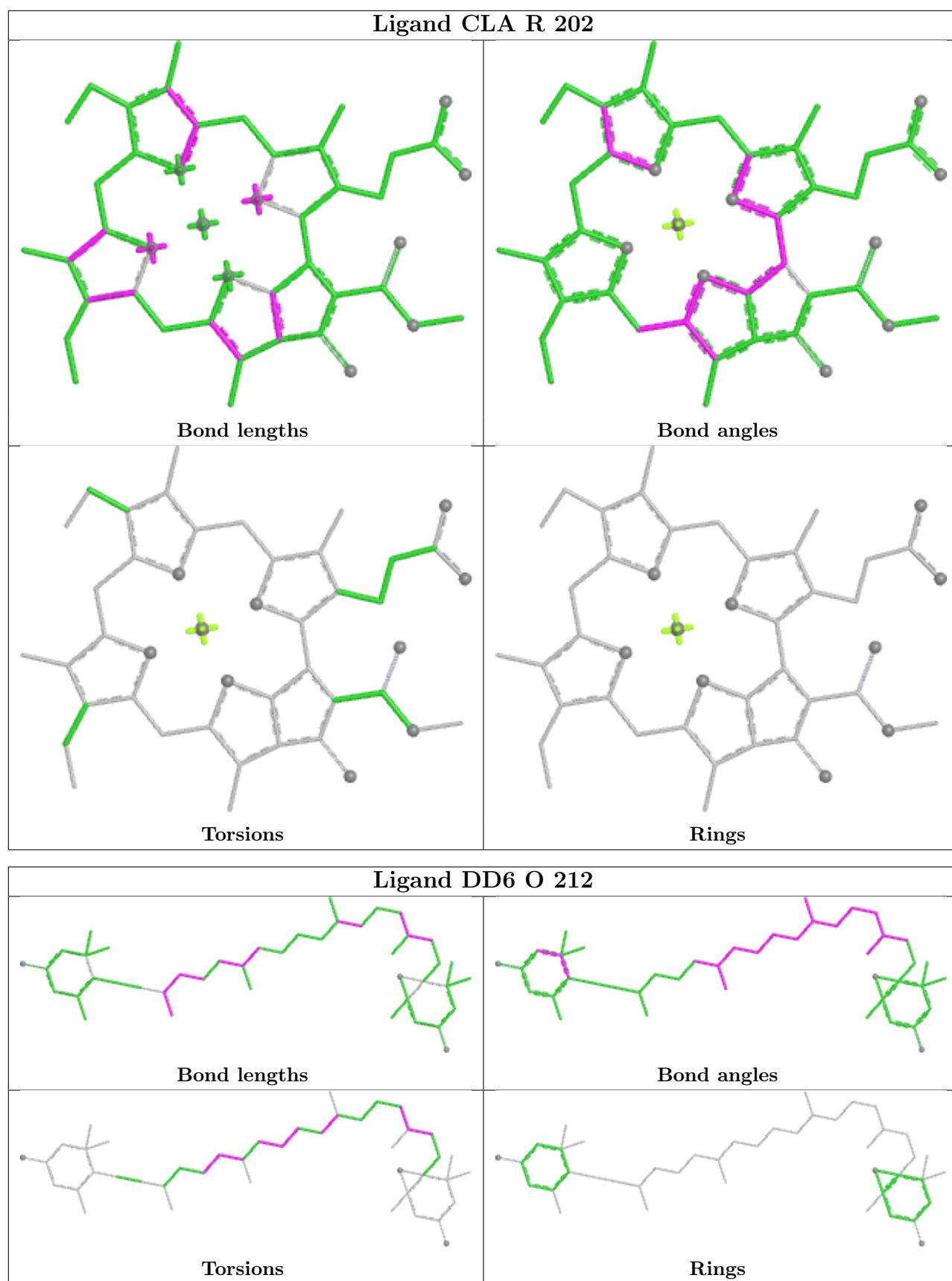
Ligand CLA B 818



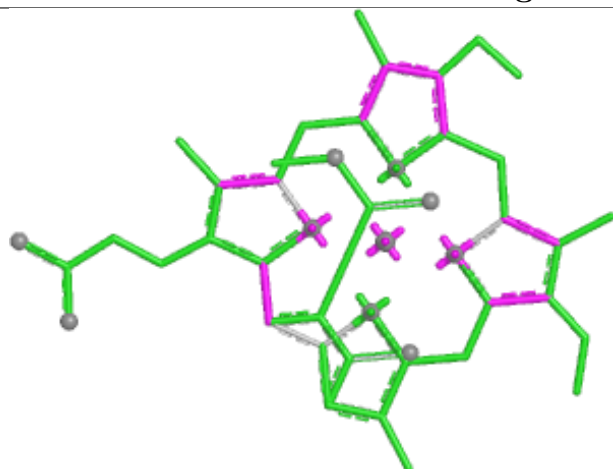
Ligand DD6 T 213



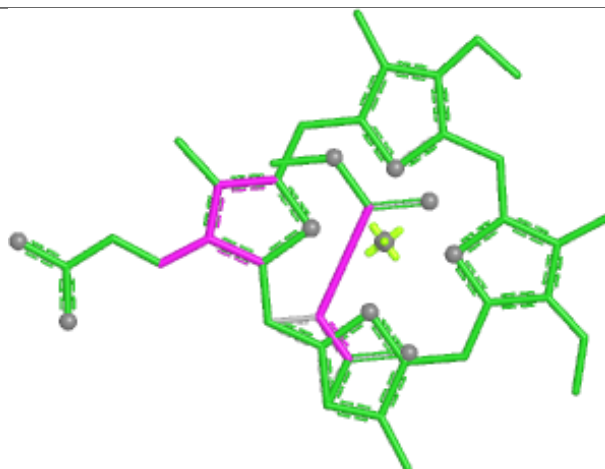




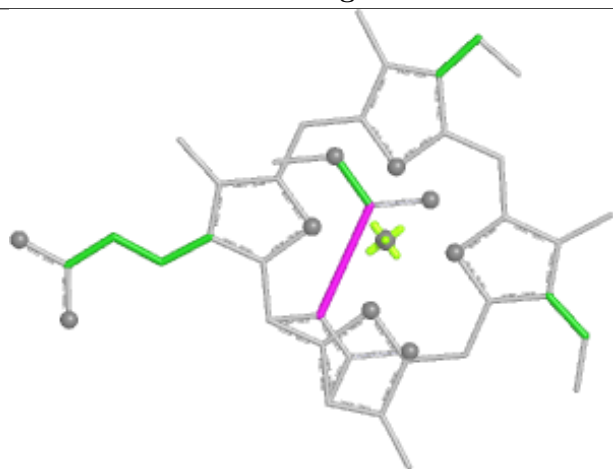
Ligand KC1 U 213



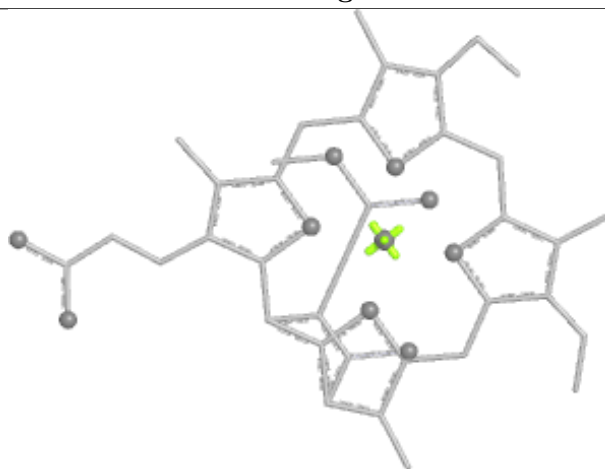
Bond lengths



Bond angles

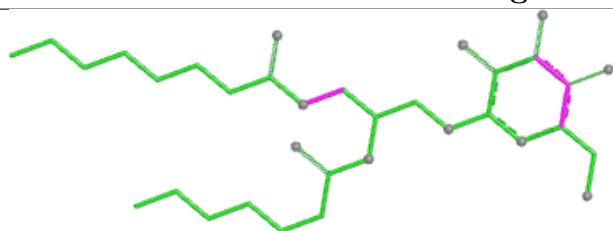


Torsions

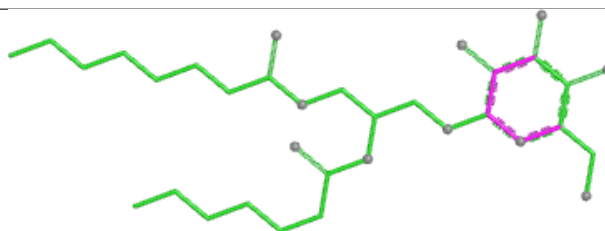


Rings

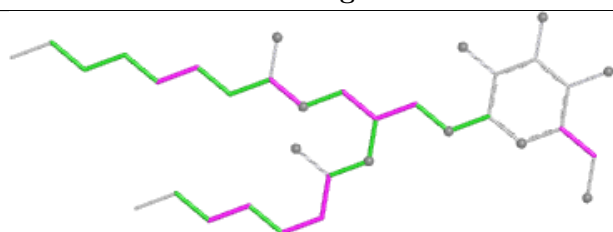
Ligand LMG P 301



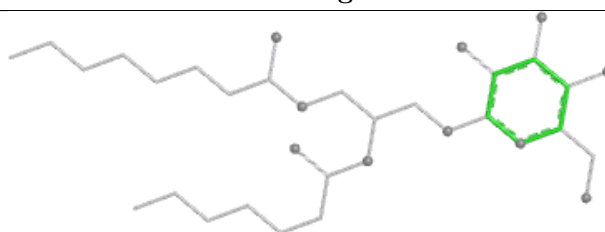
Bond lengths



Bond angles

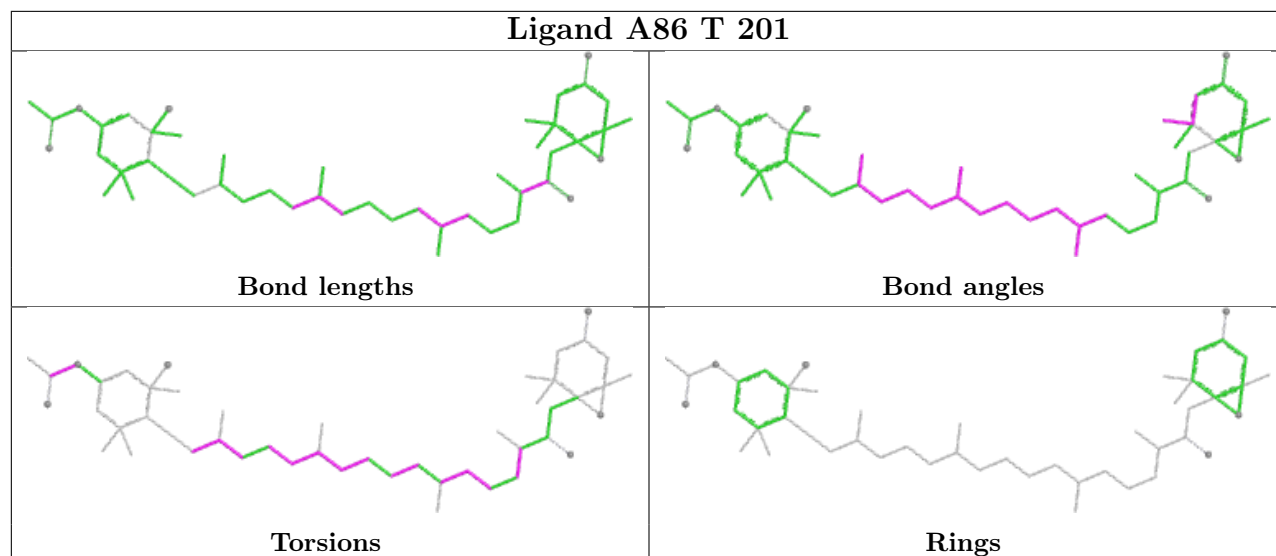


Torsions

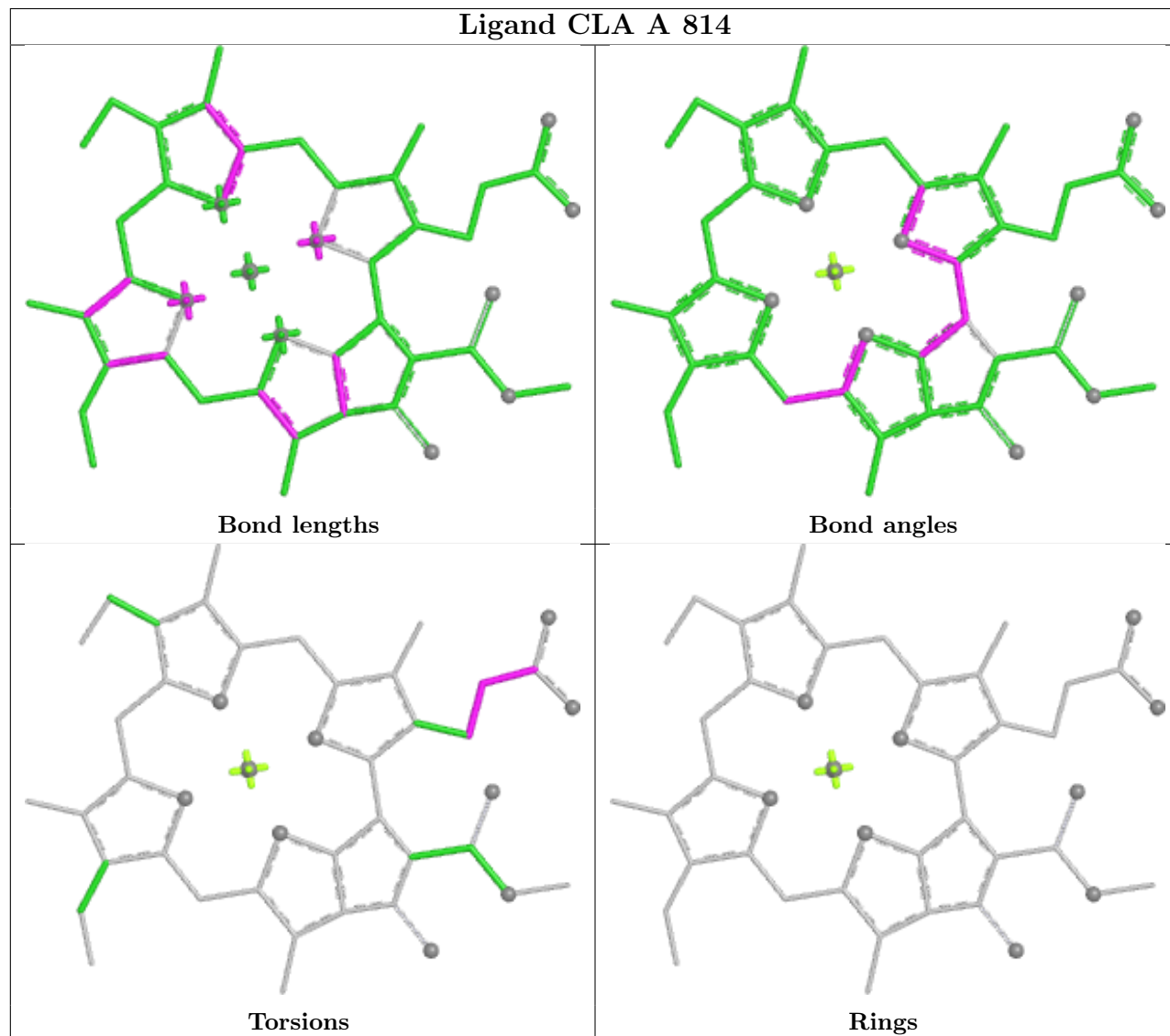


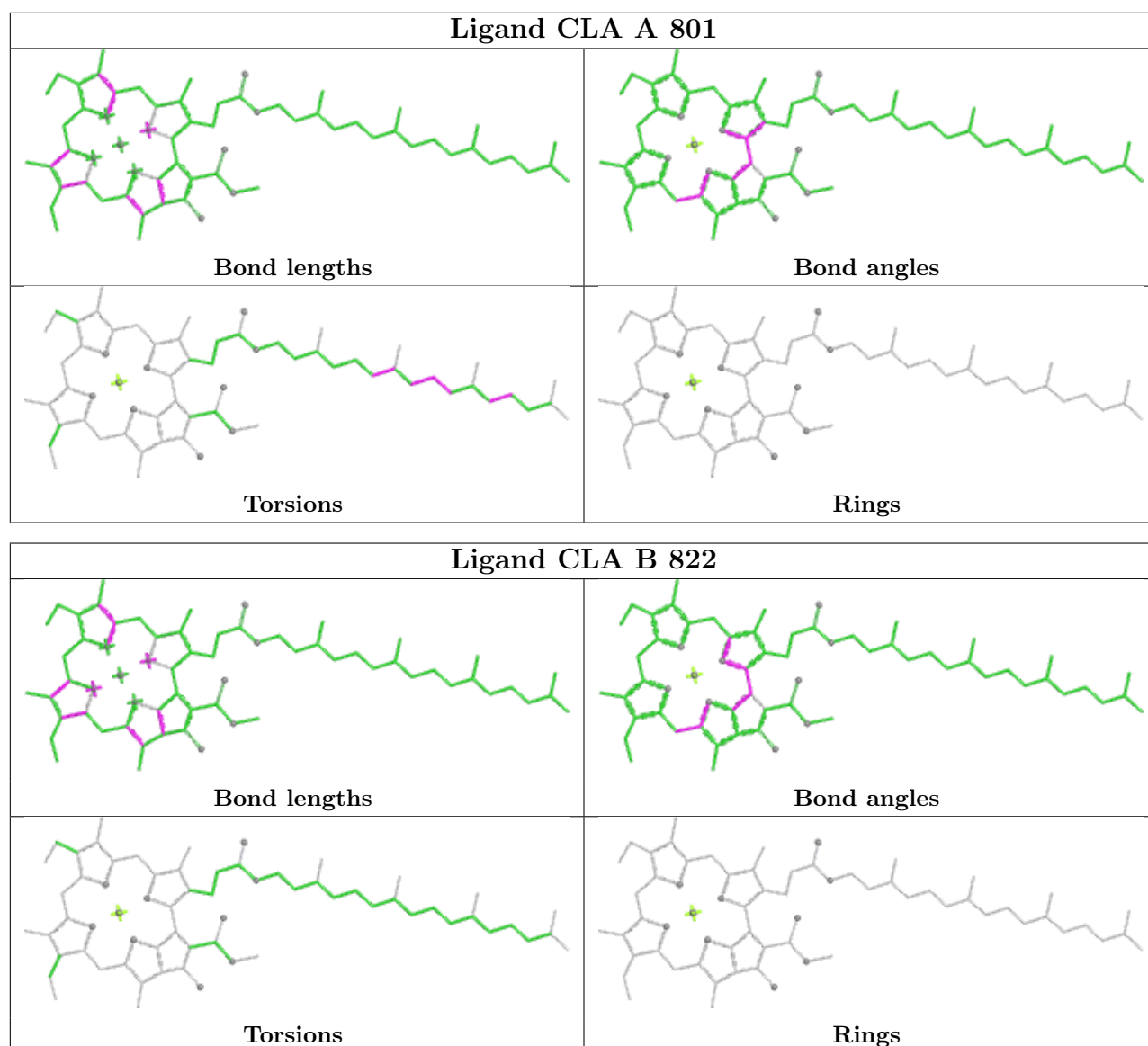
Rings

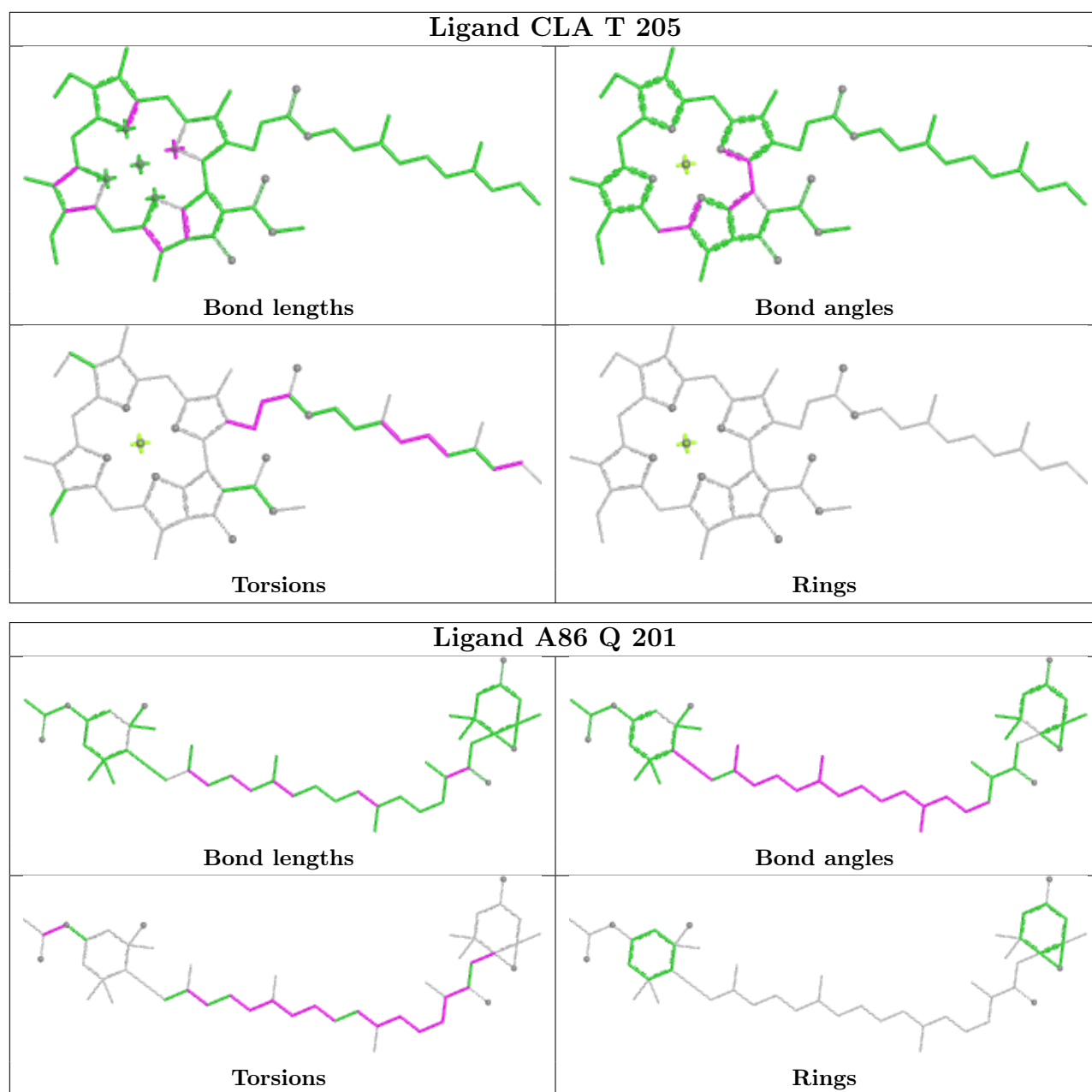
Ligand A86 T 201

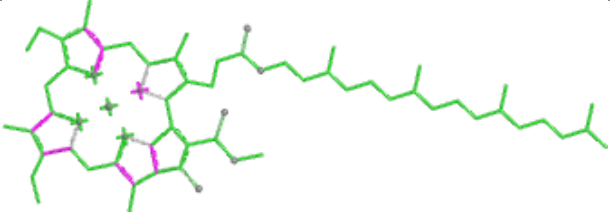
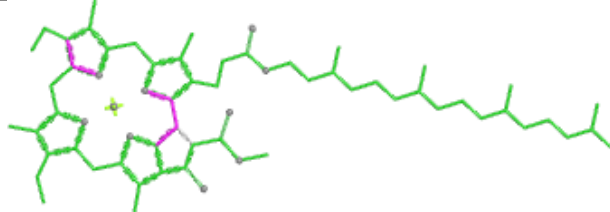
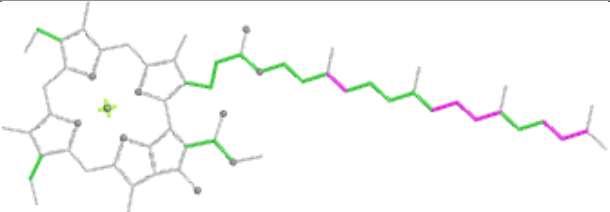
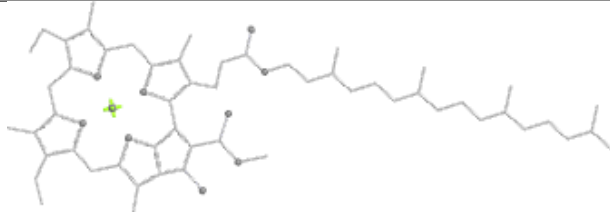
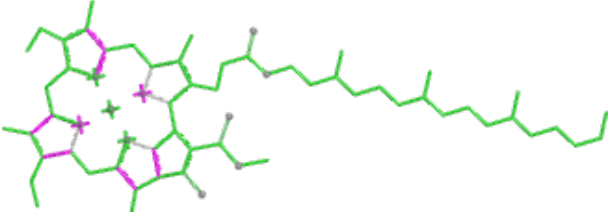
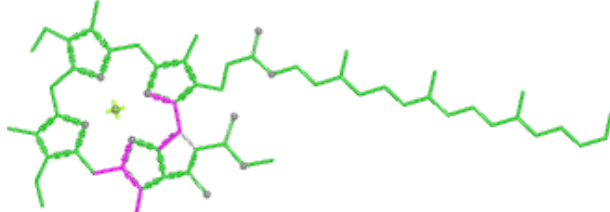
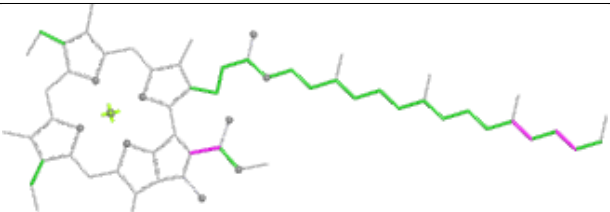
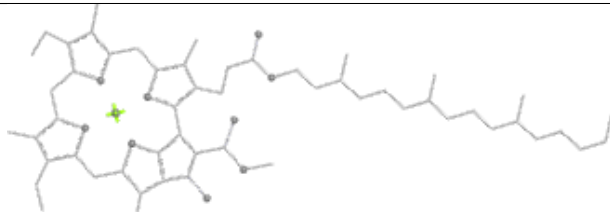

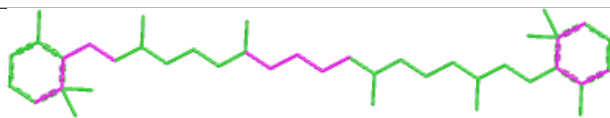
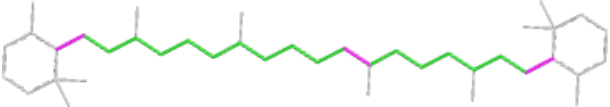
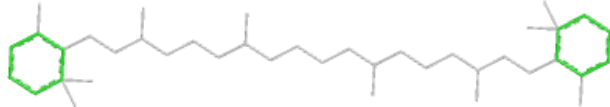


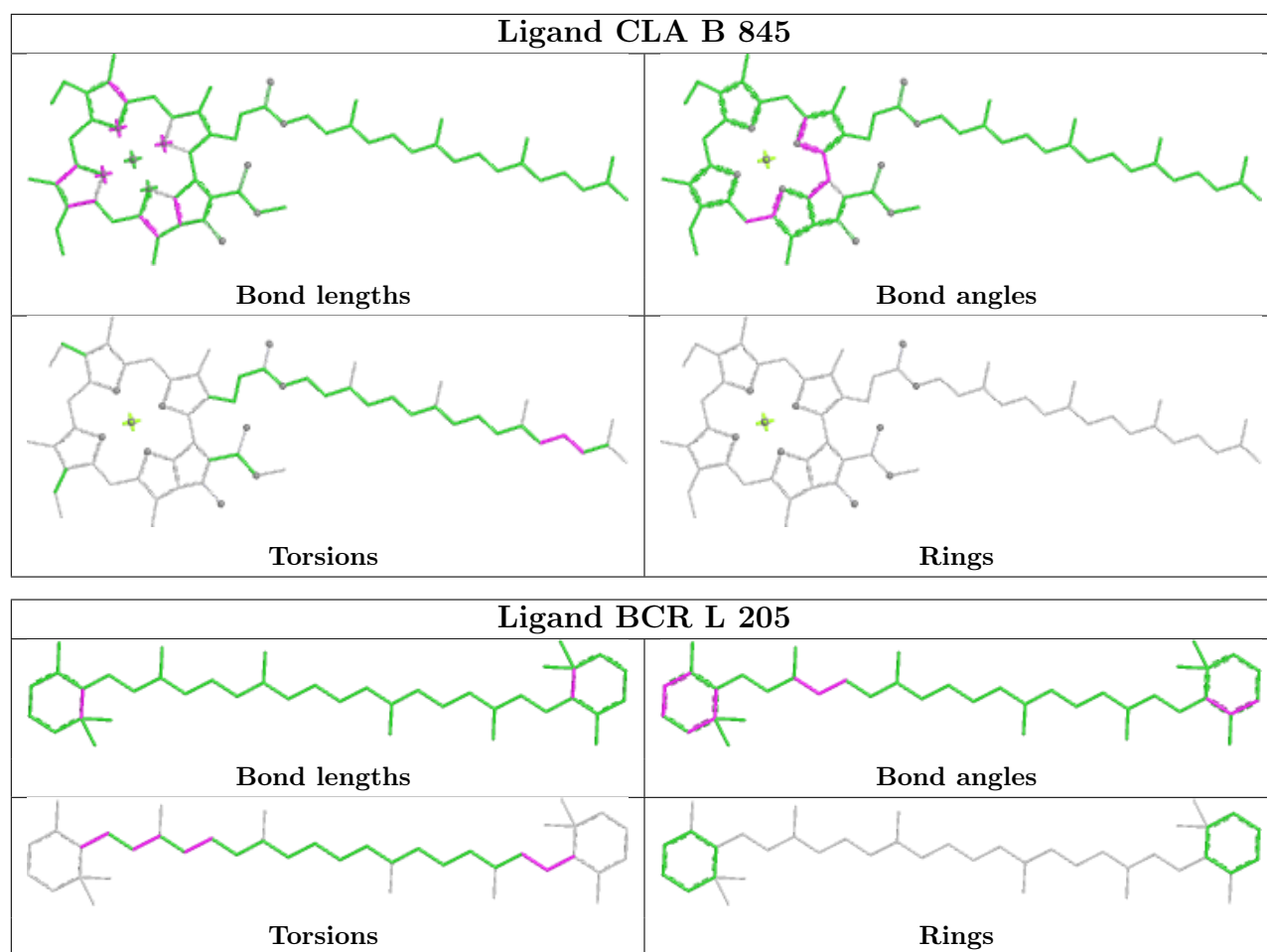
Ligand CLA A 814



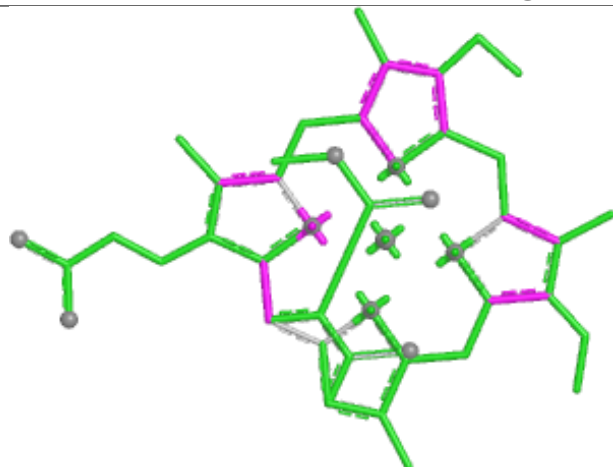




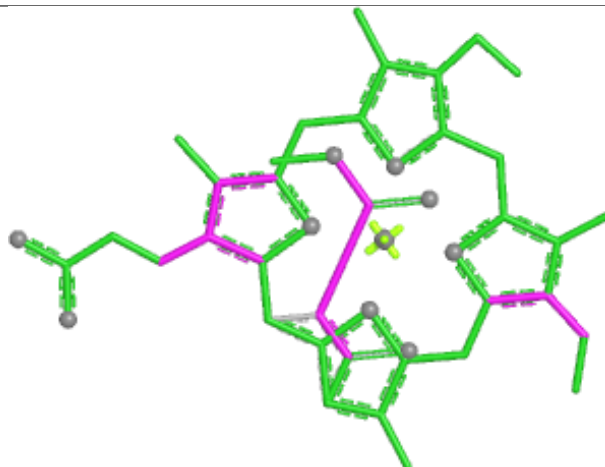
Ligand CLA B 817	
 <p>Bond lengths</p>	 <p>Bond angles</p>
 <p>Torsions</p>	 <p>Rings</p>
Ligand CLA B 821	
 <p>Bond lengths</p>	 <p>Bond angles</p>
 <p>Torsions</p>	 <p>Rings</p>
Ligand BCR B 840	
 <p>Bond lengths</p>	 <p>Bond angles</p>
 <p>Torsions</p>	 <p>Rings</p>



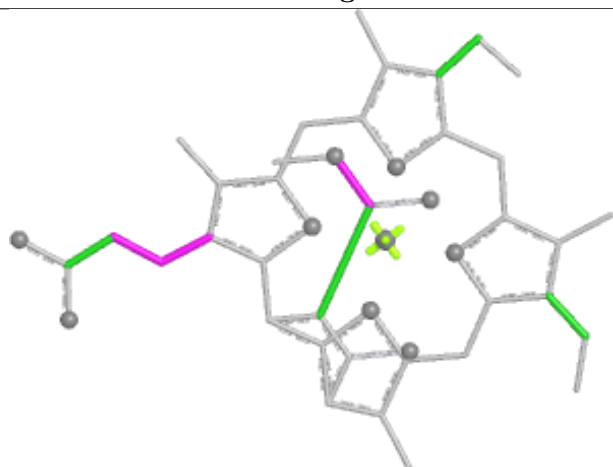
Ligand KC1 S 210



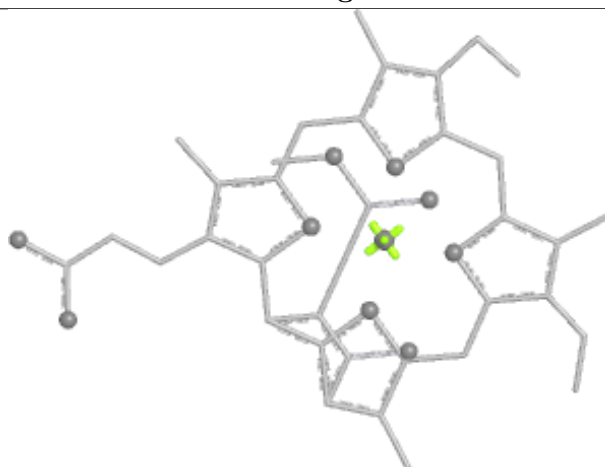
Bond lengths



Bond angles

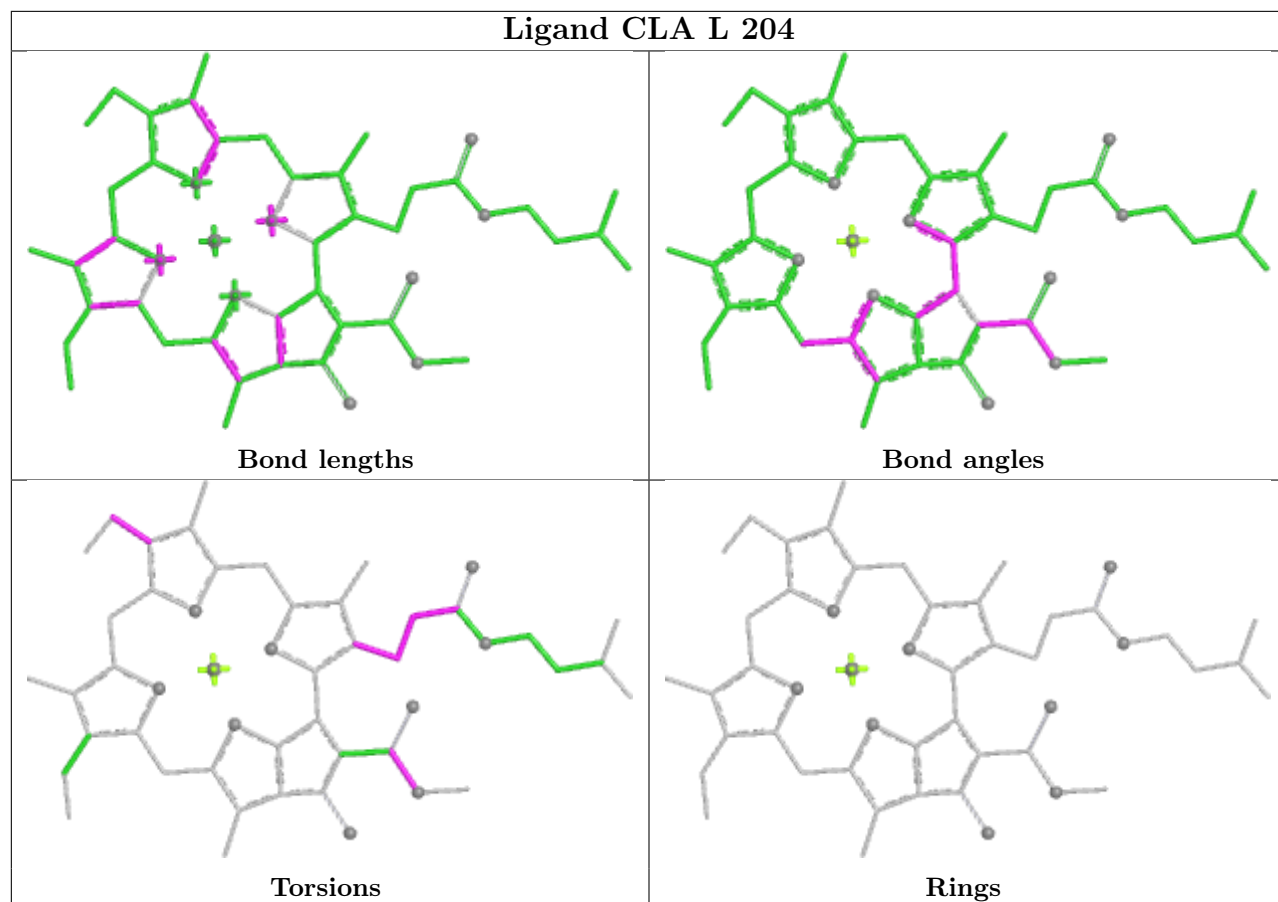


Torsions

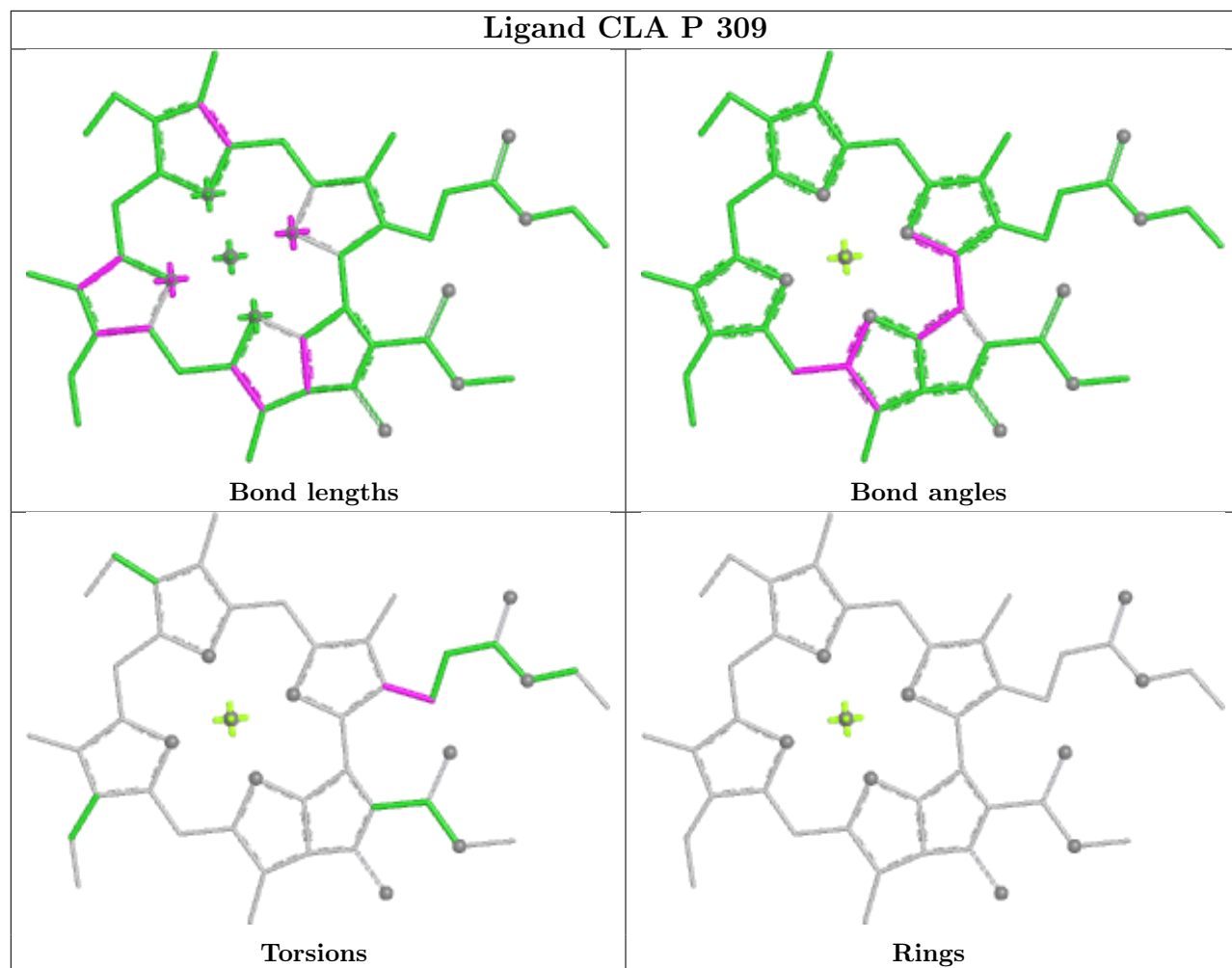


Rings

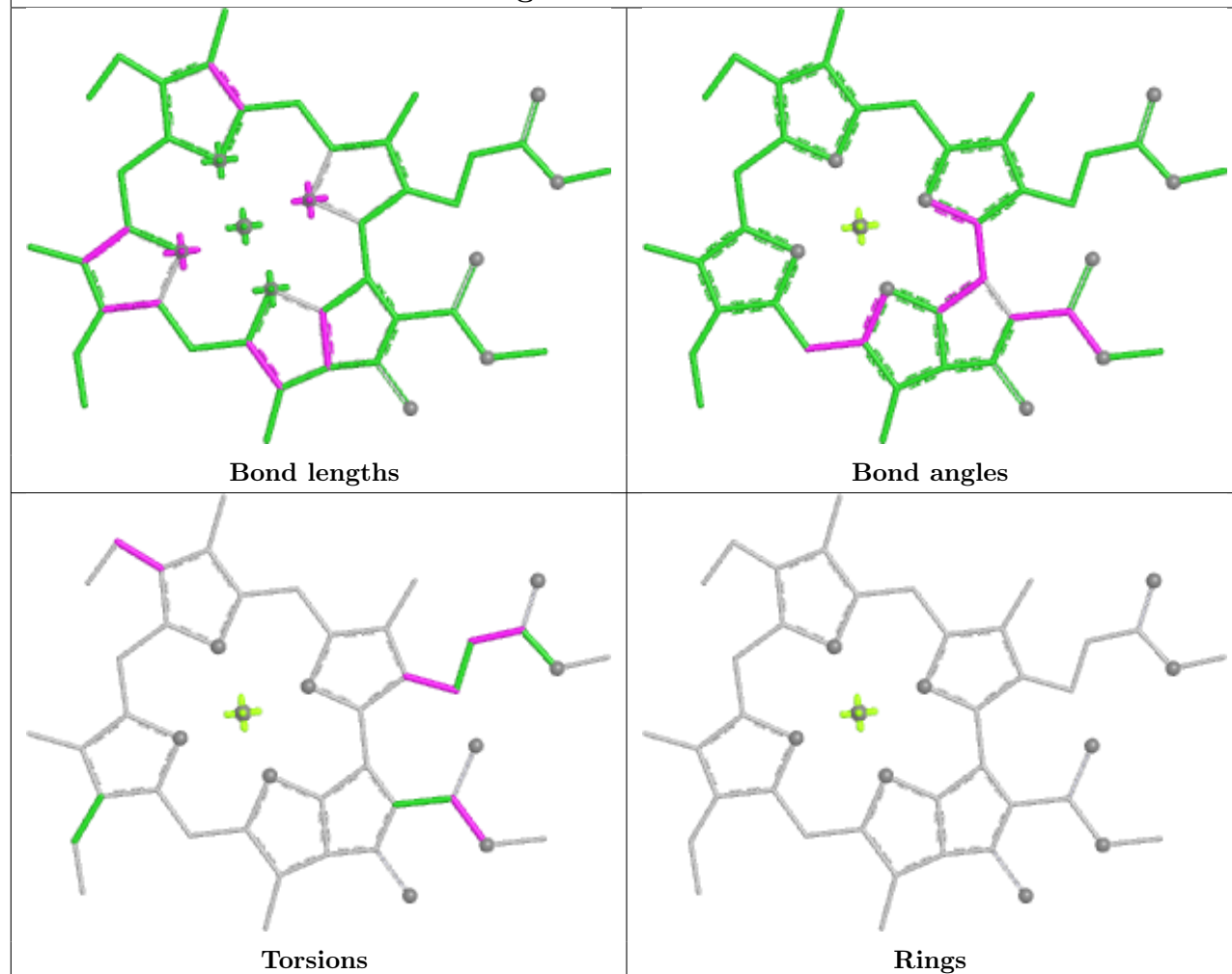
Ligand CLA L 204



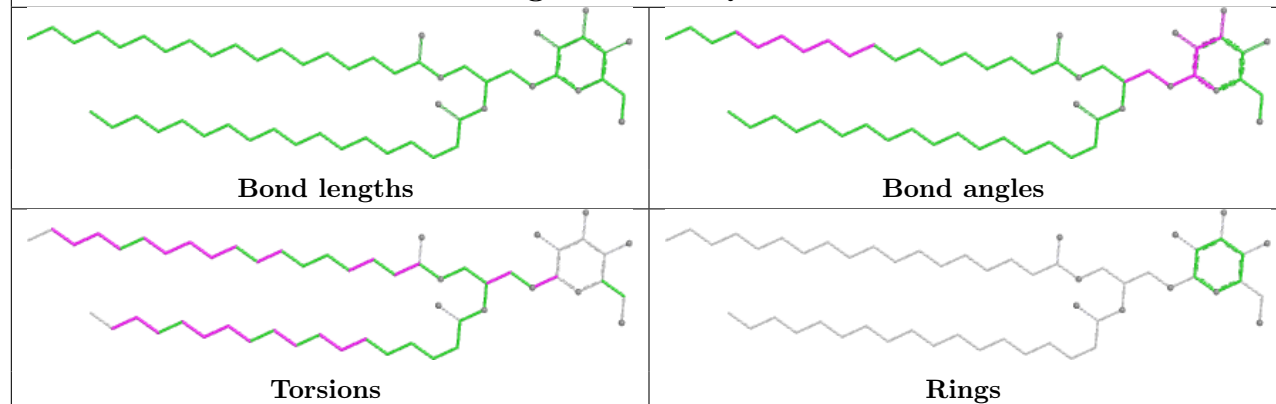
Ligand CLA P 309

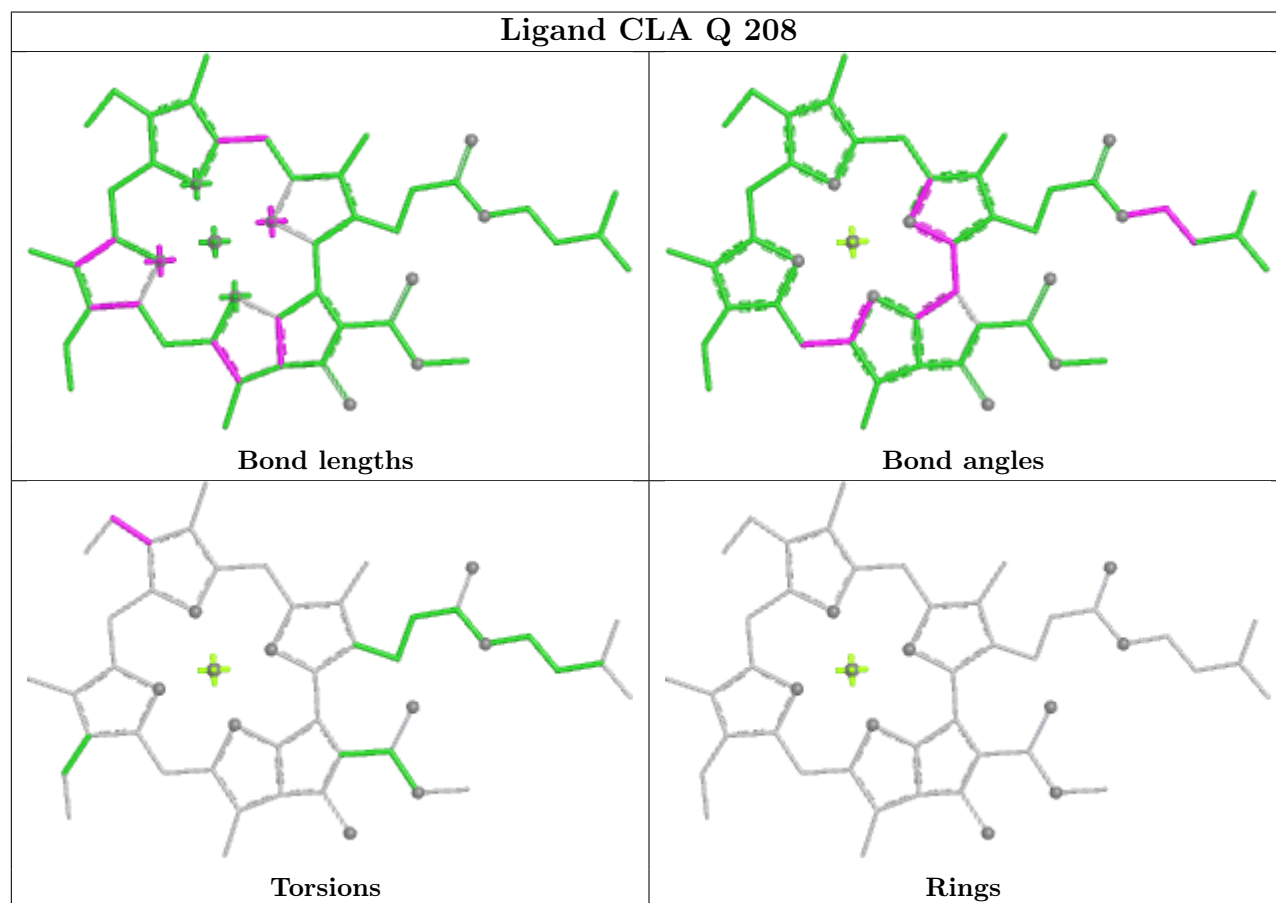
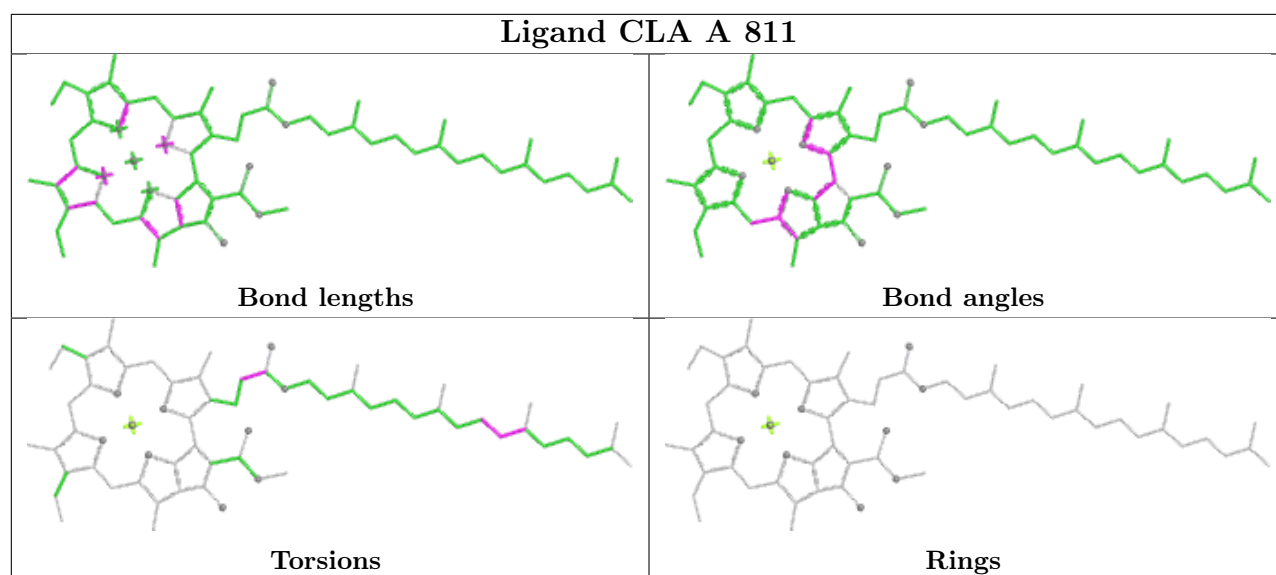


Ligand CLA S 205

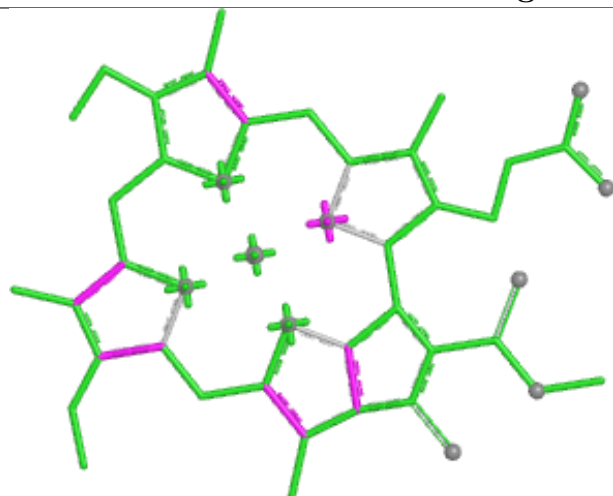


Ligand LMG Q 217

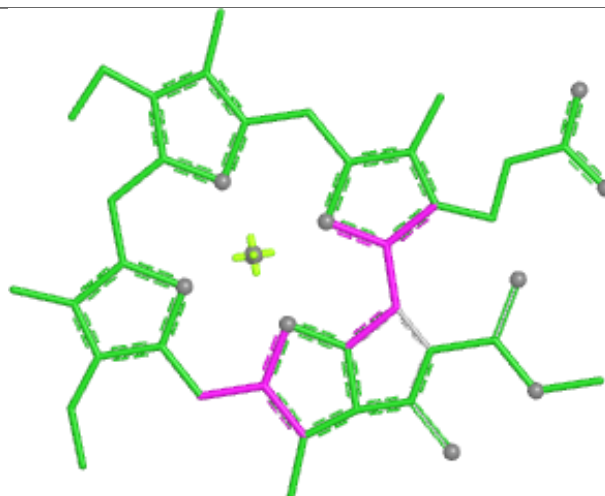




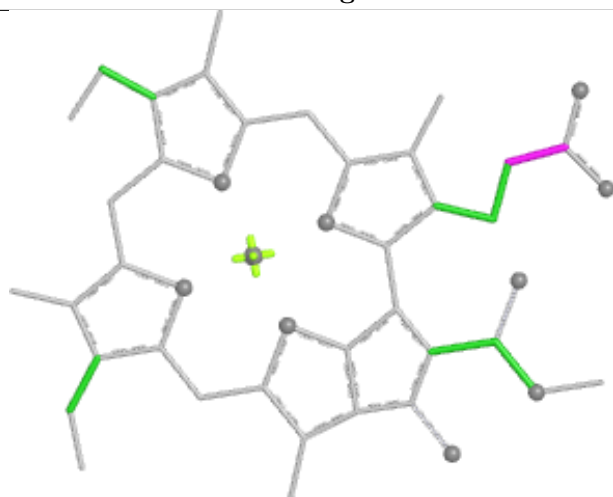
Ligand CLA S 206



Bond lengths



Bond angles

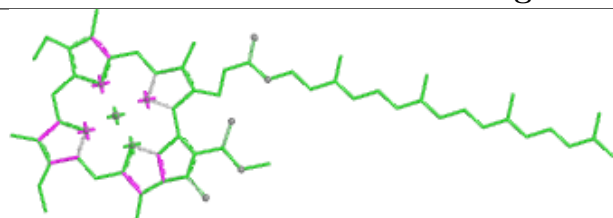


Torsions

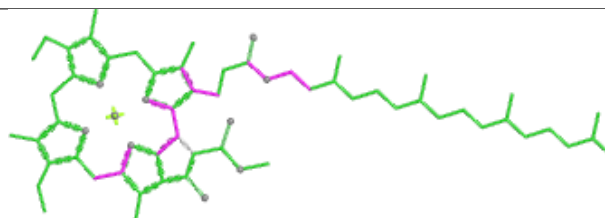


Rings

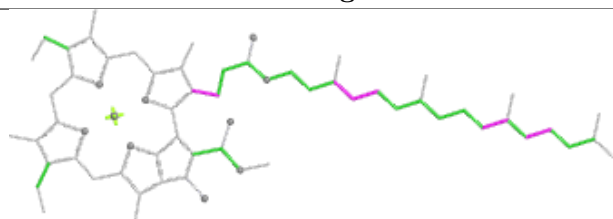
Ligand CLA B 846



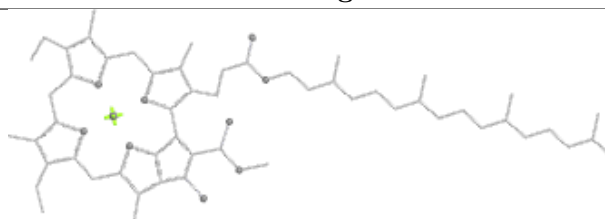
Bond lengths



Bond angles

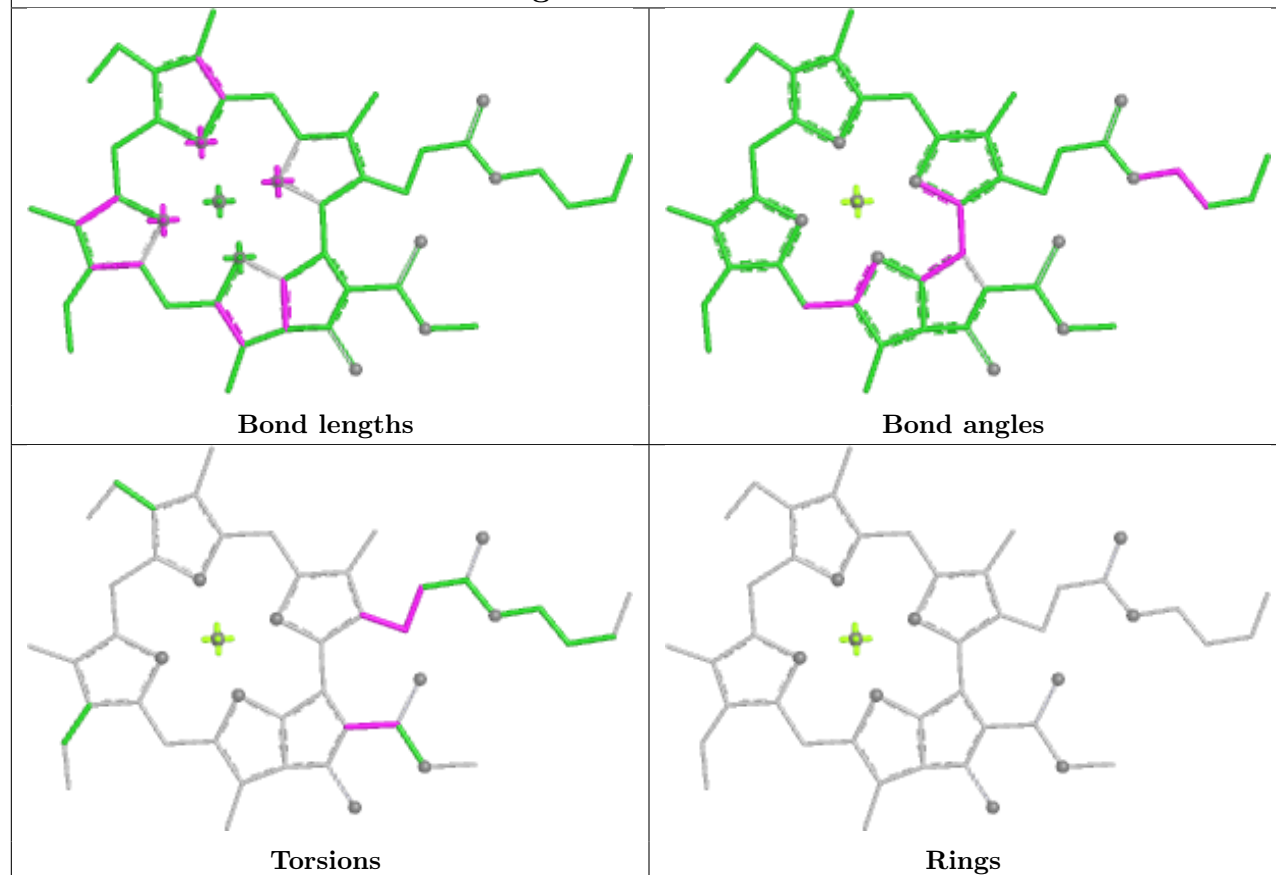


Torsions

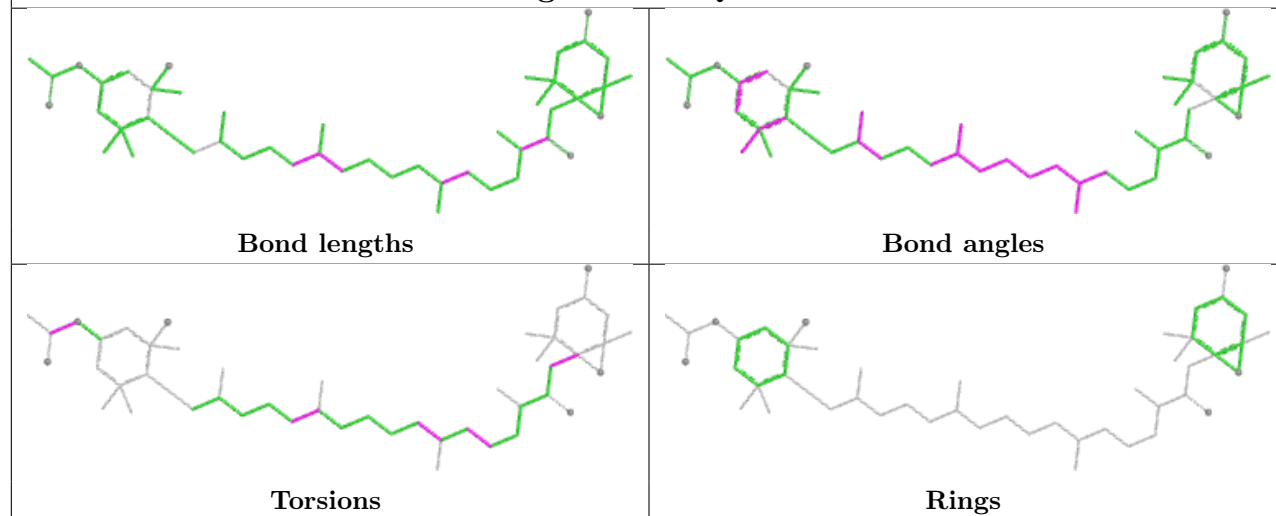


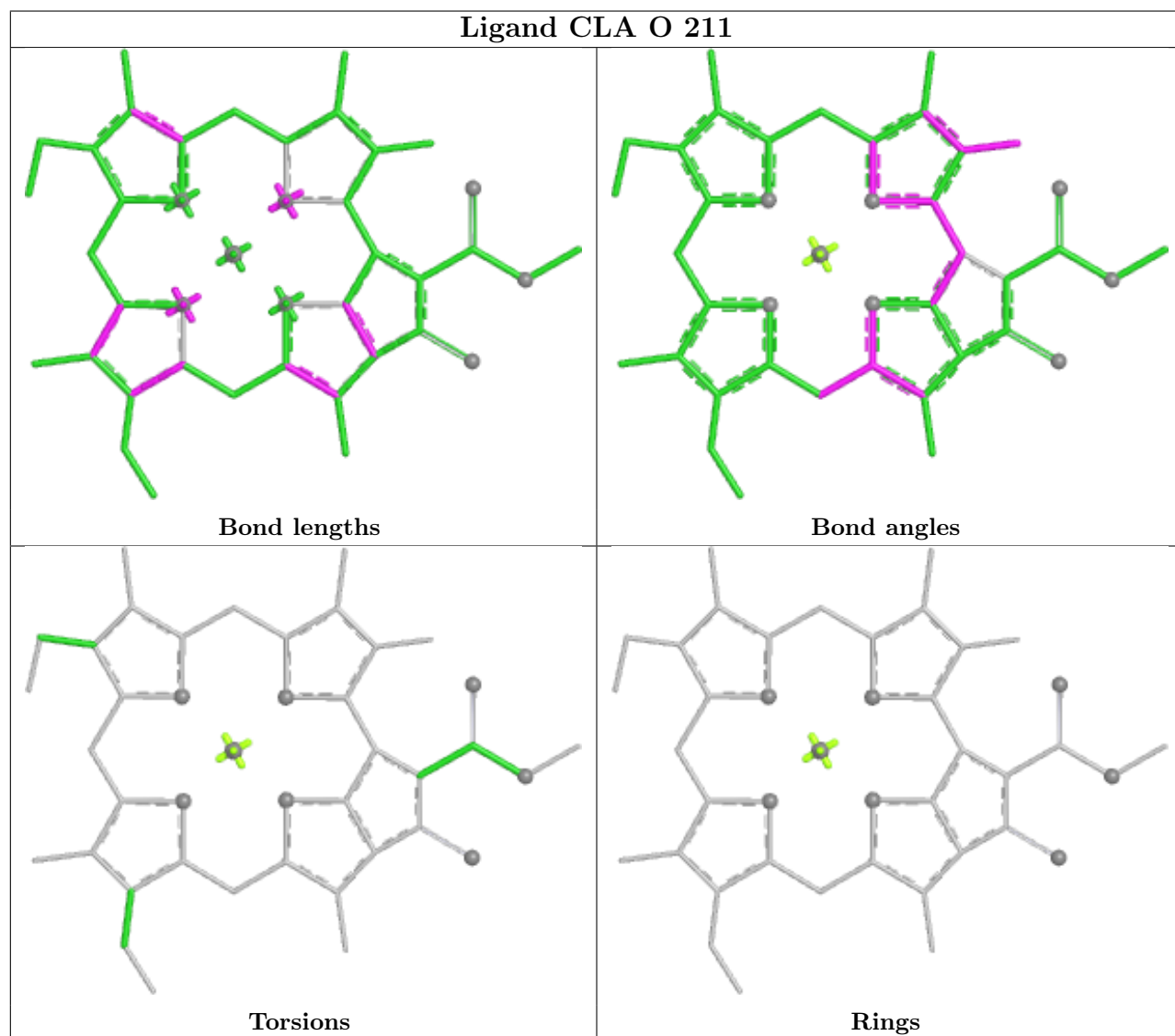
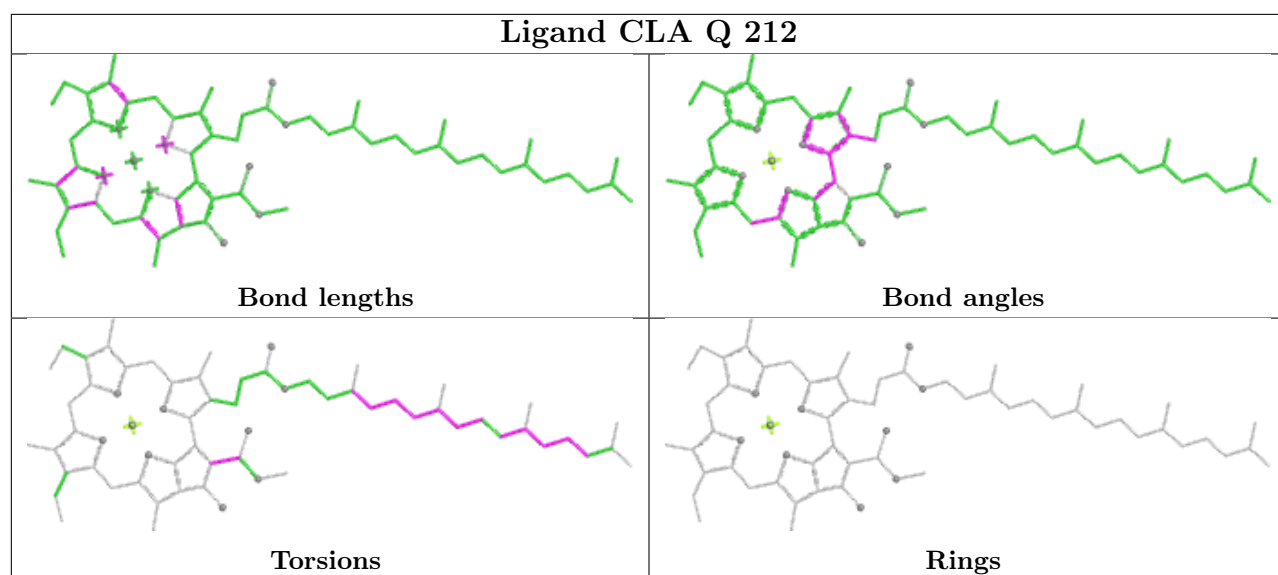
Rings

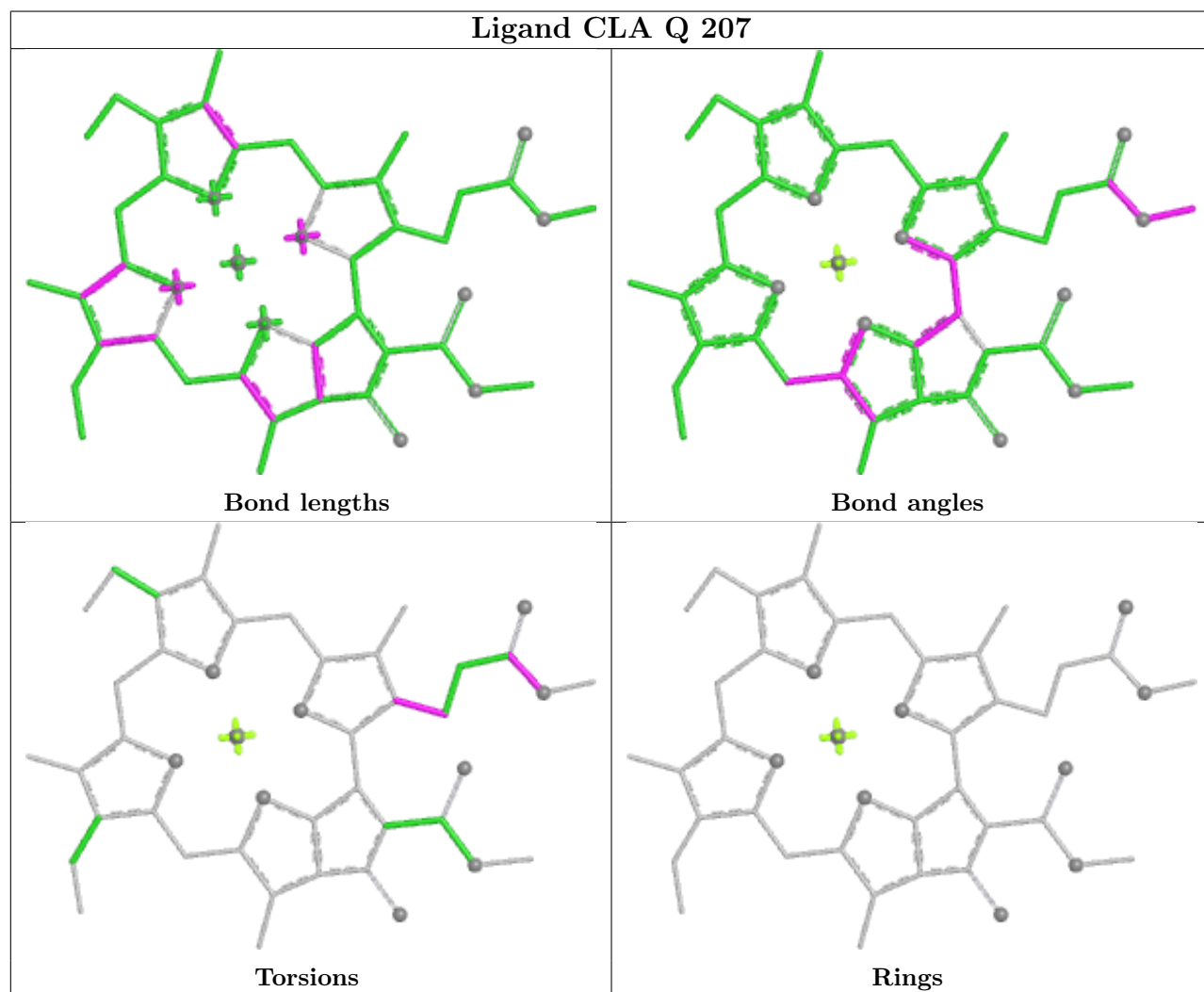
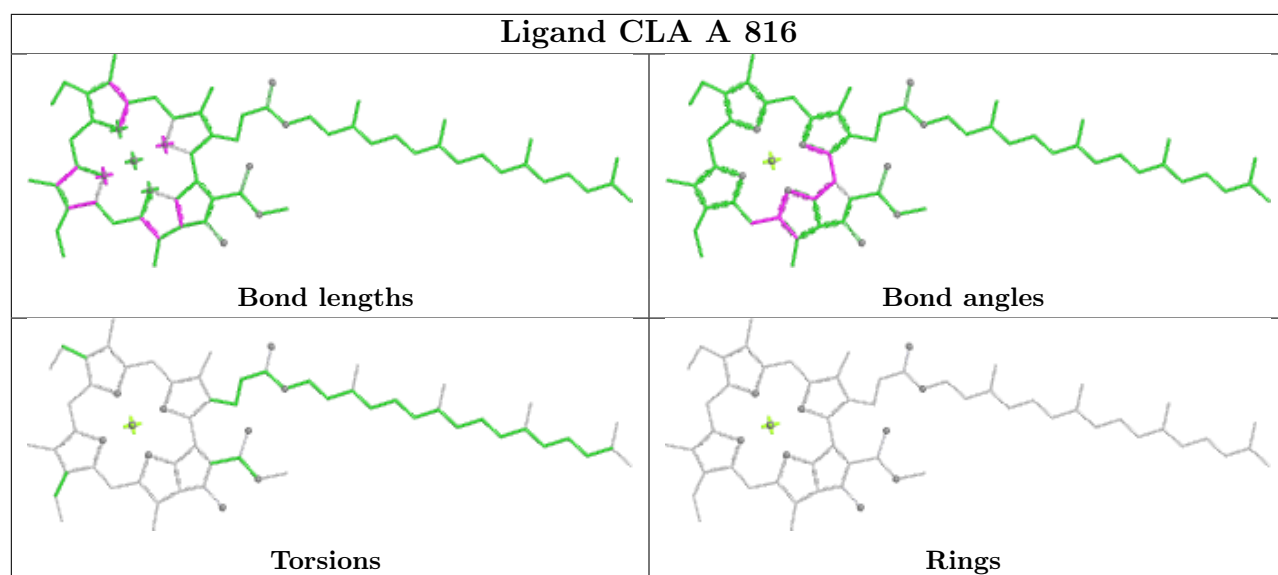
Ligand CLA B 813

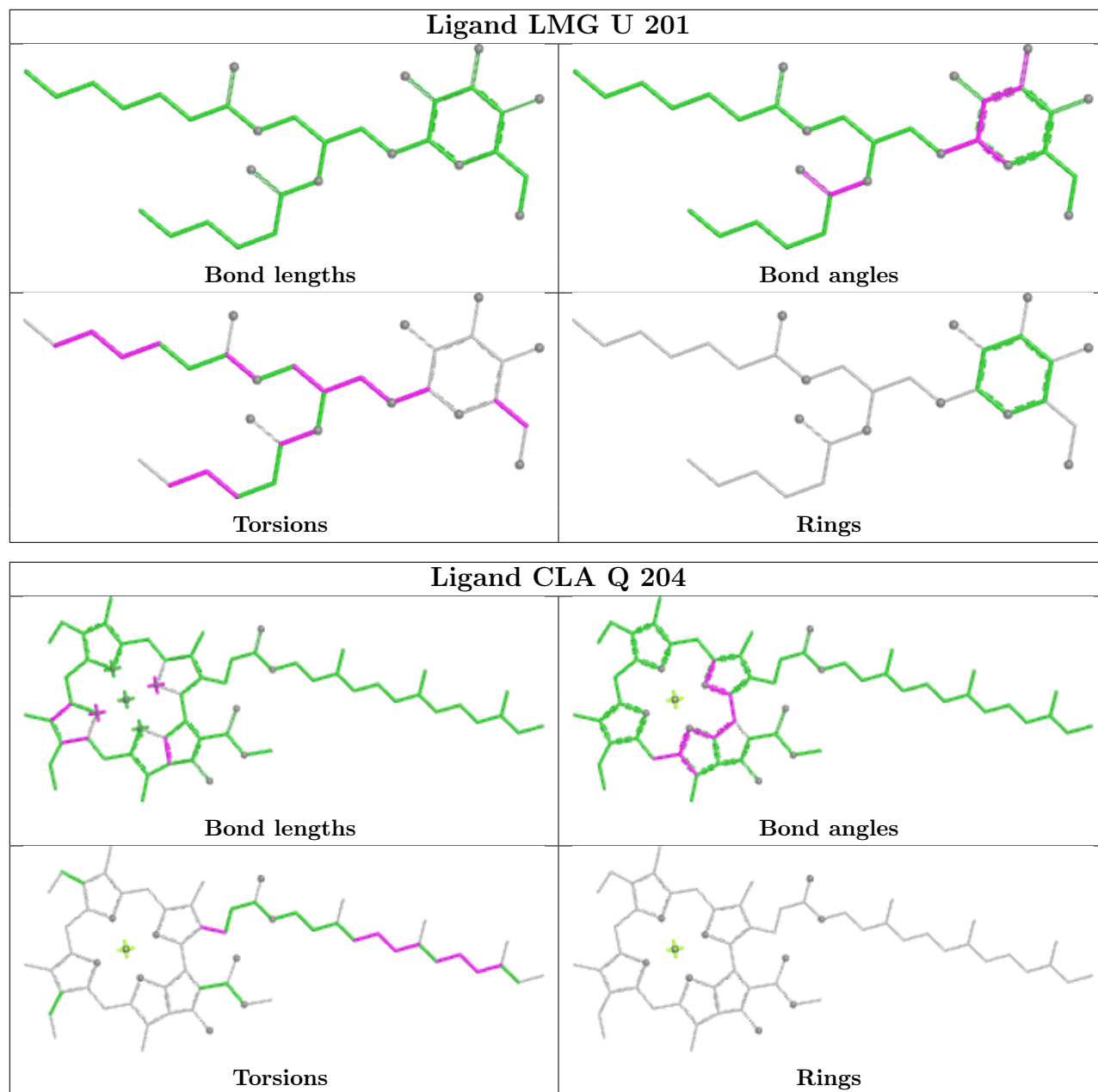


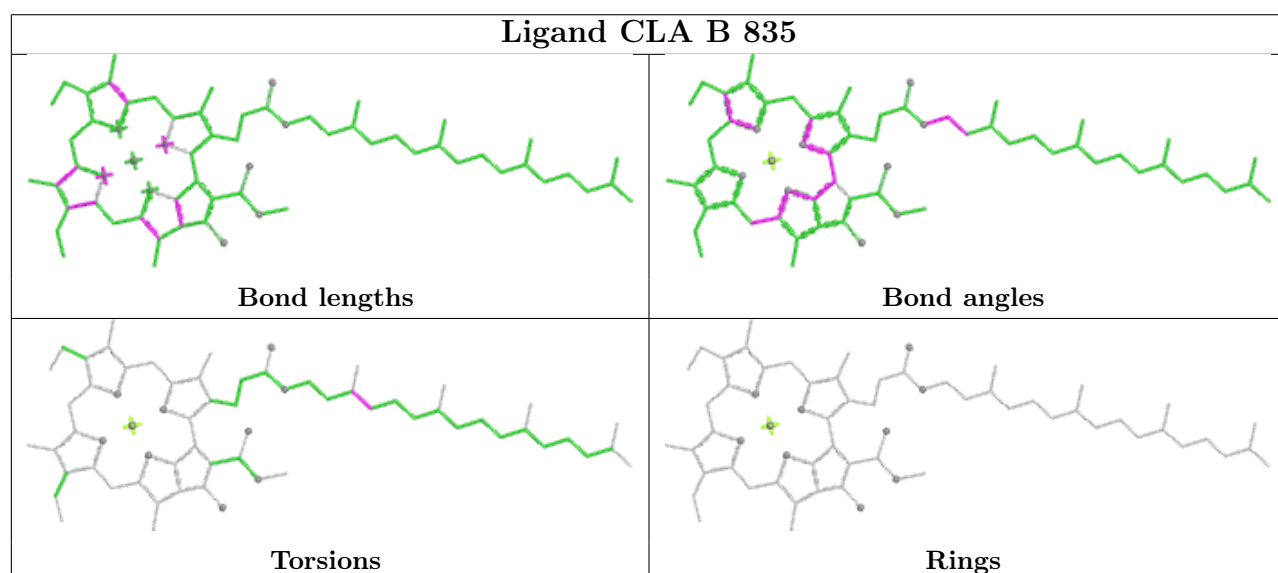
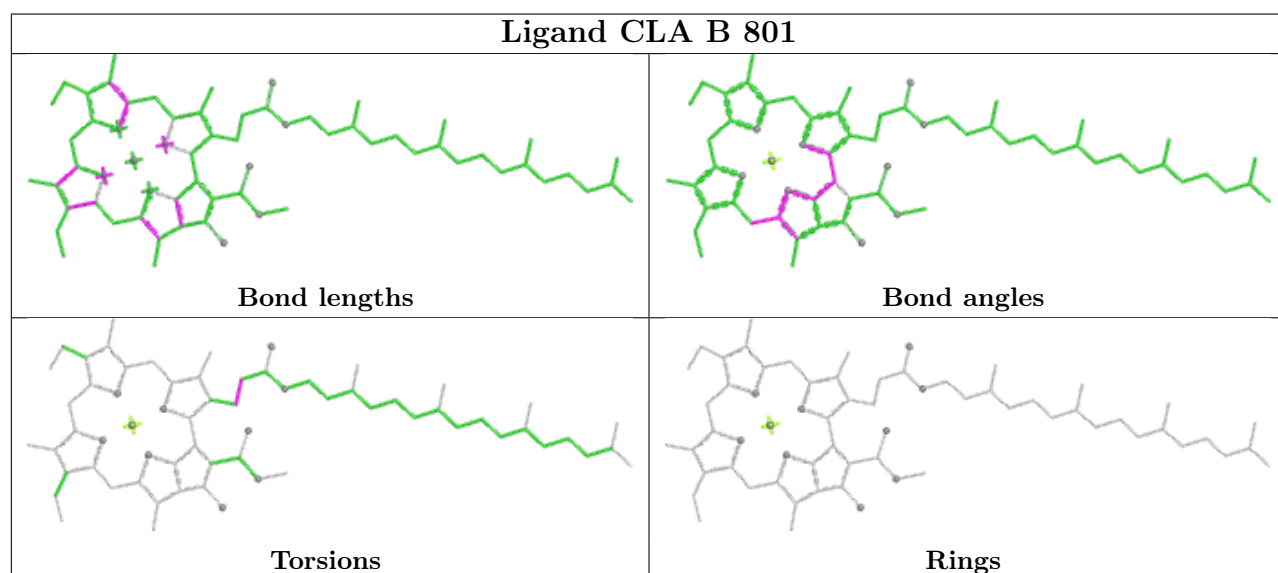
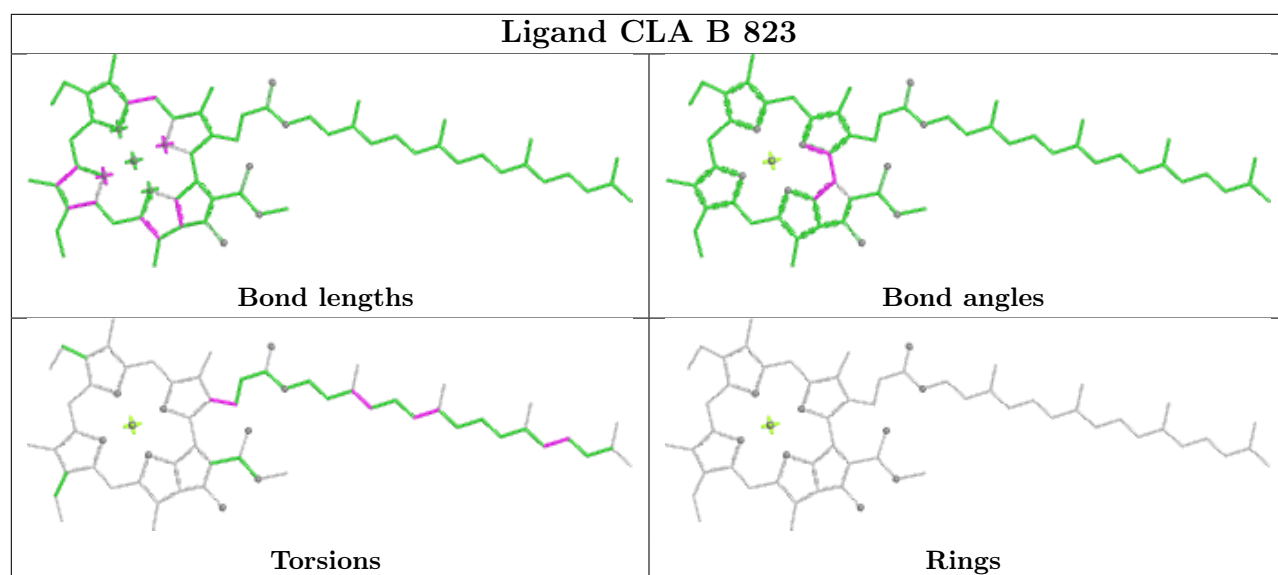
Ligand A86 Q 214

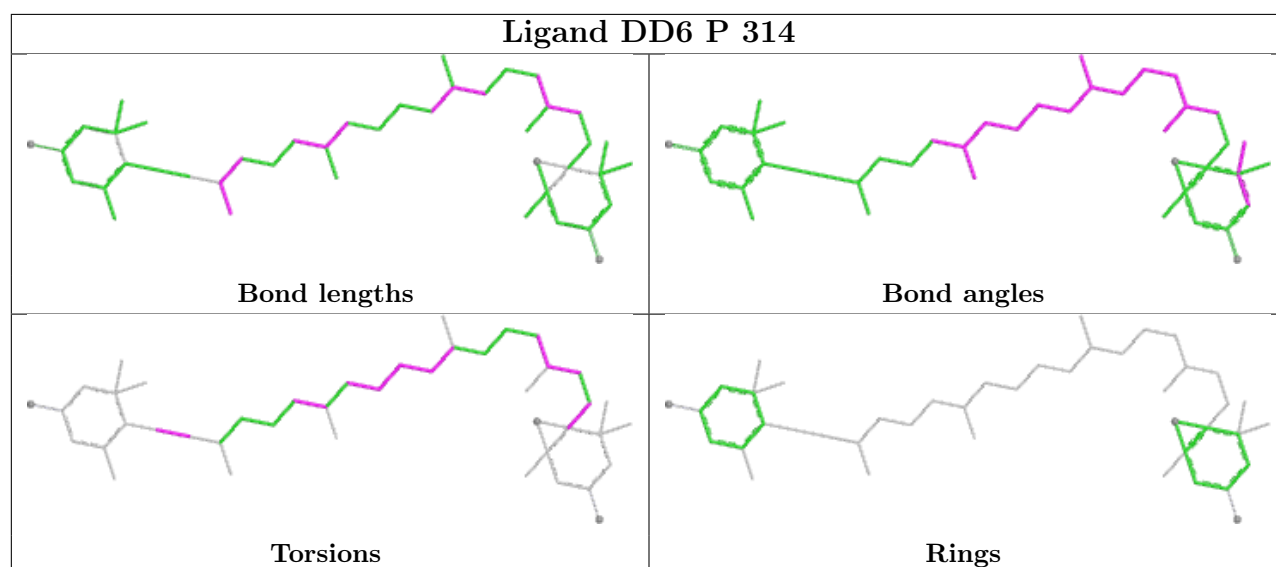
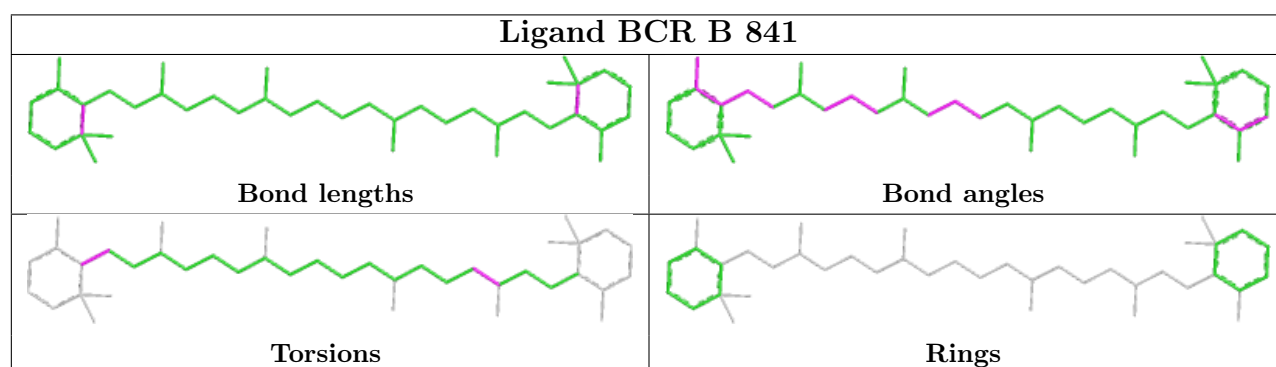
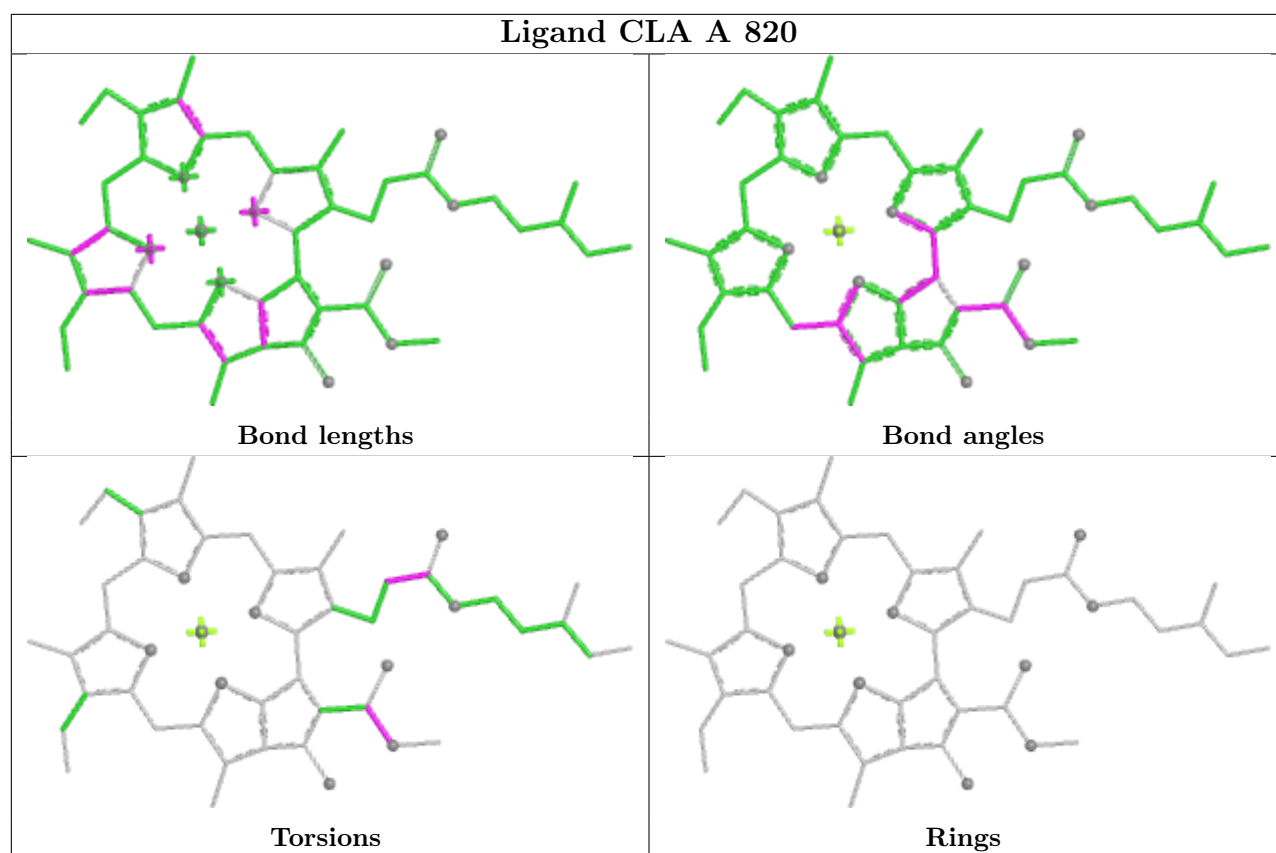


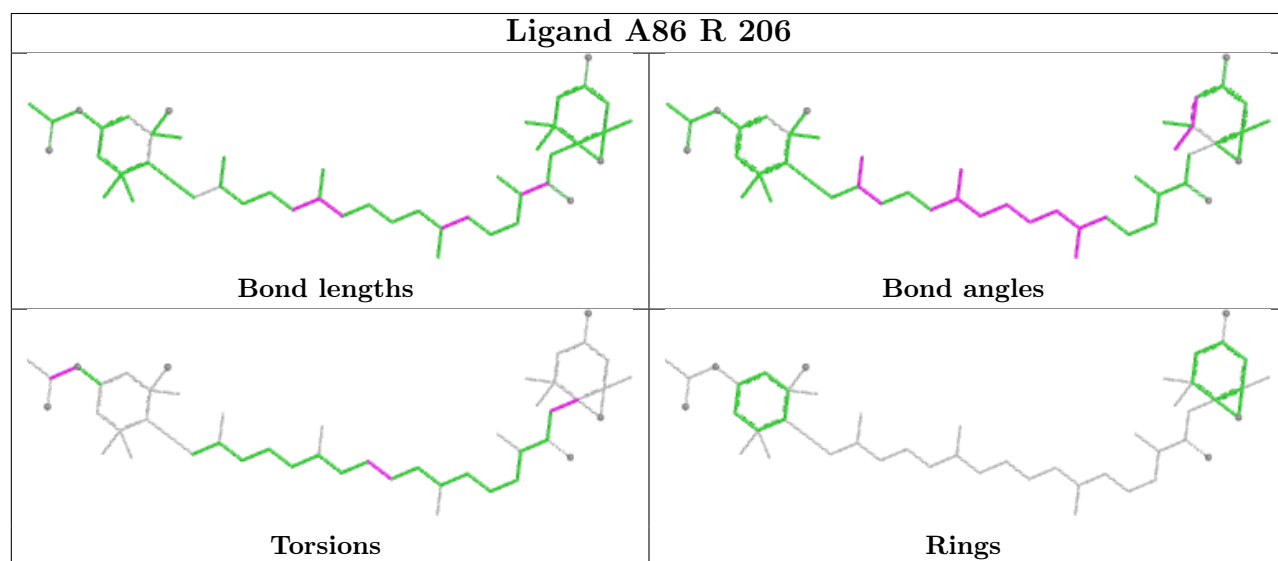
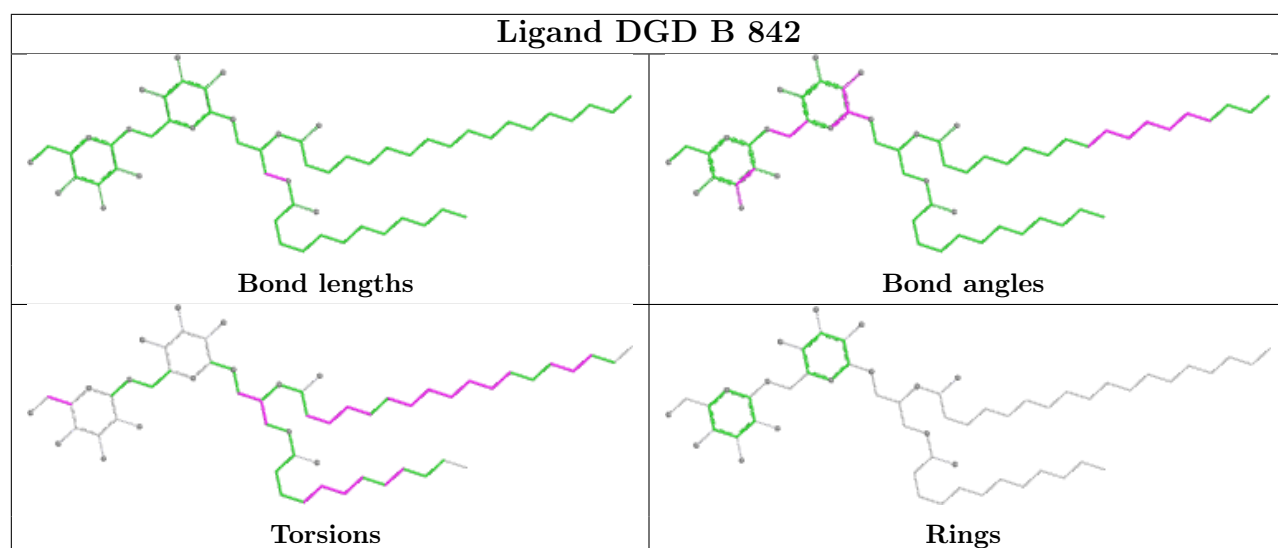




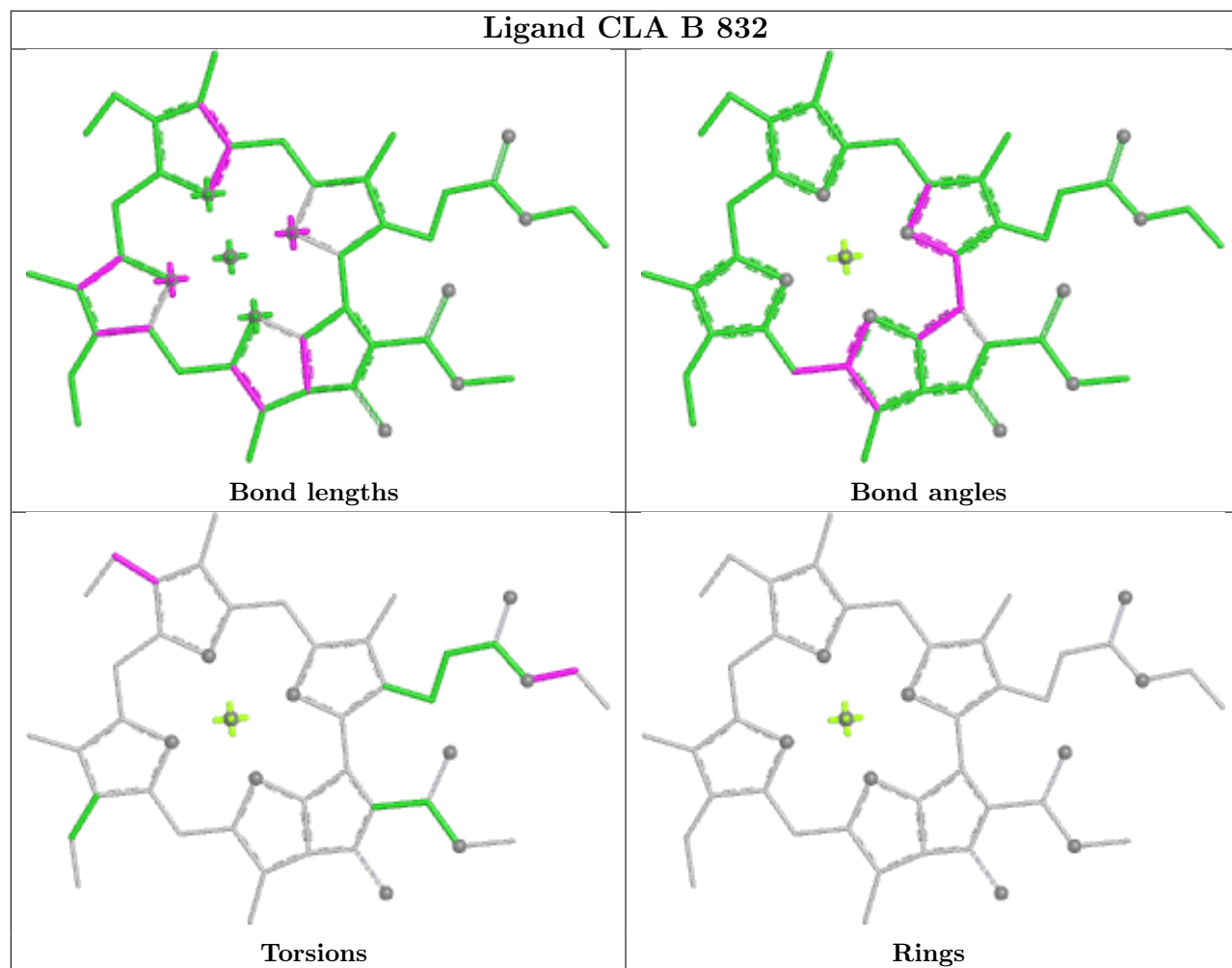


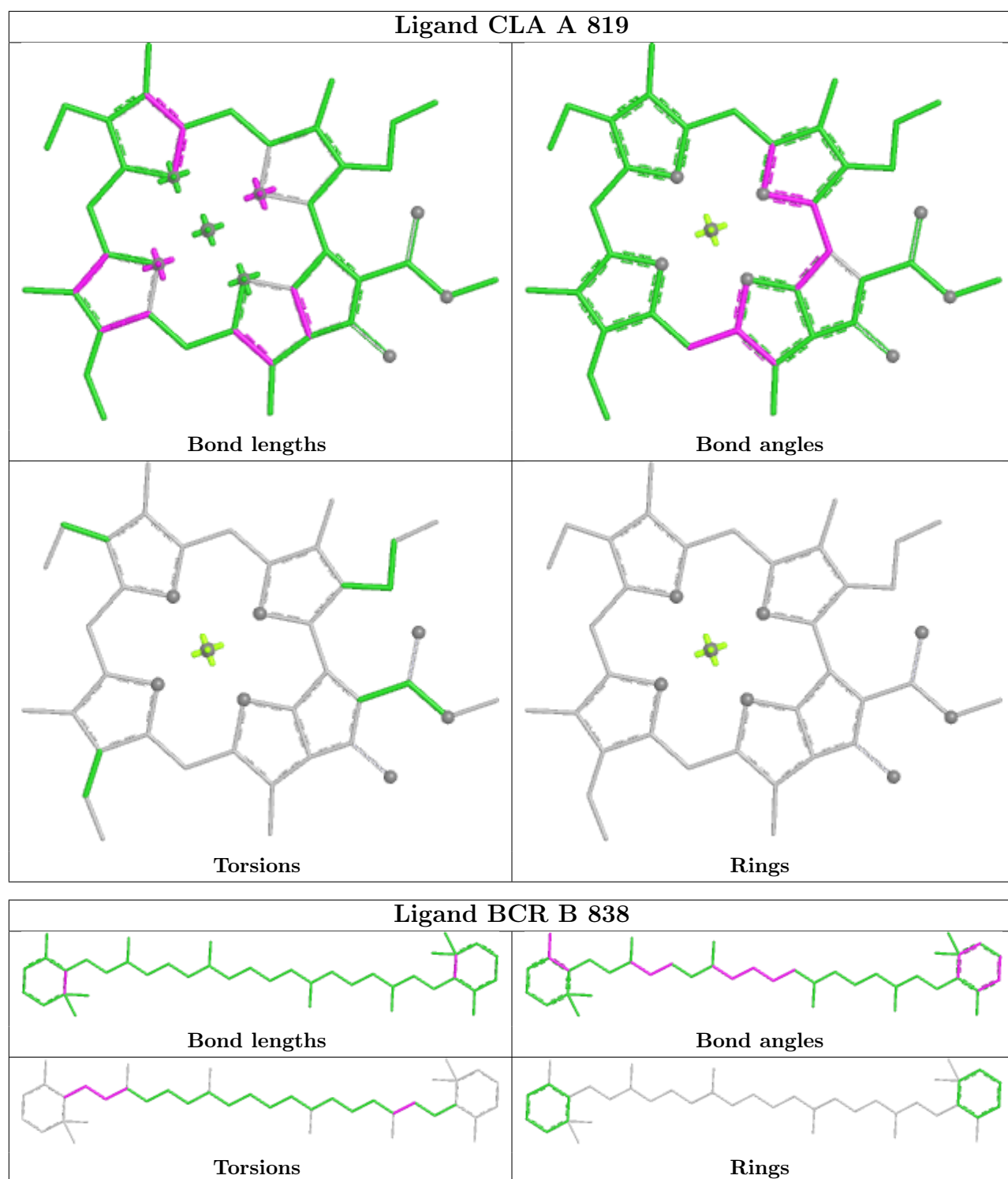


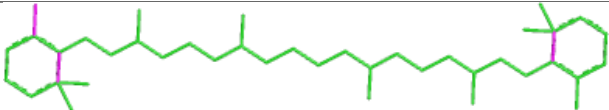
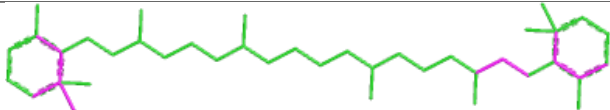
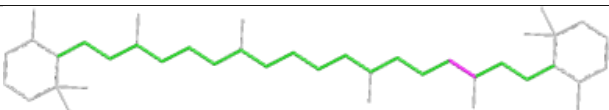
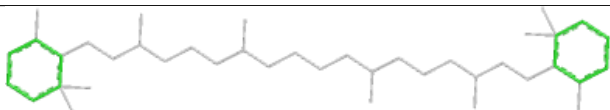




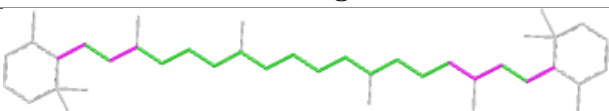
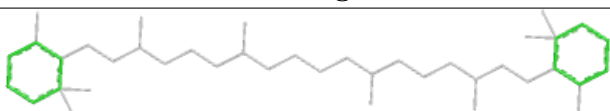


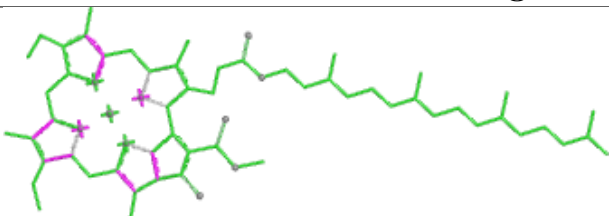
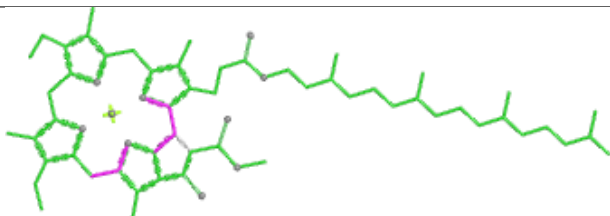
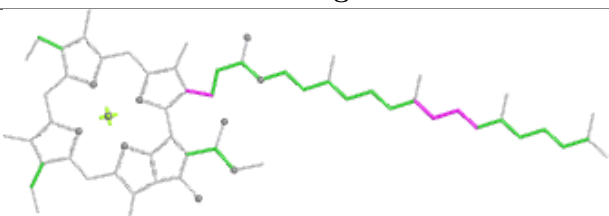
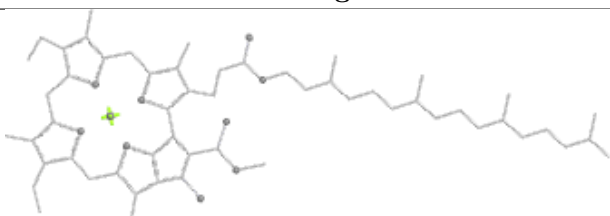
Ligand CLA B 832

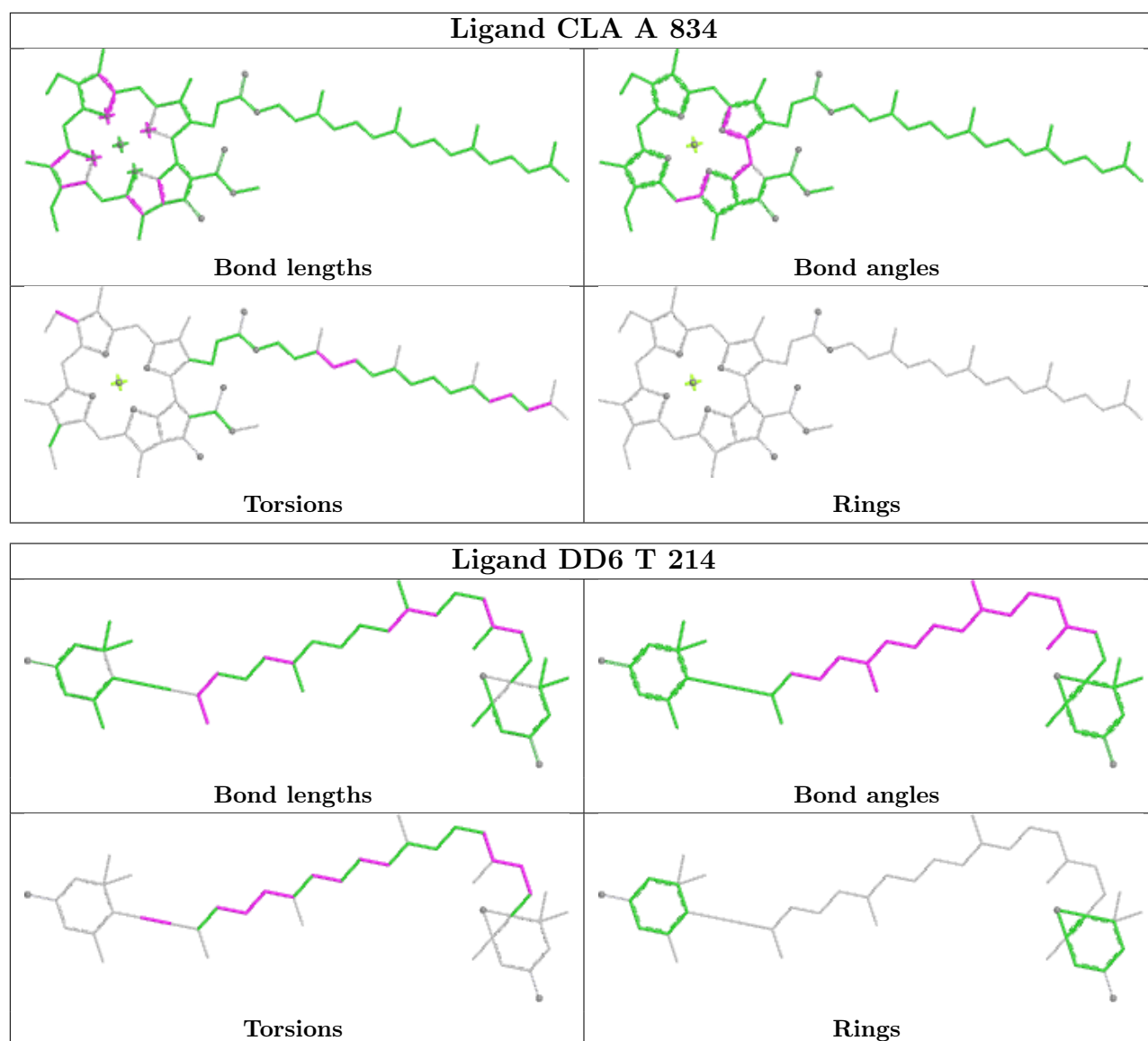




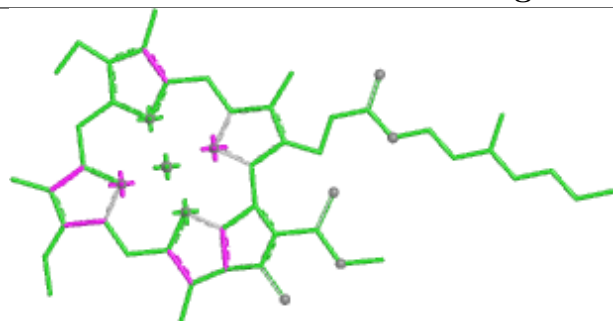
Ligand BCR B 839	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand BCR I 102	
	
Bond lengths	Bond angles
	
Torsions	Rings

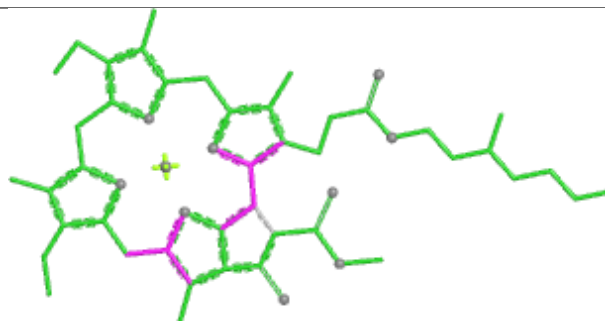
Ligand CLA B 843	
	
Bond lengths	Bond angles
	
Torsions	Rings



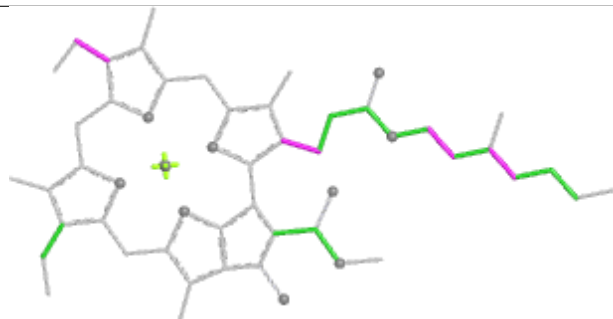
Ligand CLA B 819



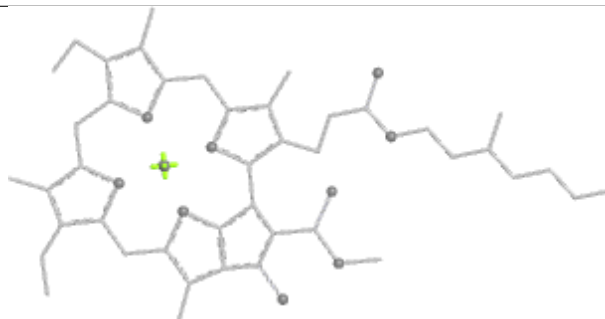
Bond lengths



Bond angles

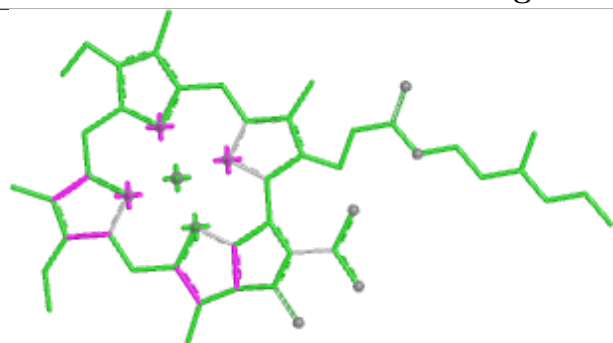


Torsions

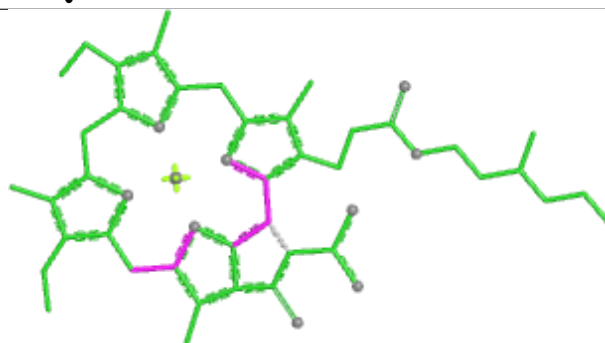


Rings

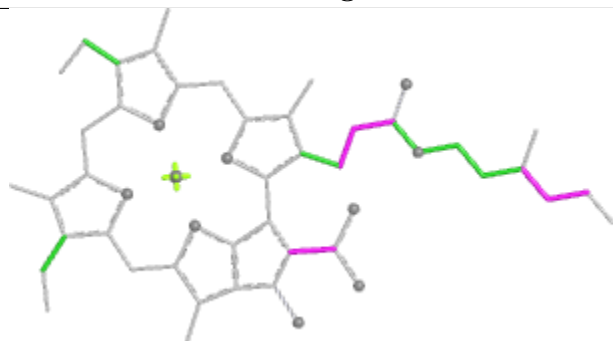
Ligand CLA Q 206



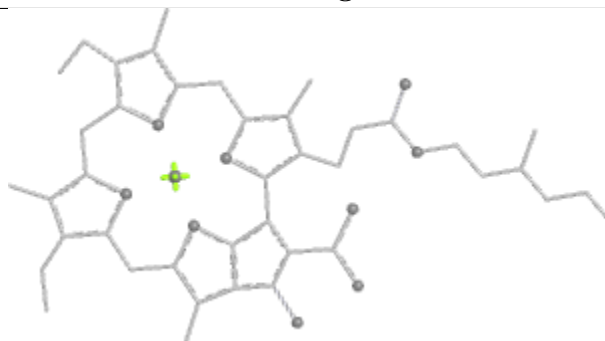
Bond lengths



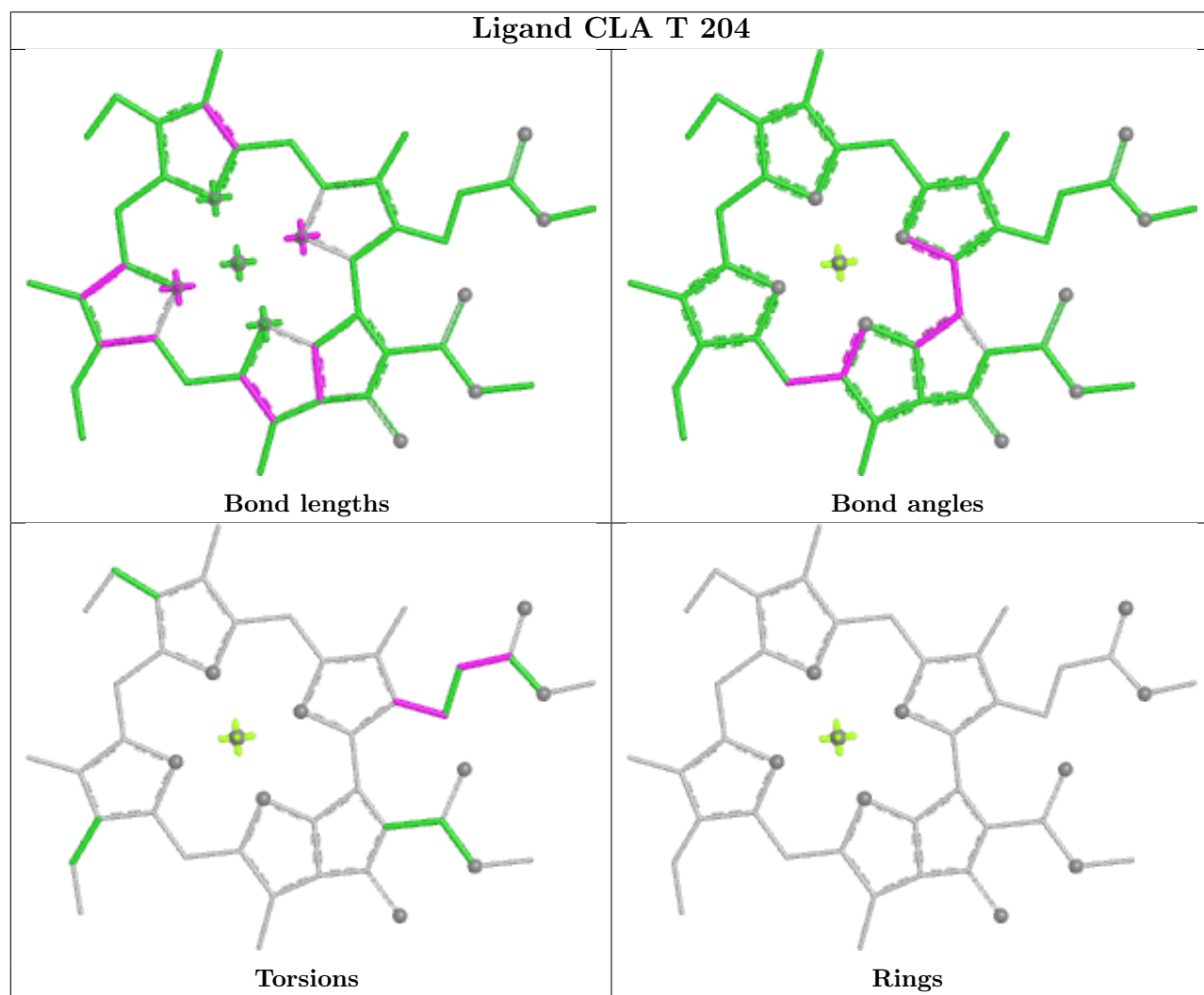
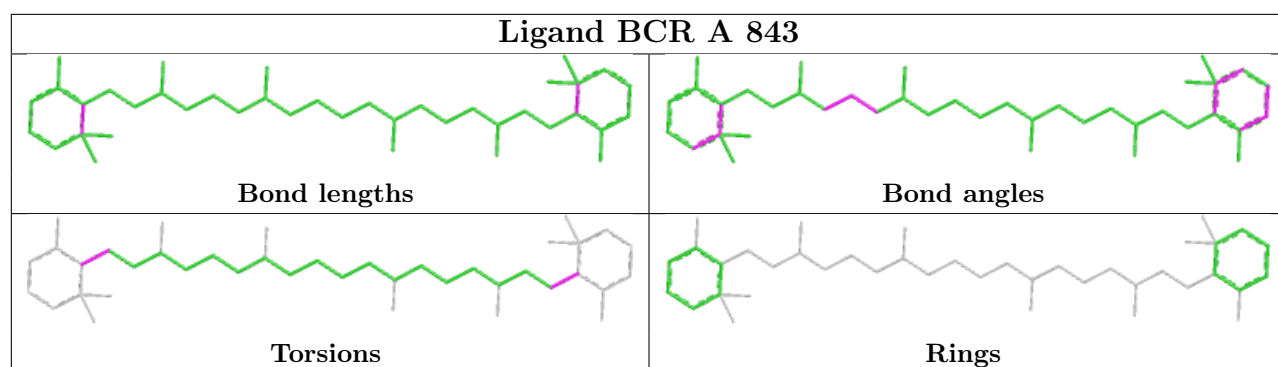
Bond angles



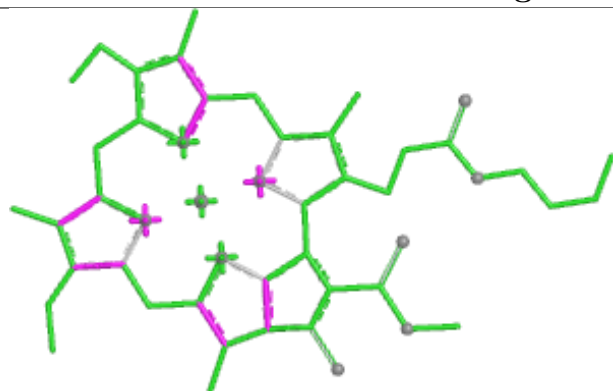
Torsions



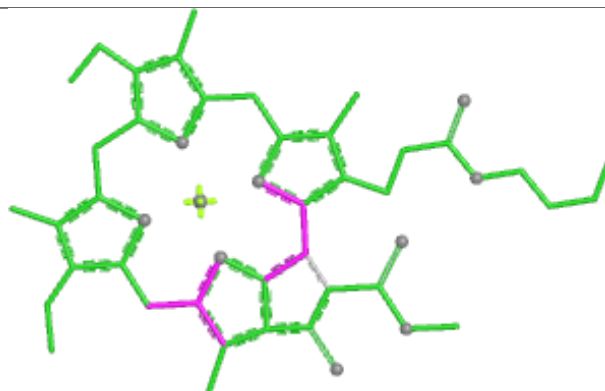
Rings



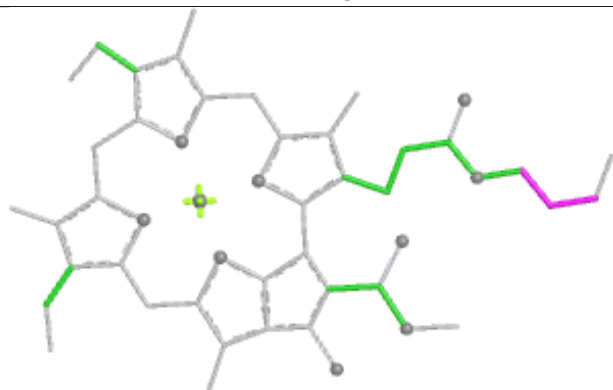
Ligand CLA B 827



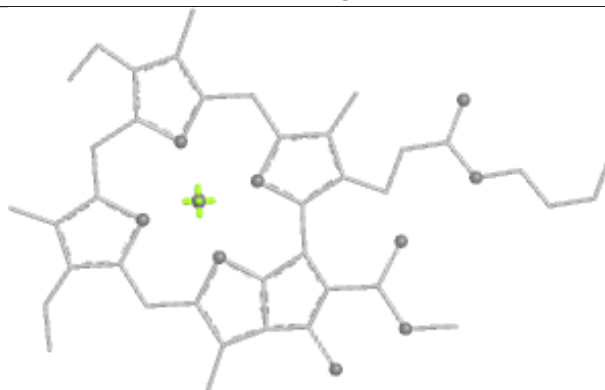
Bond lengths



Bond angles

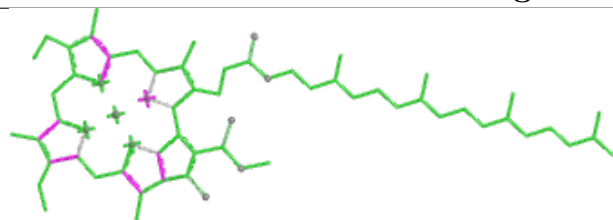


Torsions

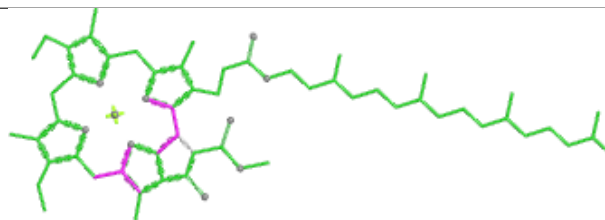


Rings

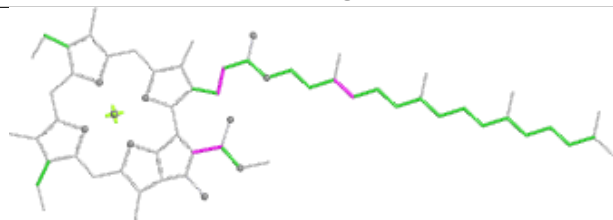
Ligand CLA A 806



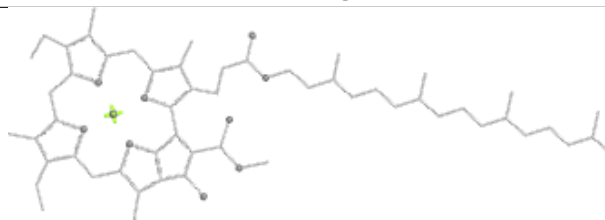
Bond lengths



Bond angles

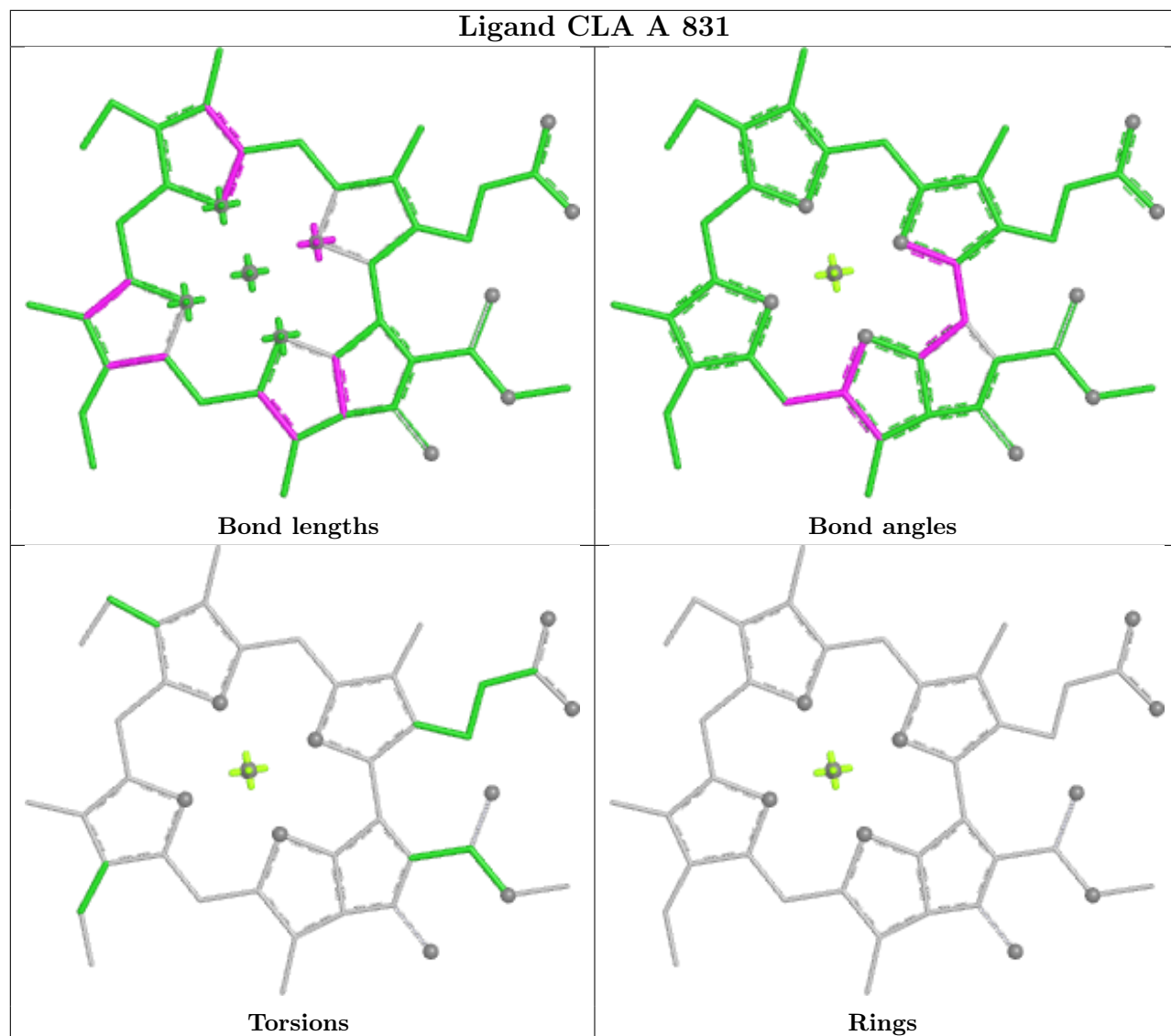


Torsions

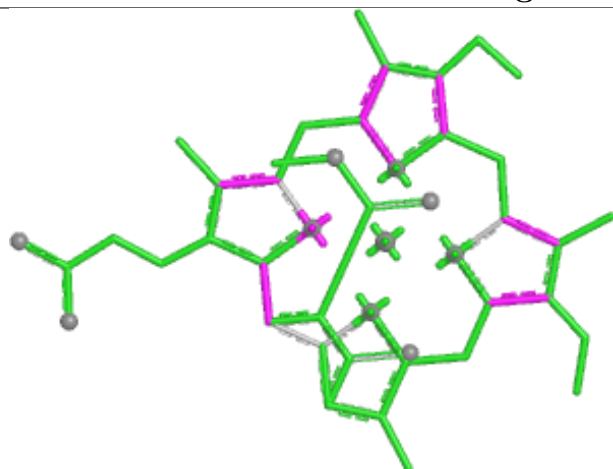


Rings

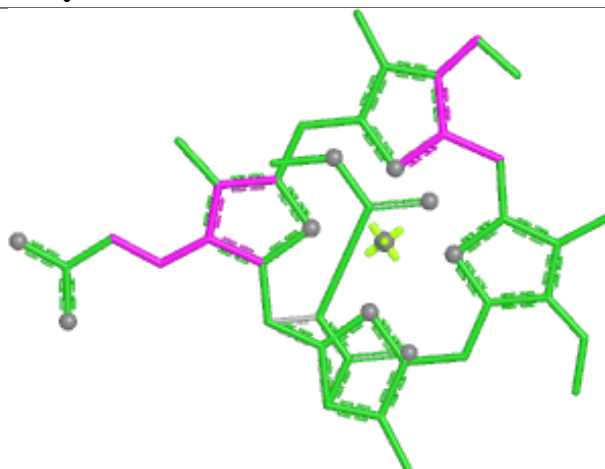
Ligand CLA A 831



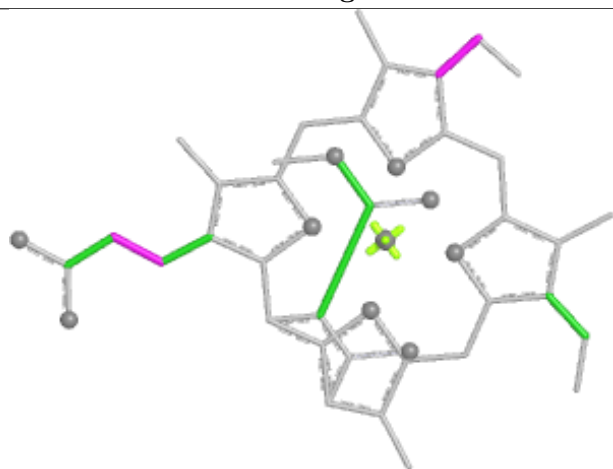
Ligand KC1 Q 210



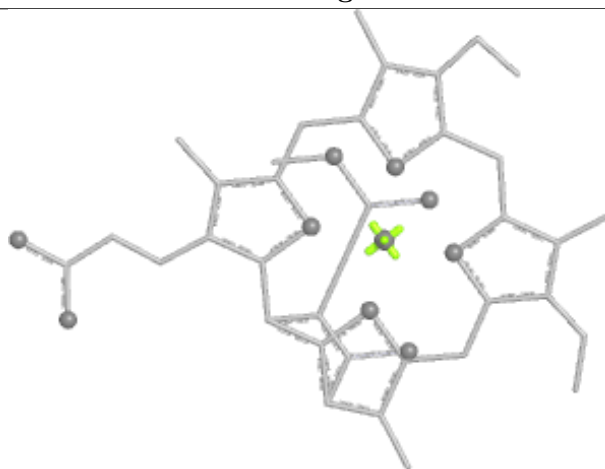
Bond lengths



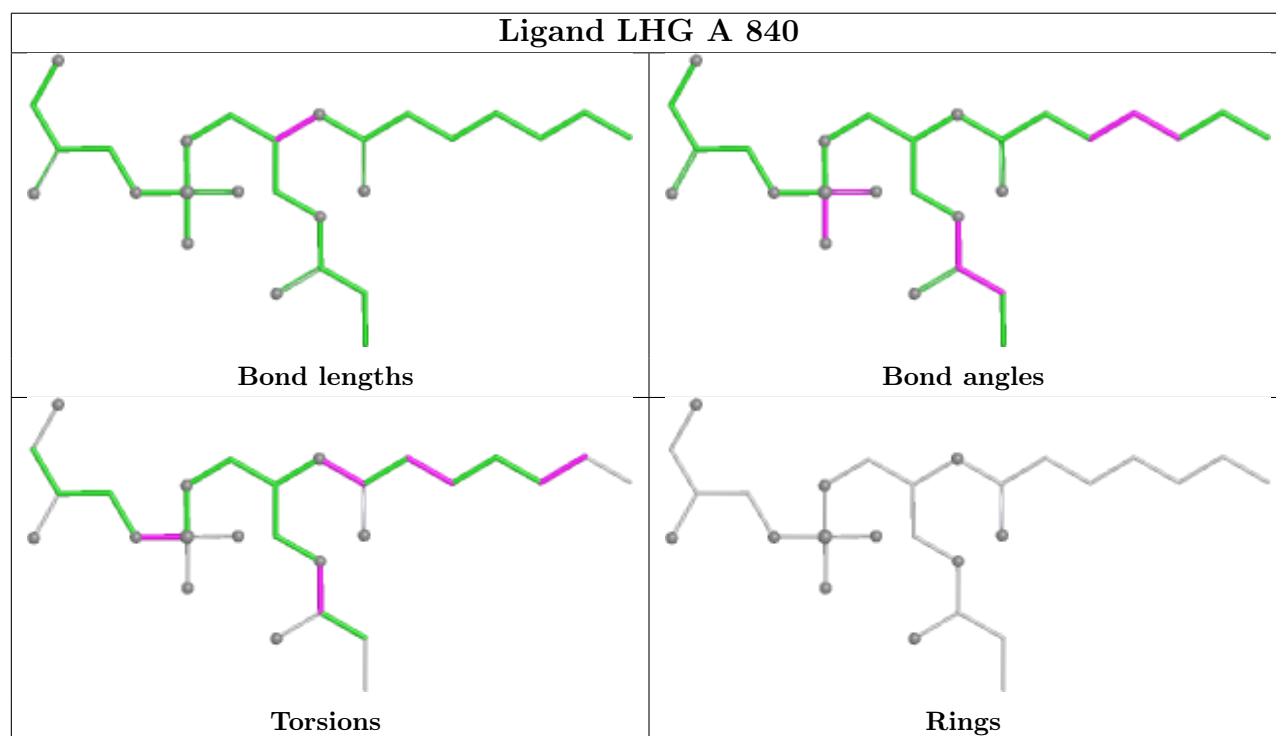
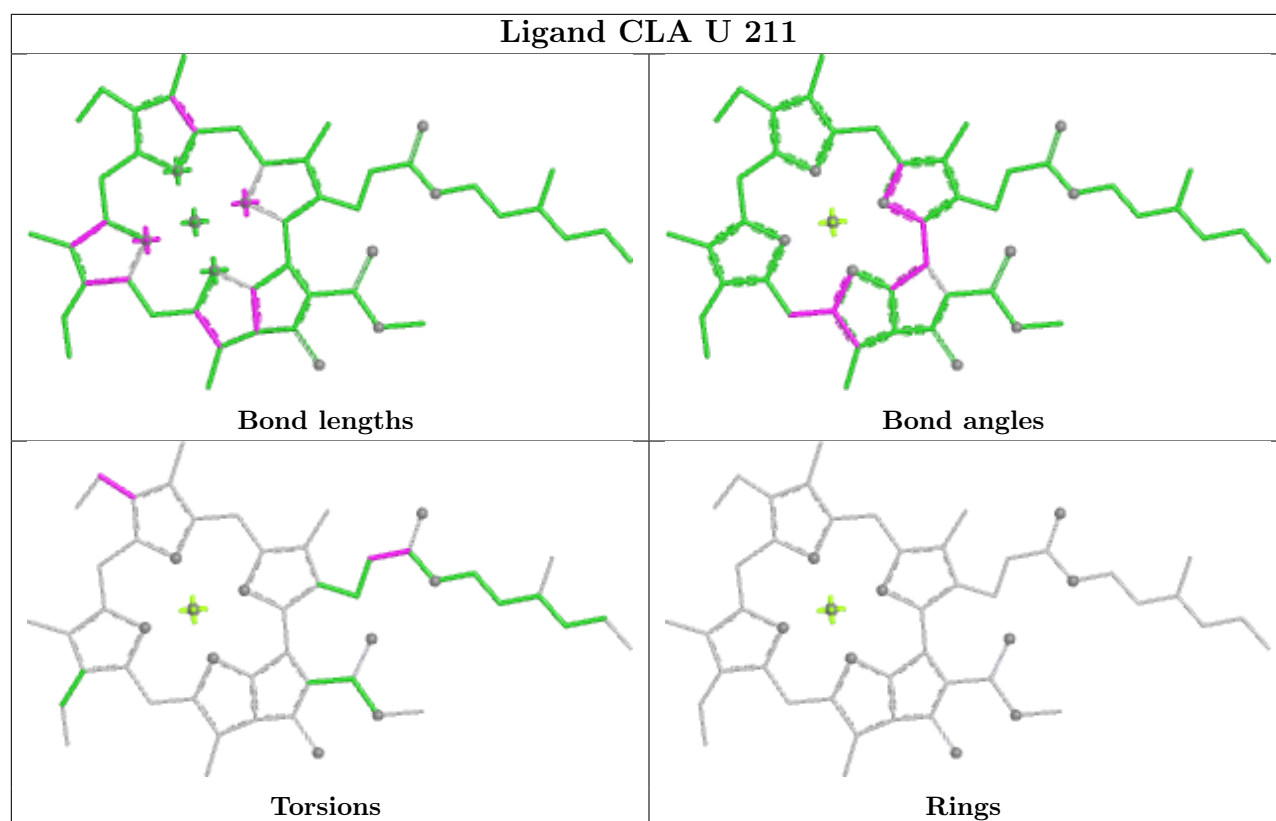
Bond angles

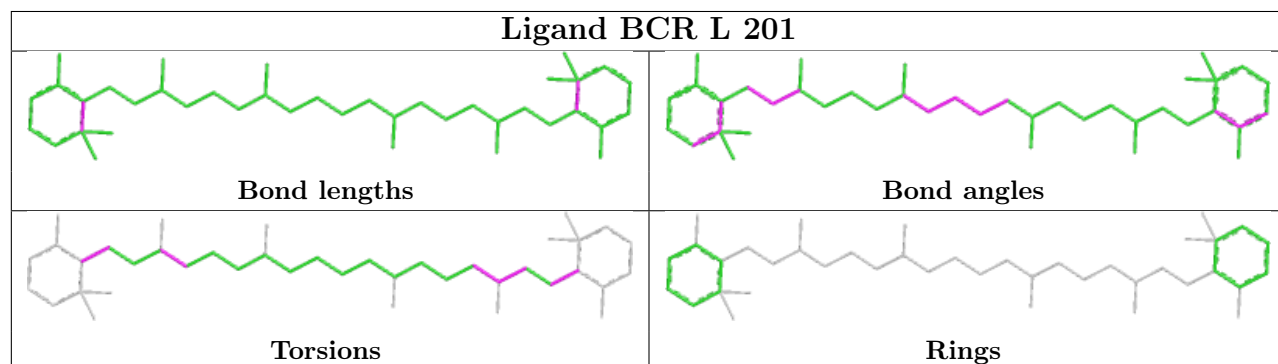
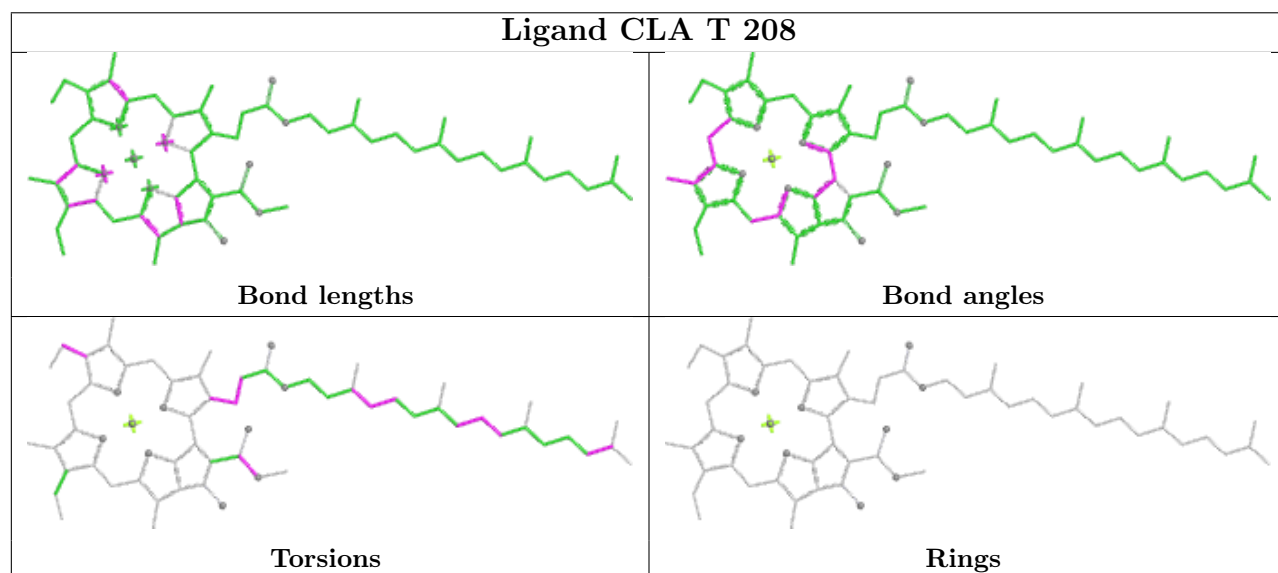
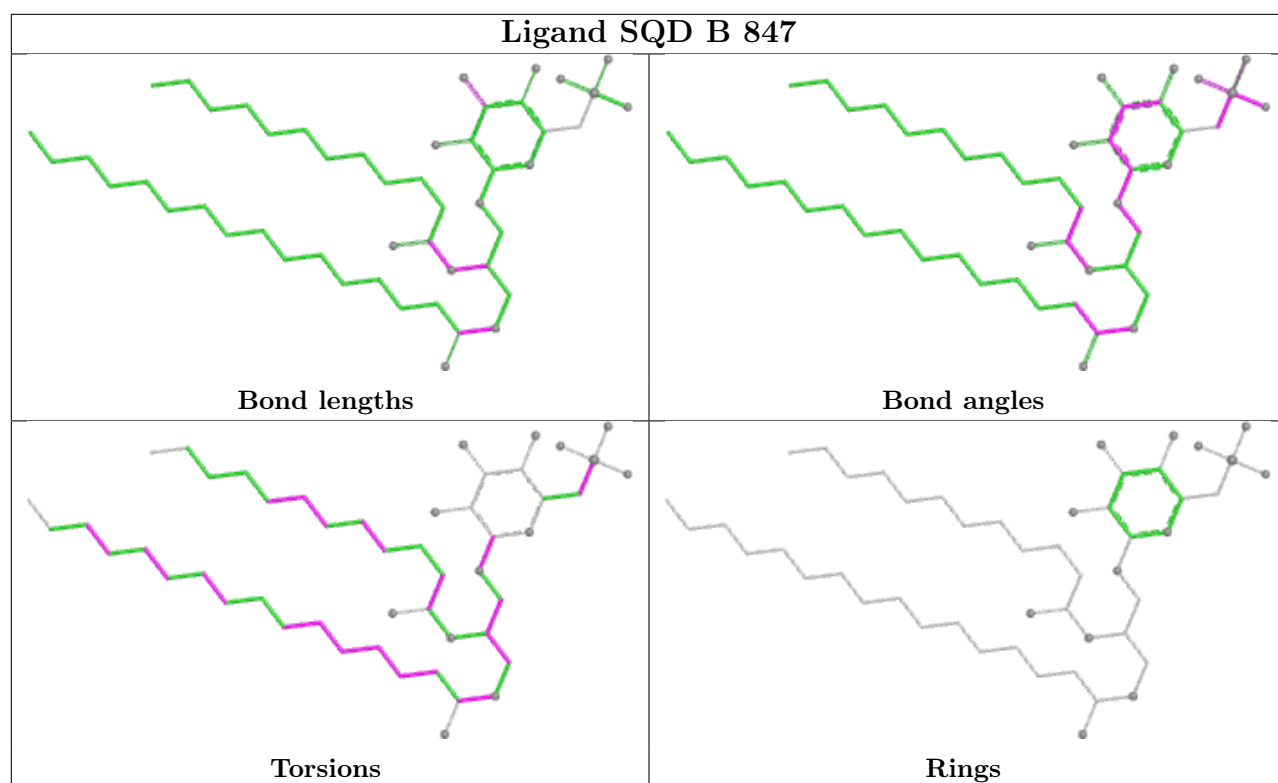


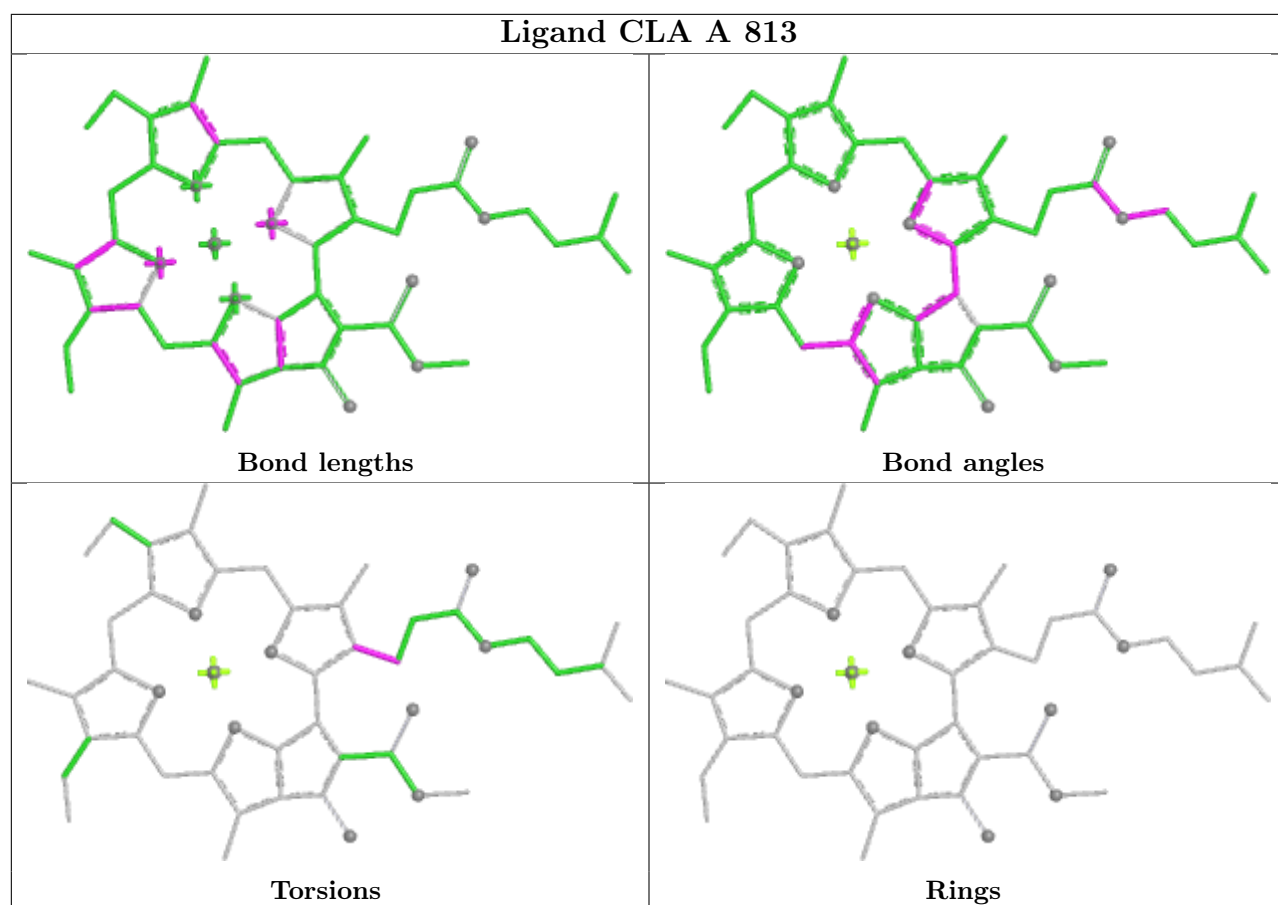
Torsions

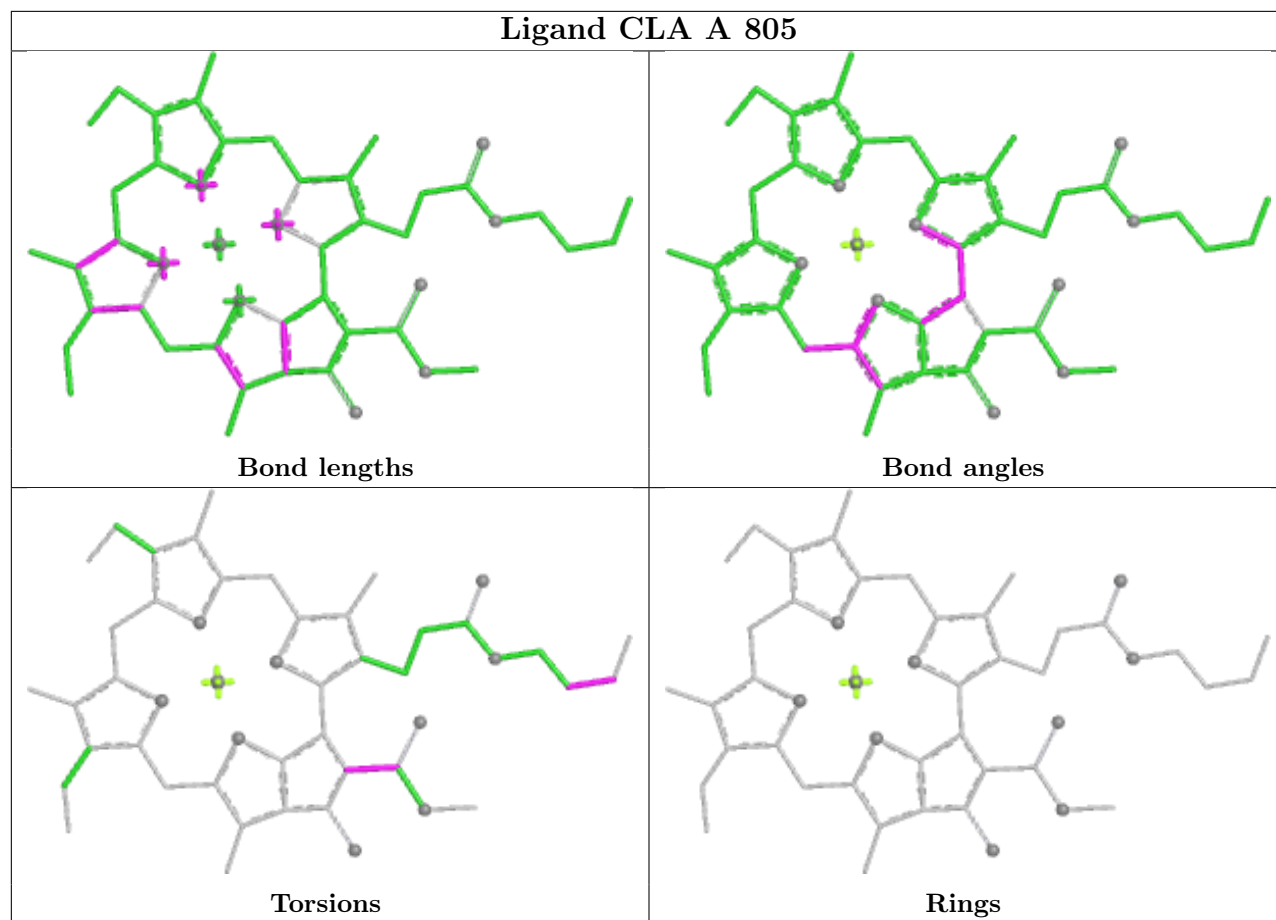


Rings

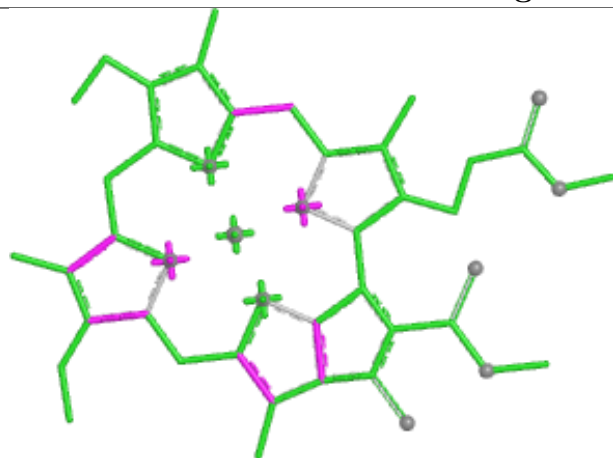




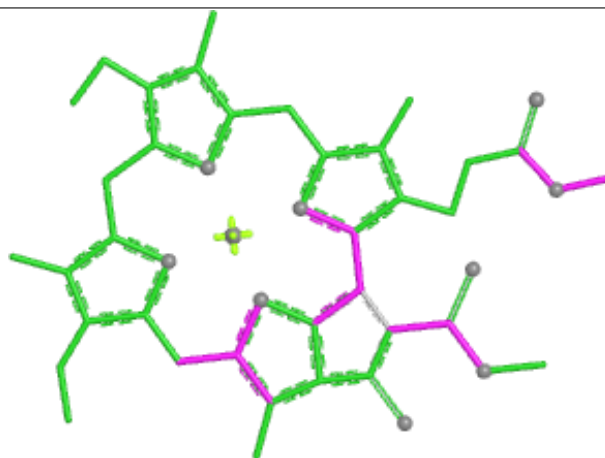




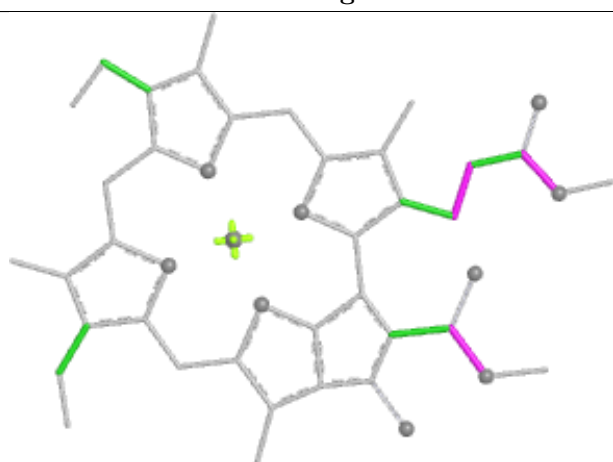
Ligand CLA T 206



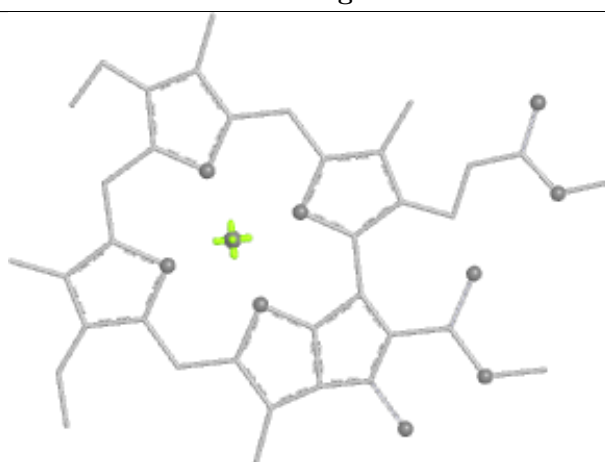
Bond lengths



Bond angles

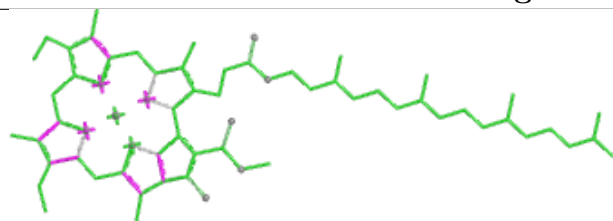


Torsions

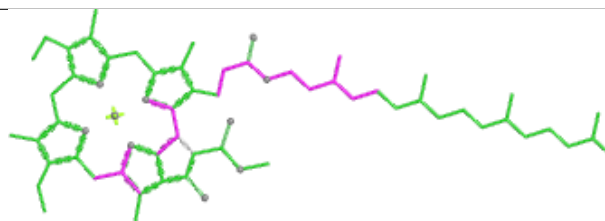


Rings

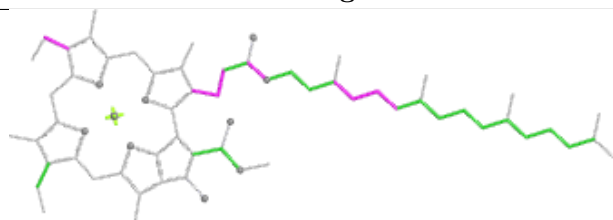
Ligand CLA B 824



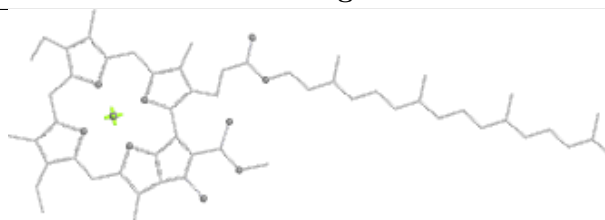
Bond lengths



Bond angles

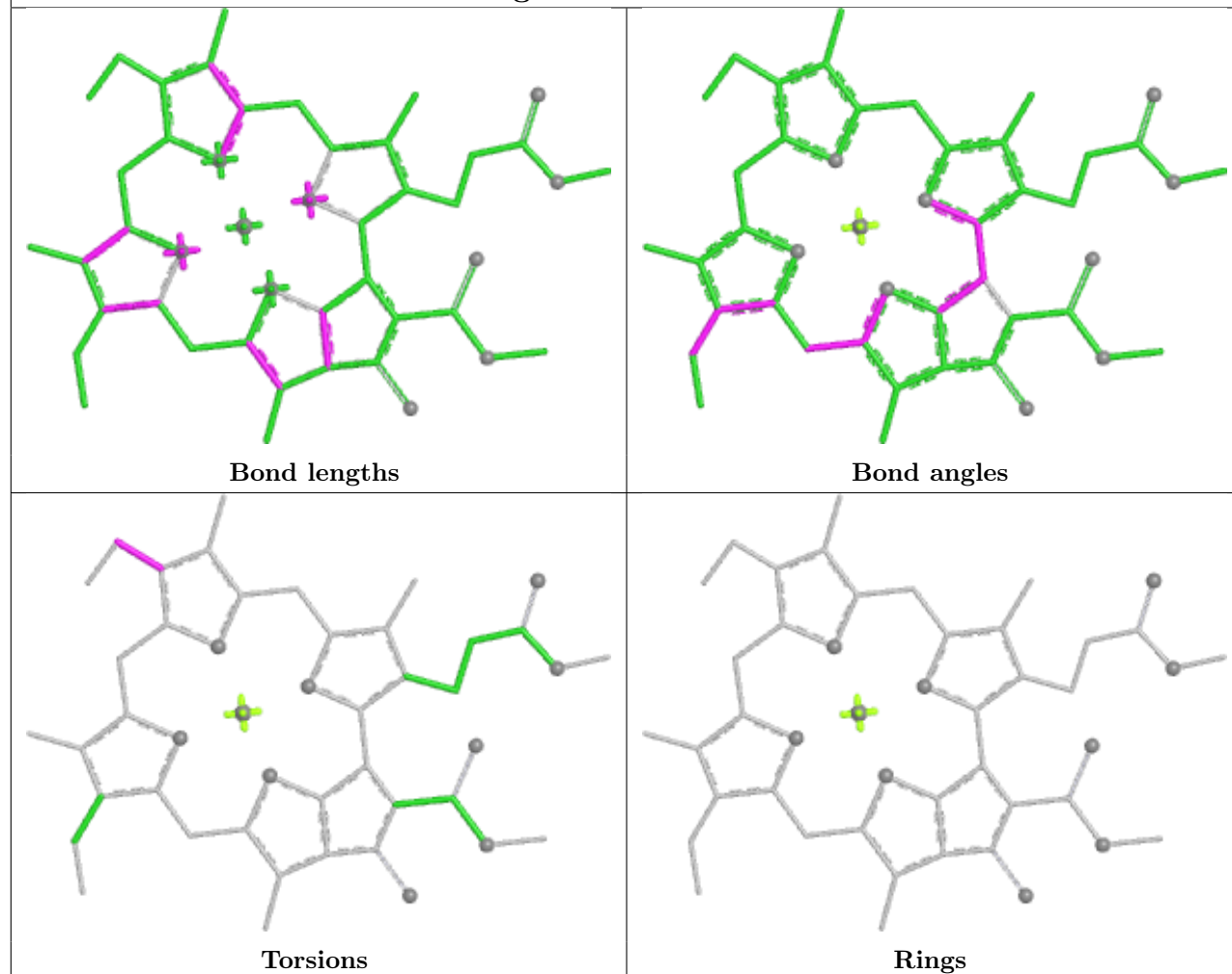


Torsions

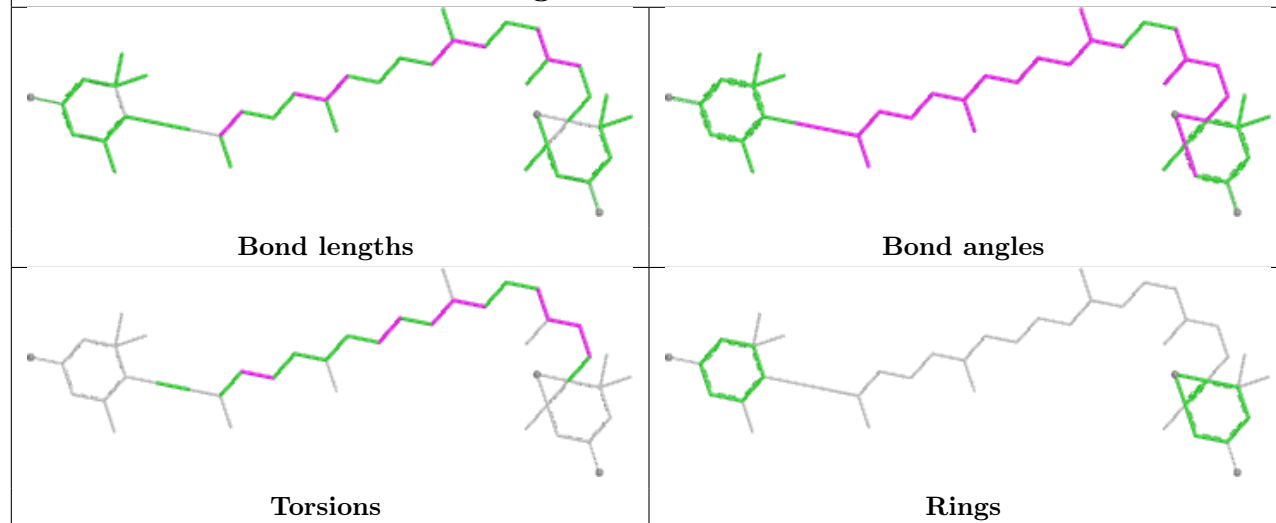


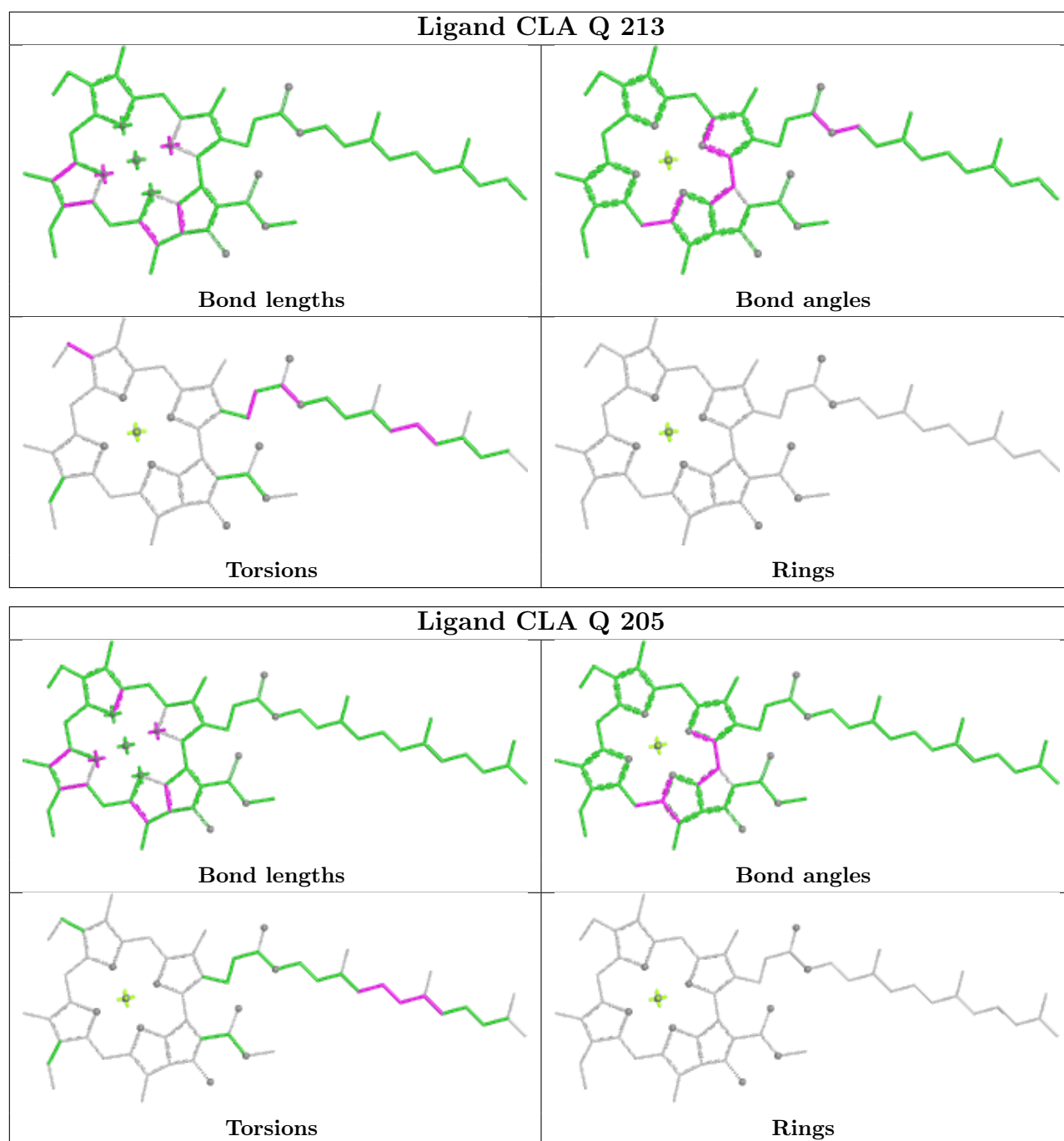
Rings

Ligand CLA O 207

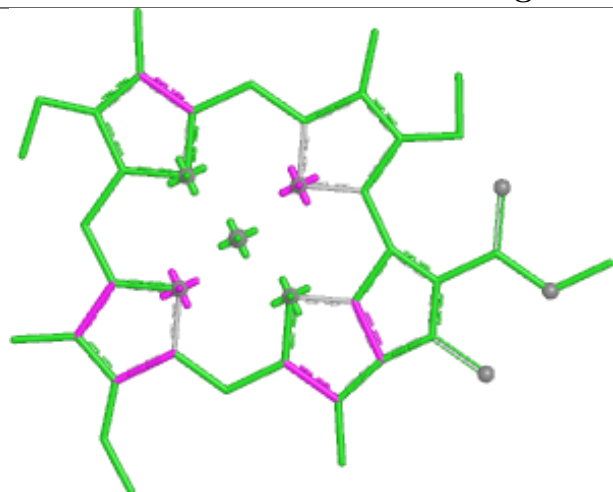


Ligand DD6 U 212

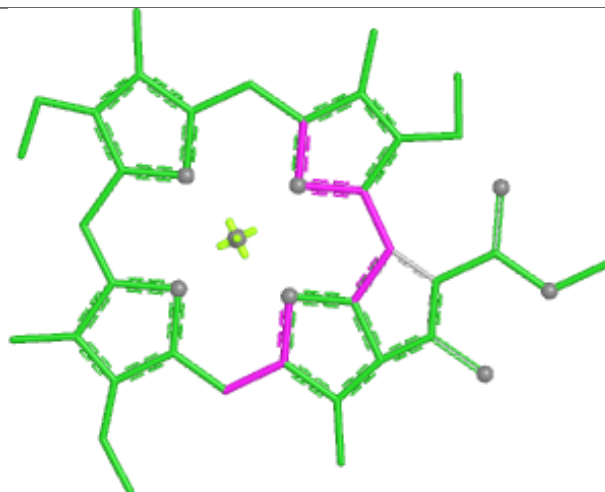




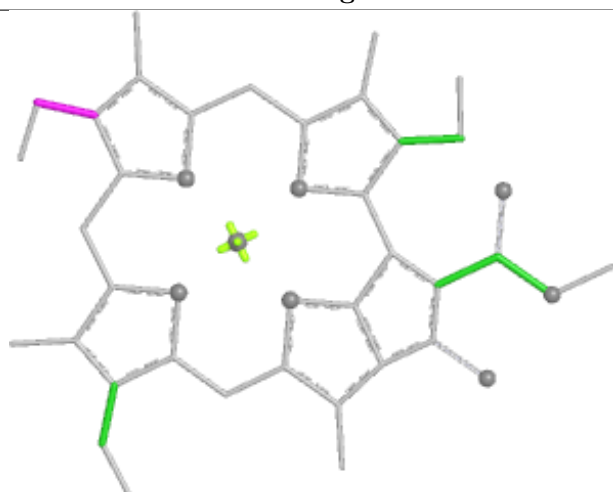
Ligand CLA T 202



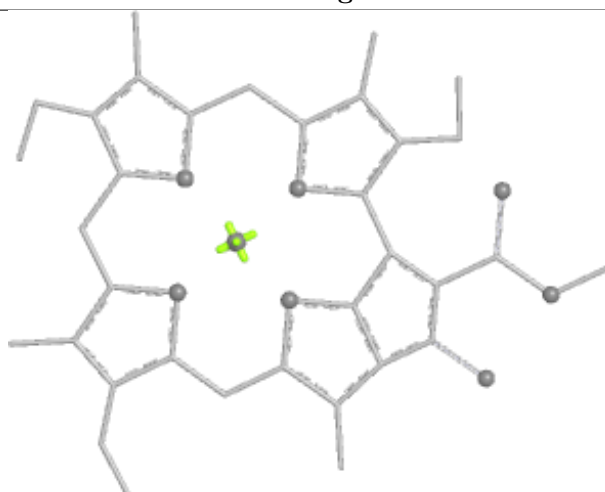
Bond lengths



Bond angles

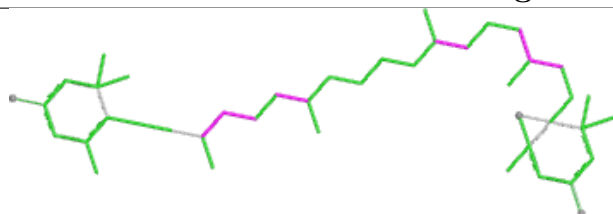


Torsions

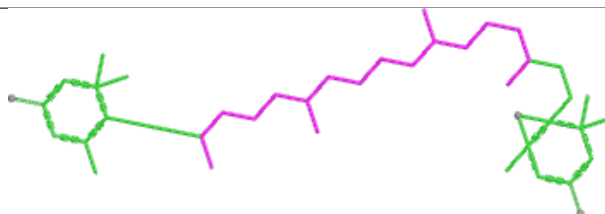


Rings

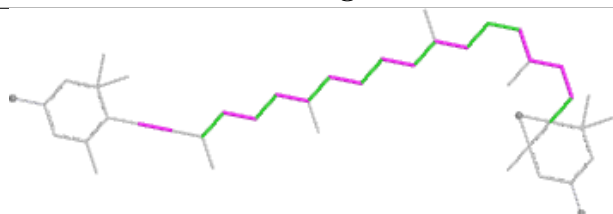
Ligand DD6 A 854



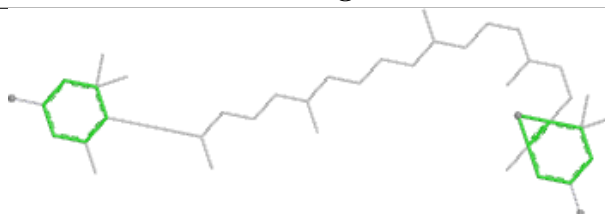
Bond lengths



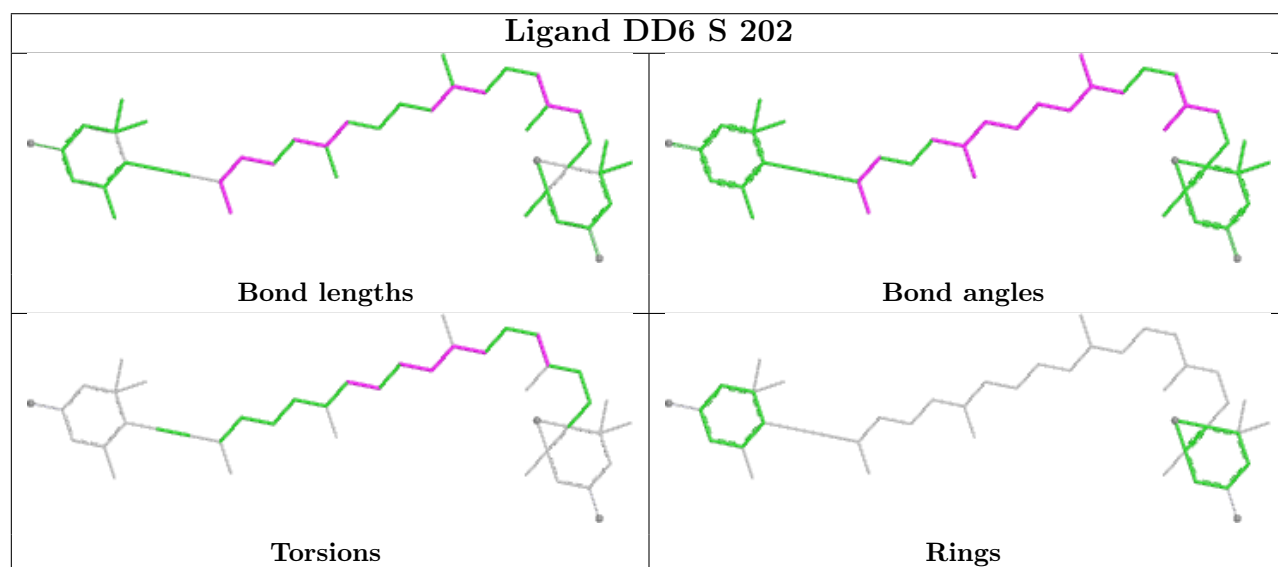
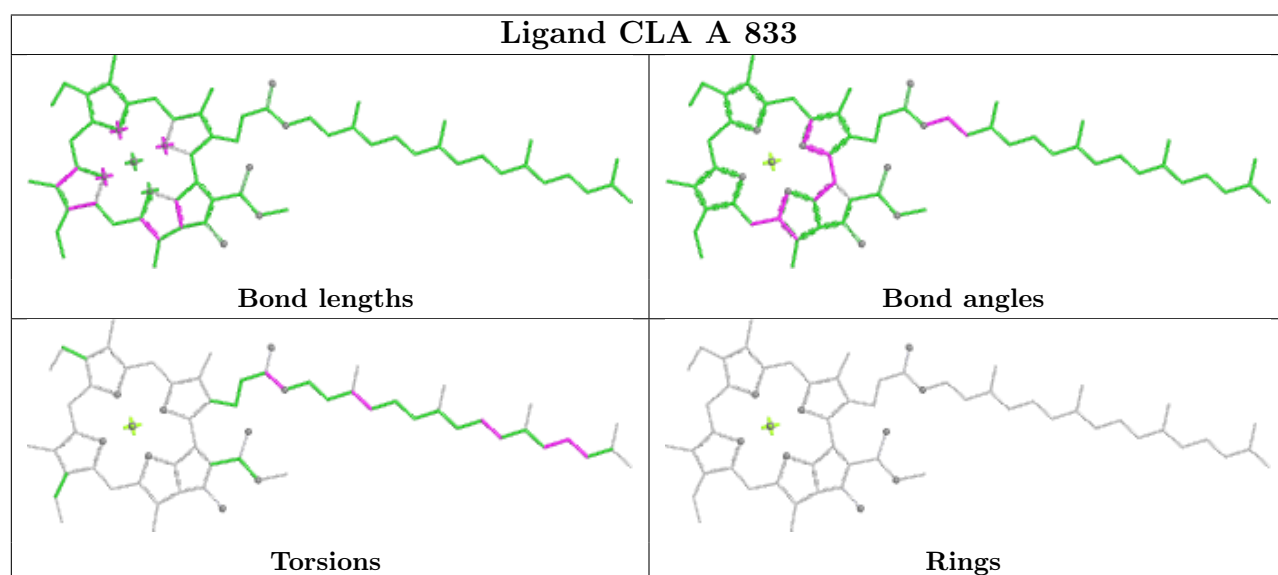
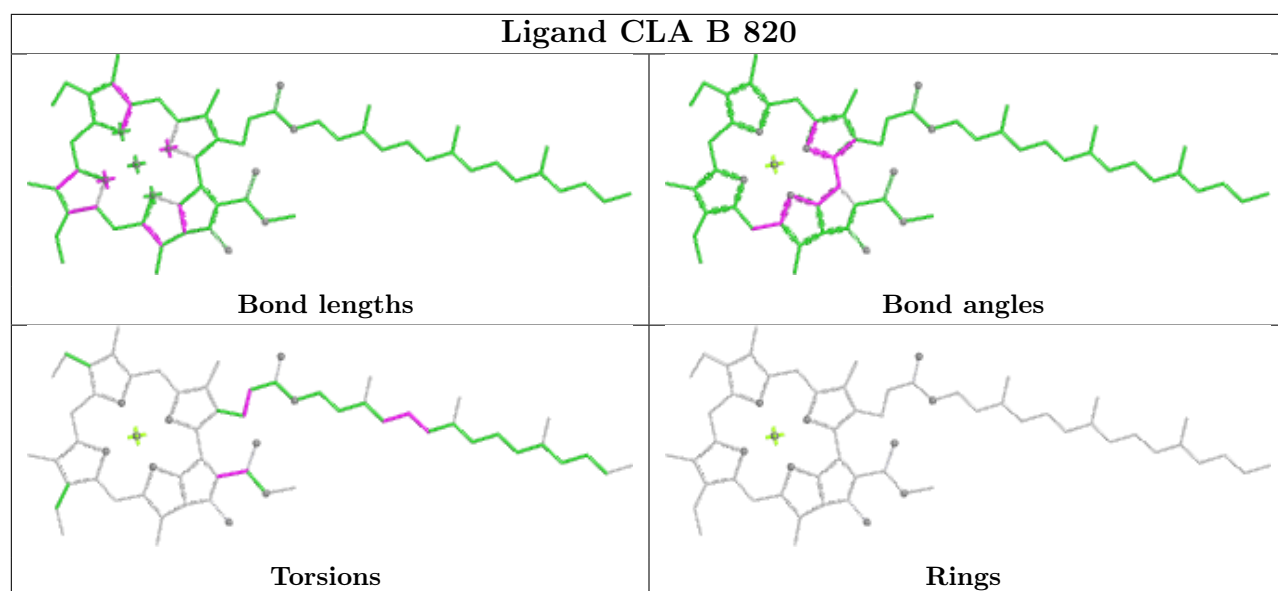
Bond angles

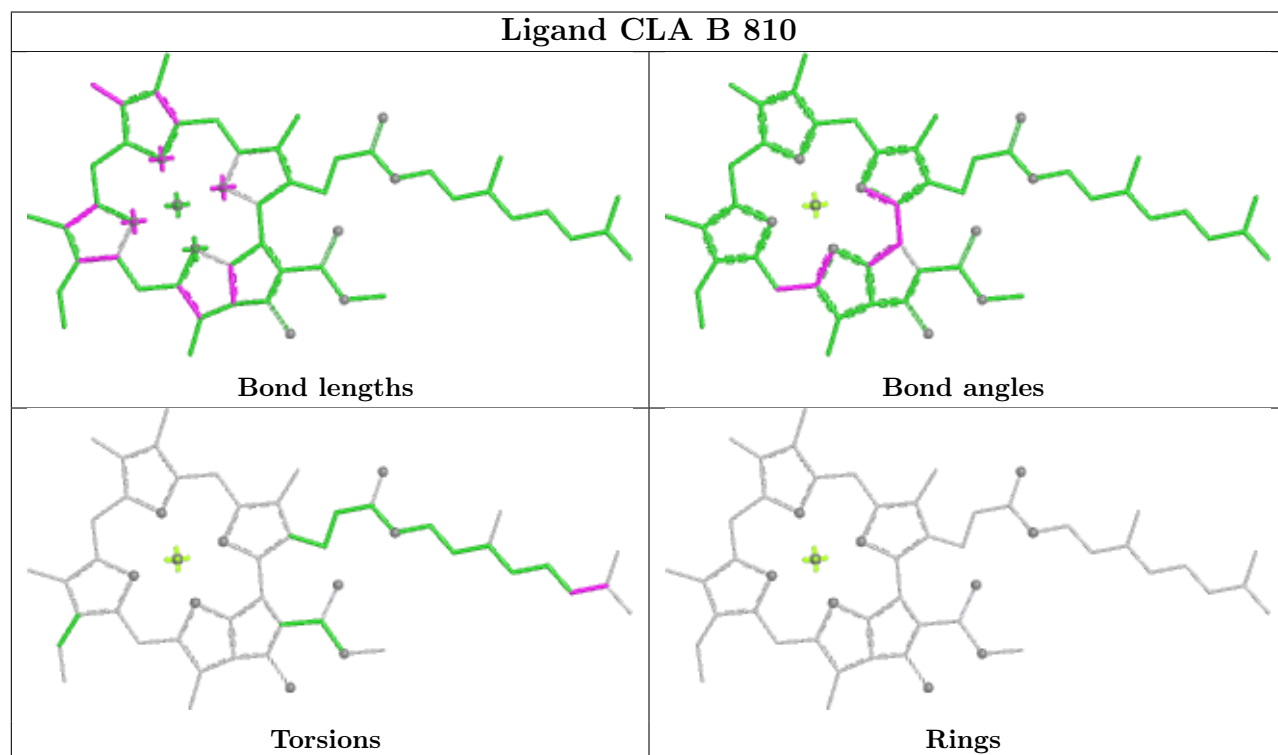
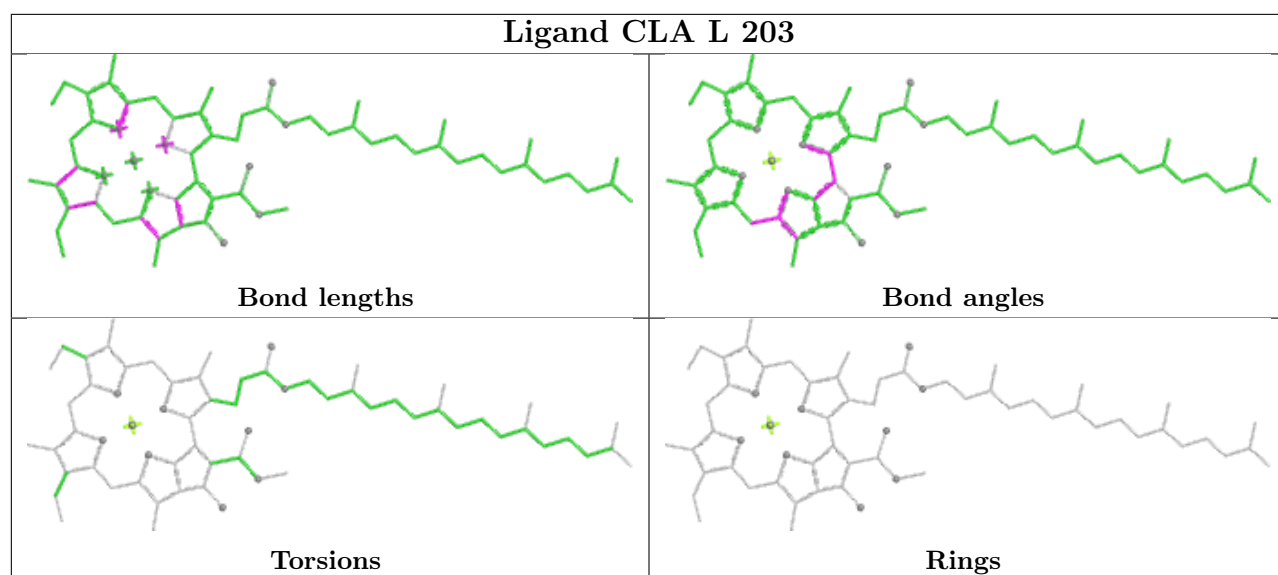


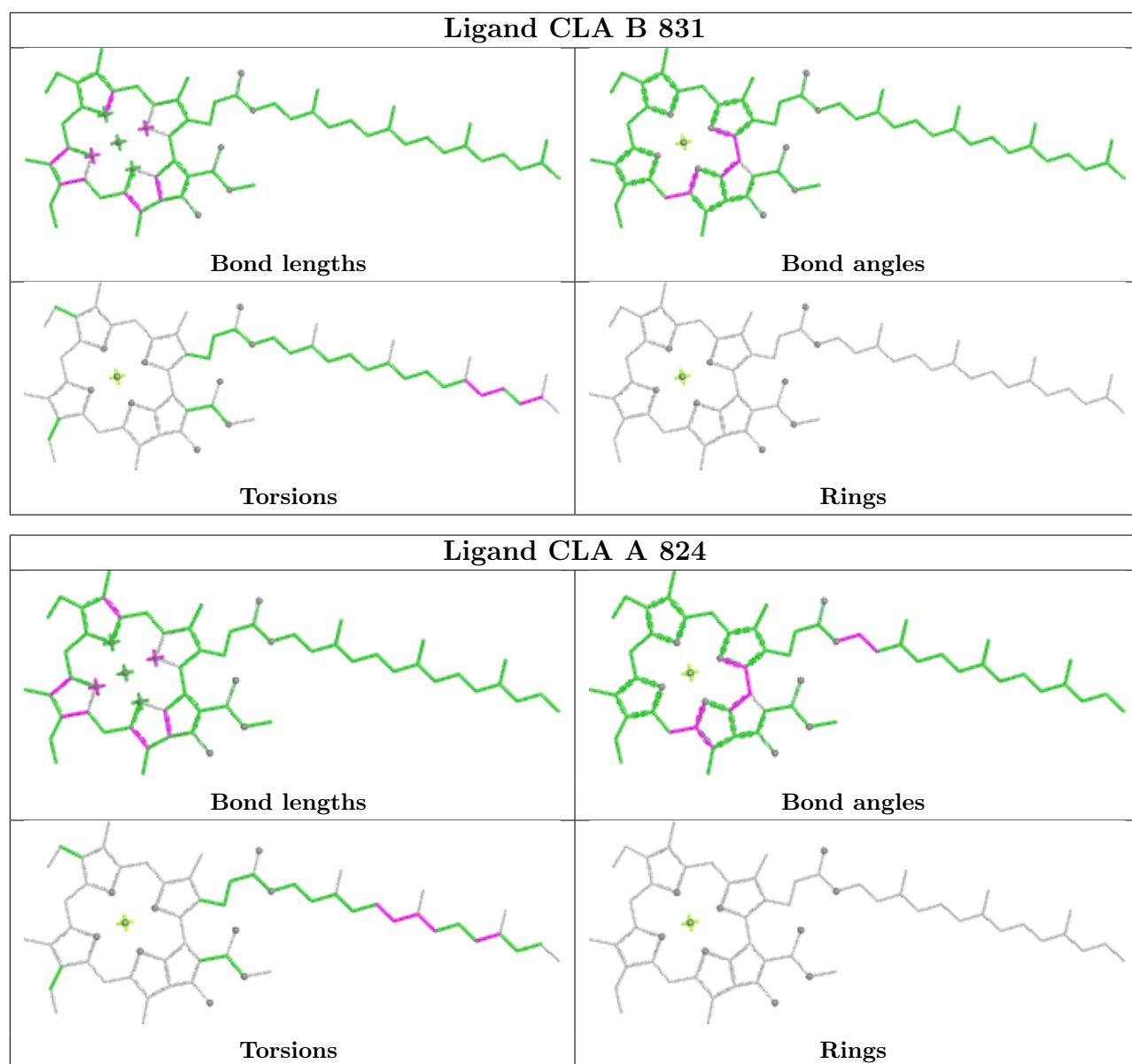
Torsions



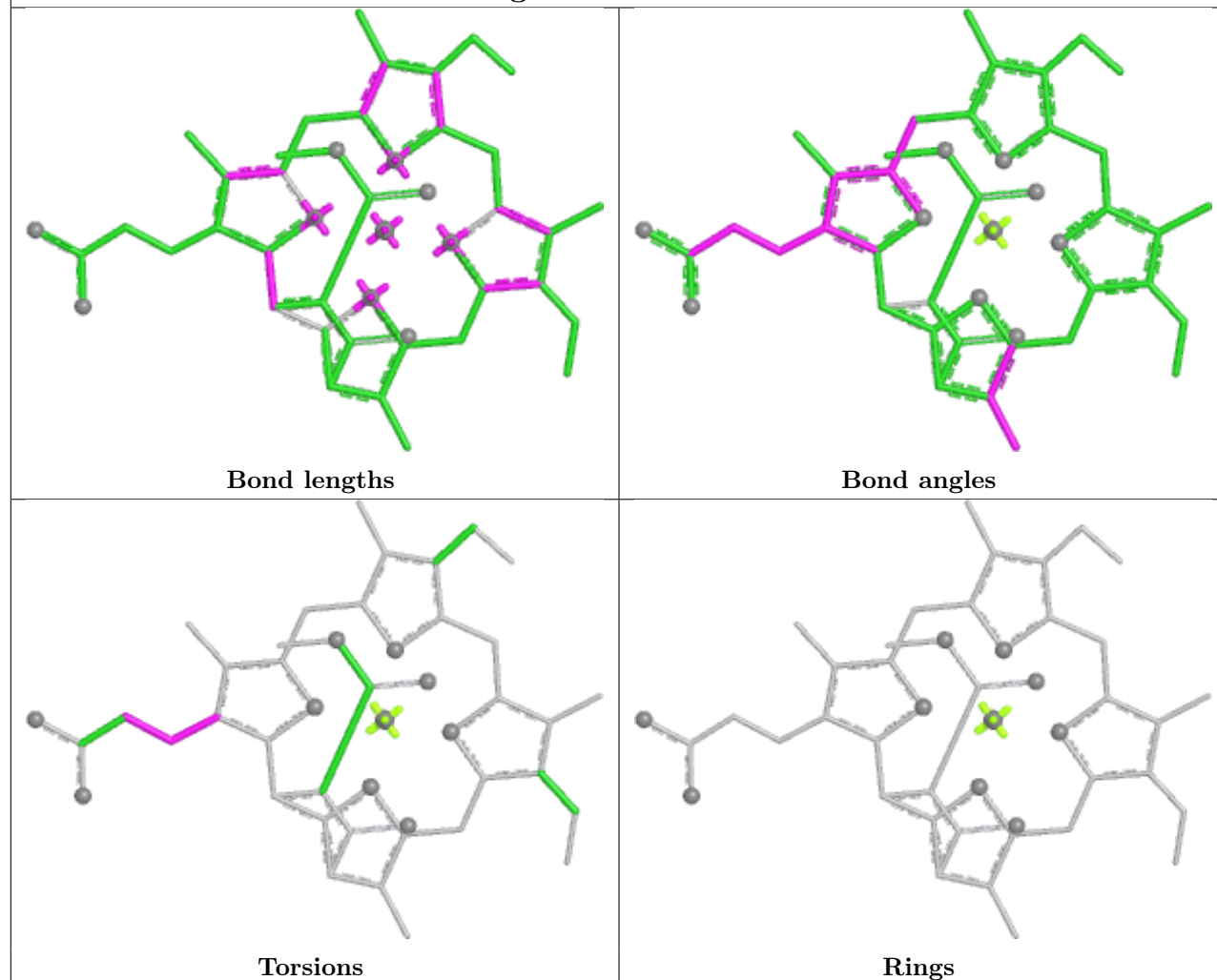
Rings



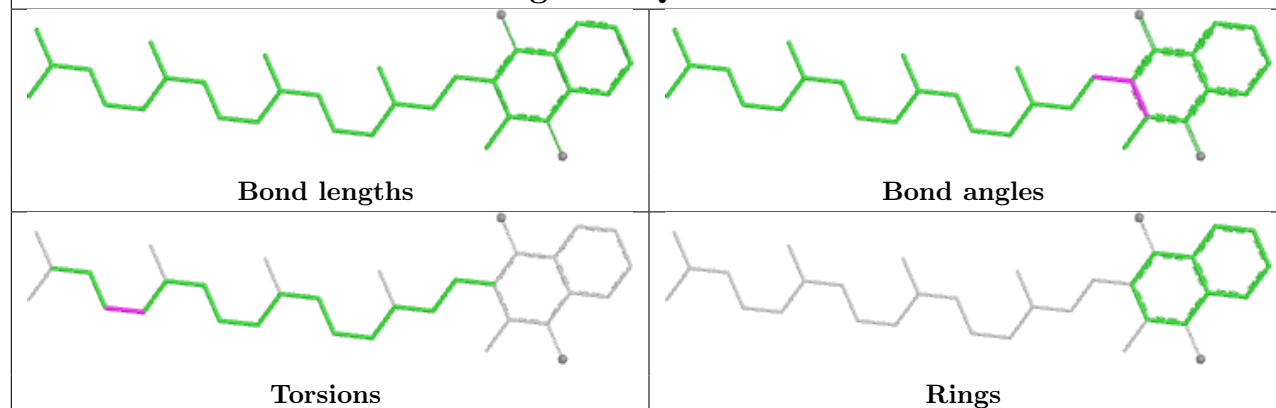




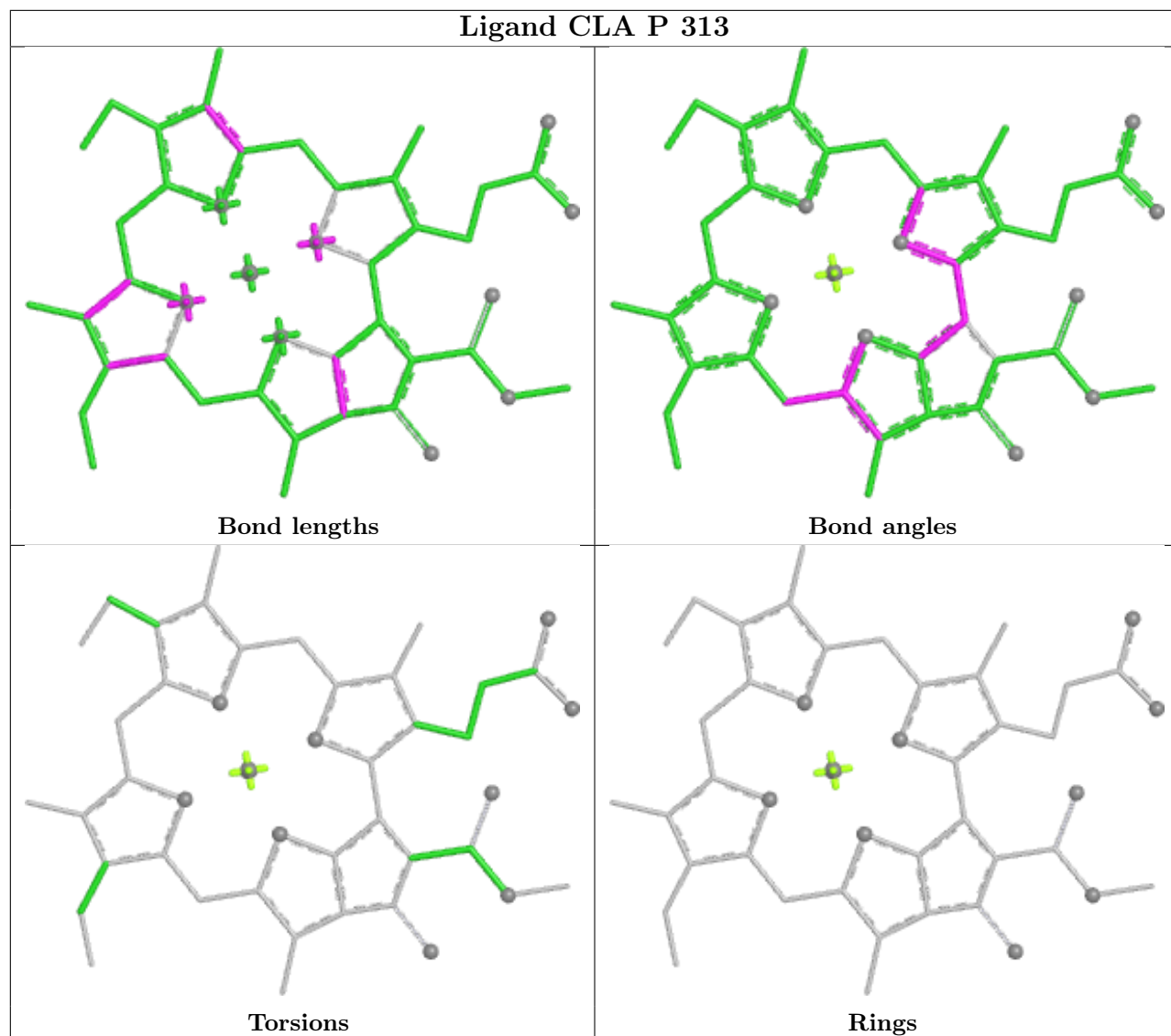
Ligand KC1 P 305



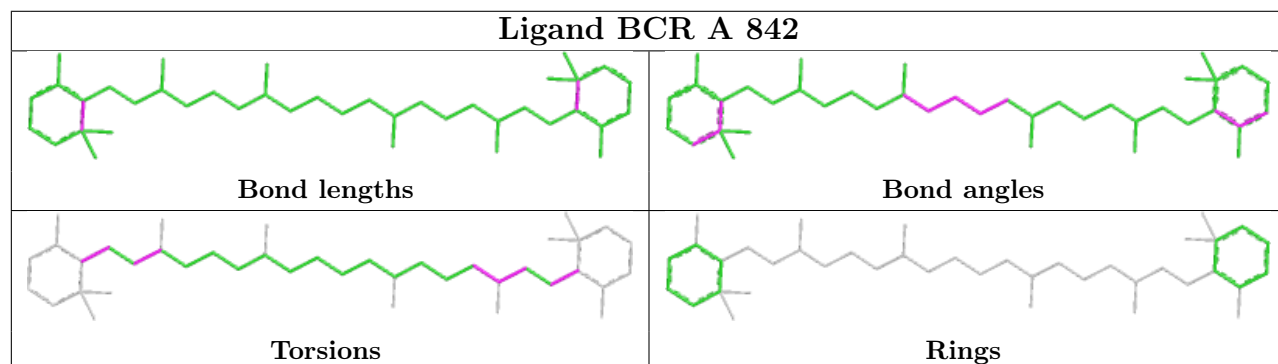
Ligand PQN B 836



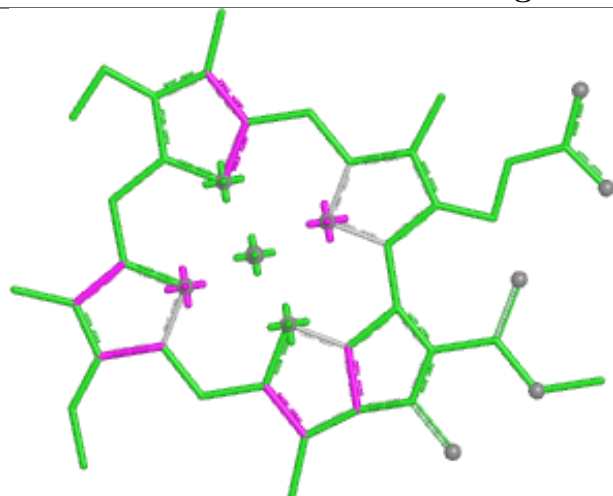
Ligand CLA P 313



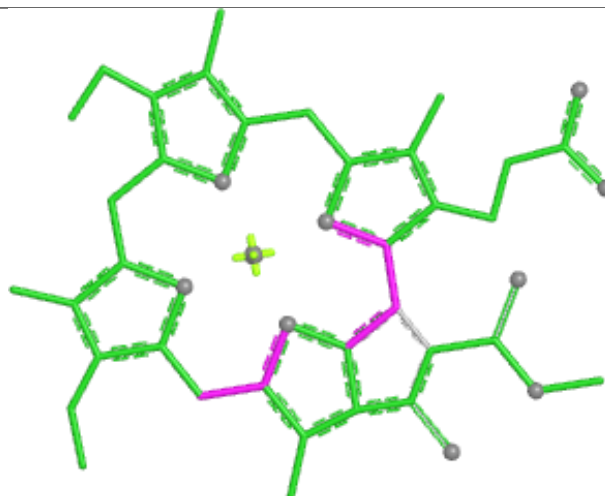
Ligand BCR A 842



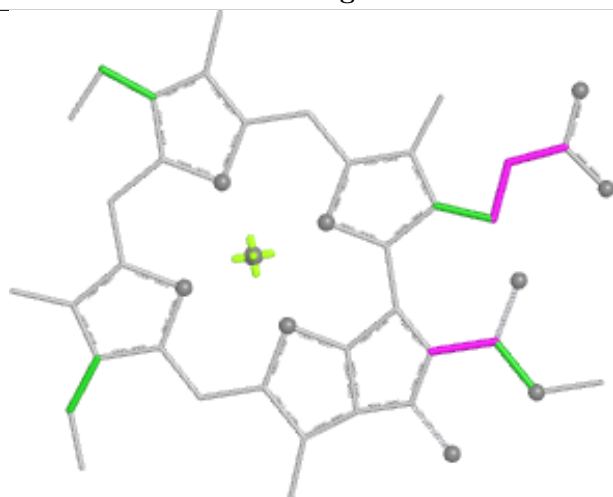
Ligand CLA B 804



Bond lengths



Bond angles

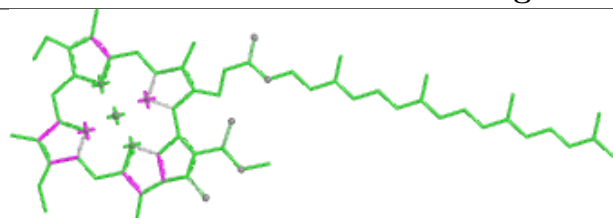


Torsions

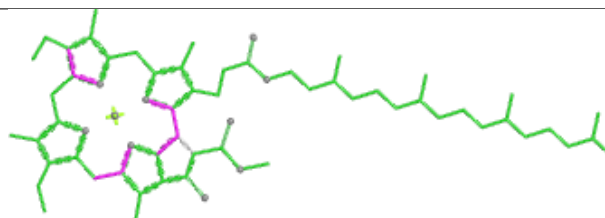


Rings

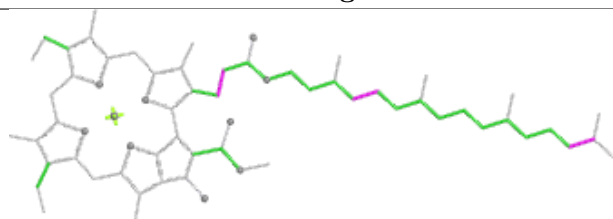
Ligand CLA A 835



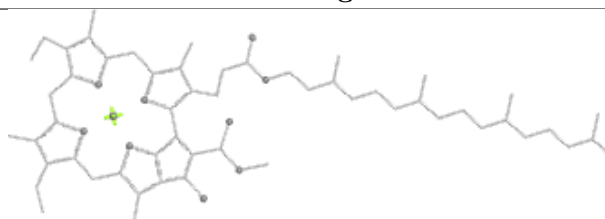
Bond lengths



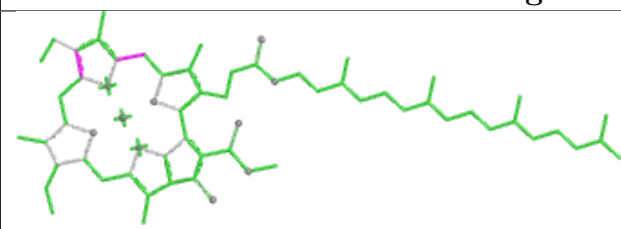
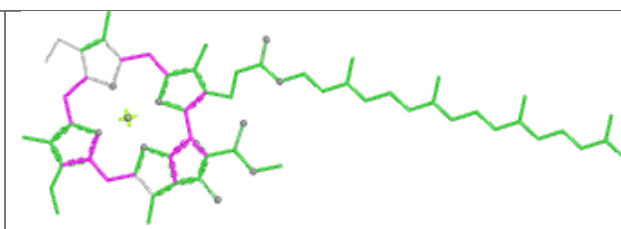
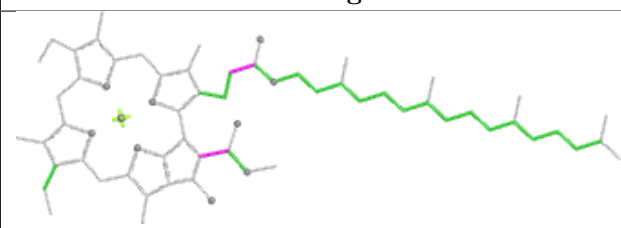
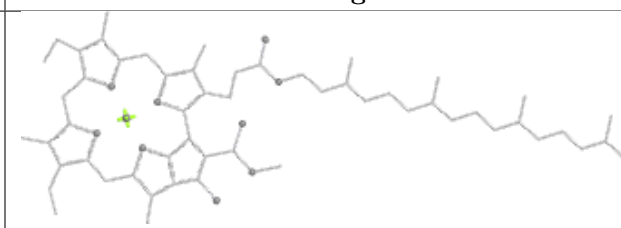
Bond angles

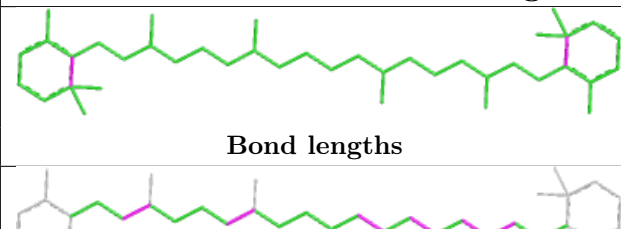
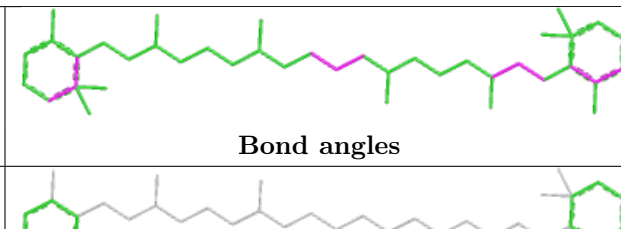
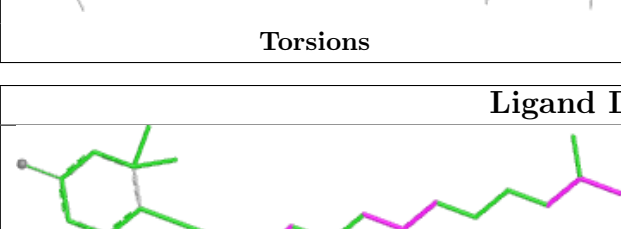



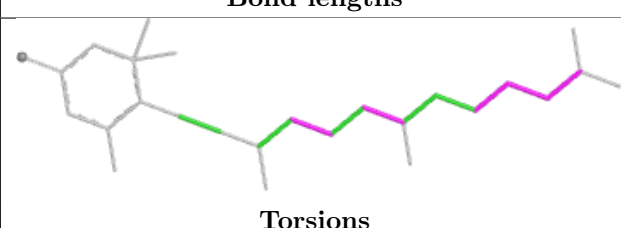
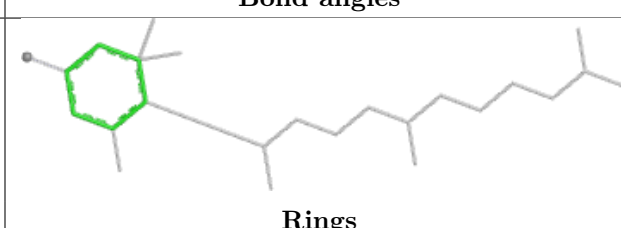


Torsions



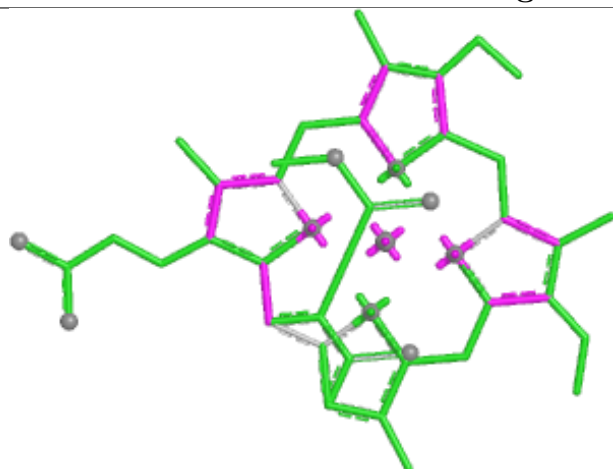
Rings

Ligand CL0 A 848	
	
Bond lengths	Bond angles
	
Torsions	Rings

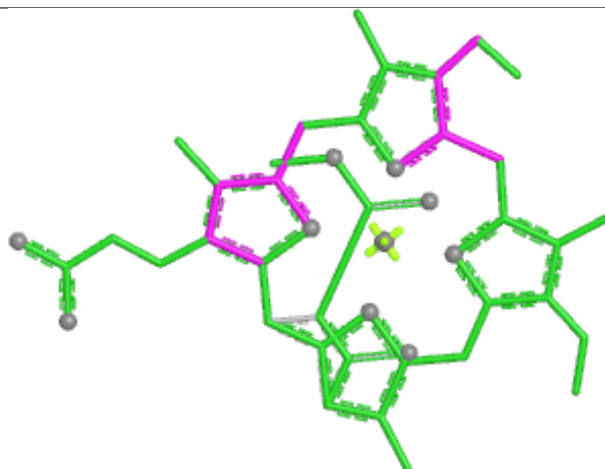
Ligand BCR J 104	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand DD6 U 214	
	
Bond lengths	Bond angles
	
Torsions	Rings

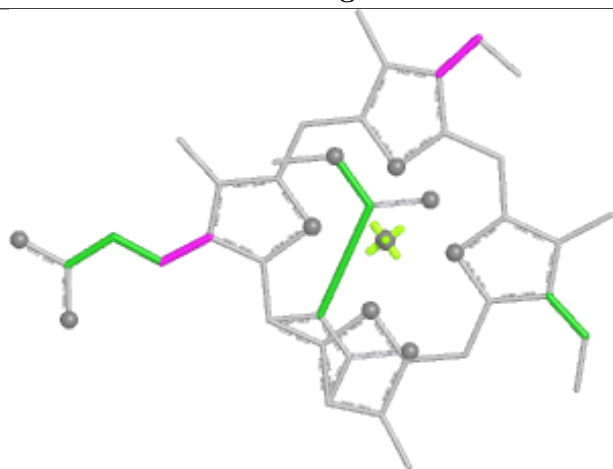
Ligand KC1 P 302



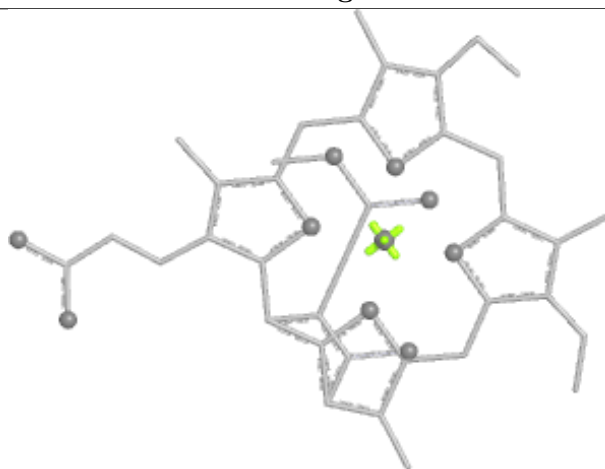
Bond lengths



Bond angles

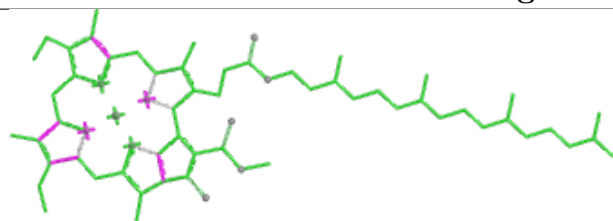


Torsions

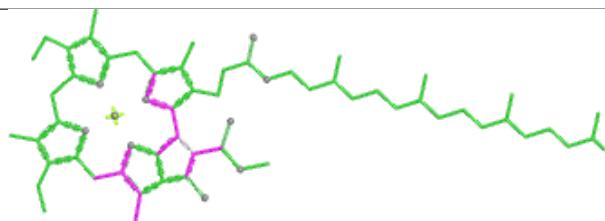


Rings

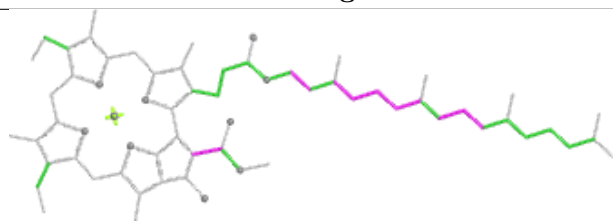
Ligand CLA S 201



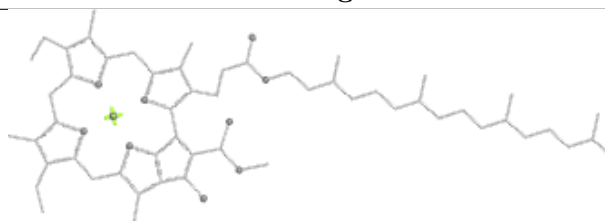
Bond lengths



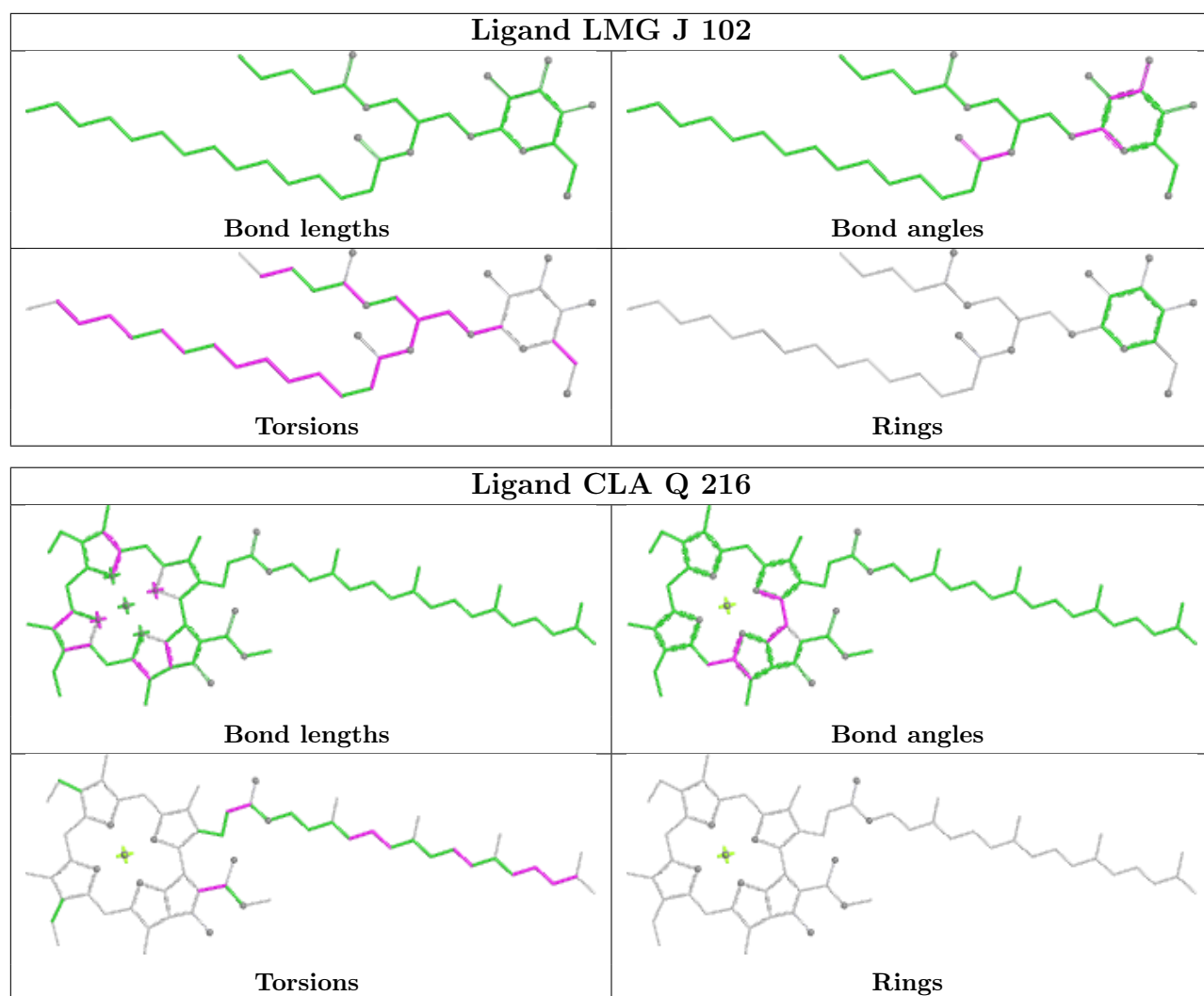
Bond angles

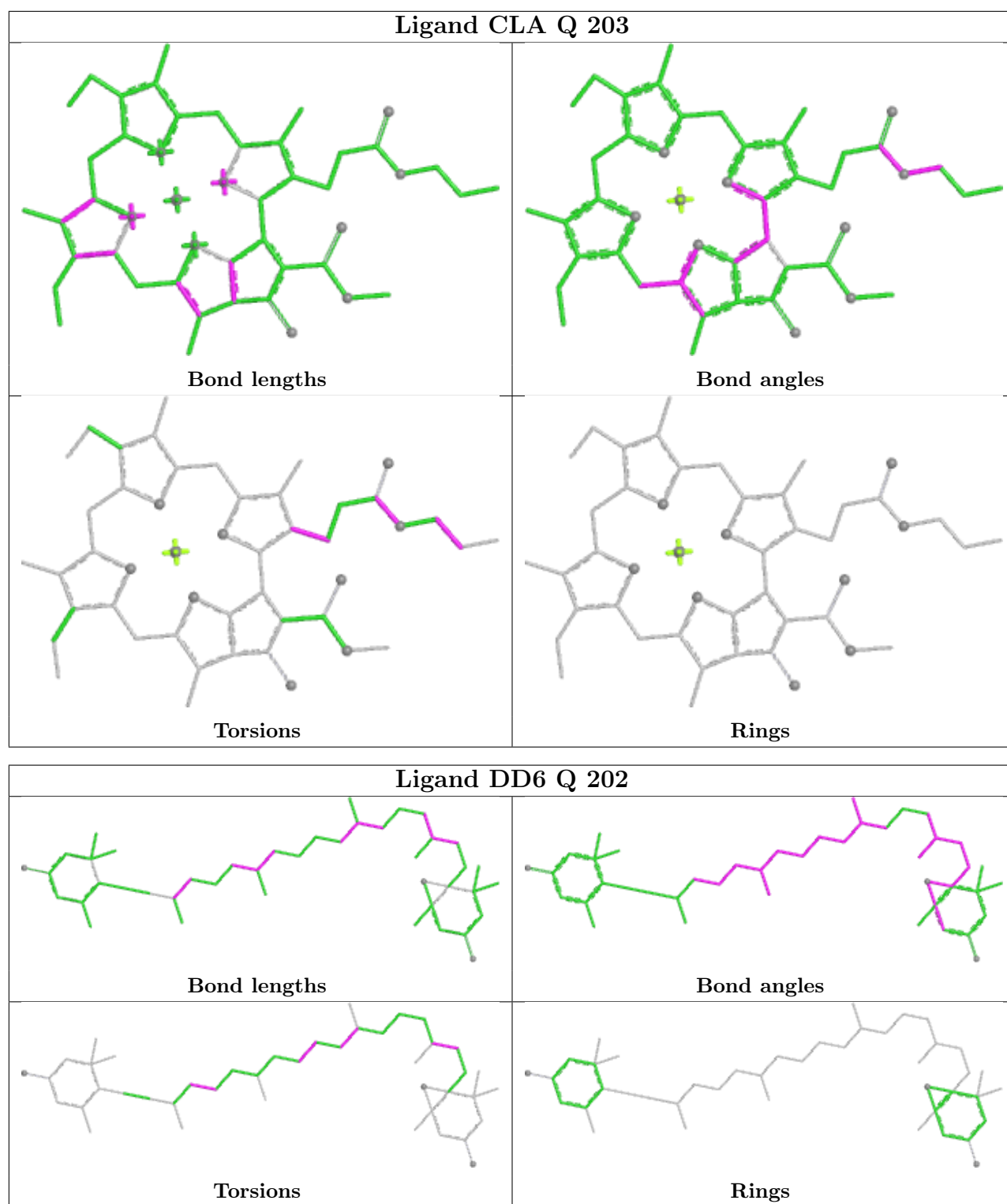


Torsions

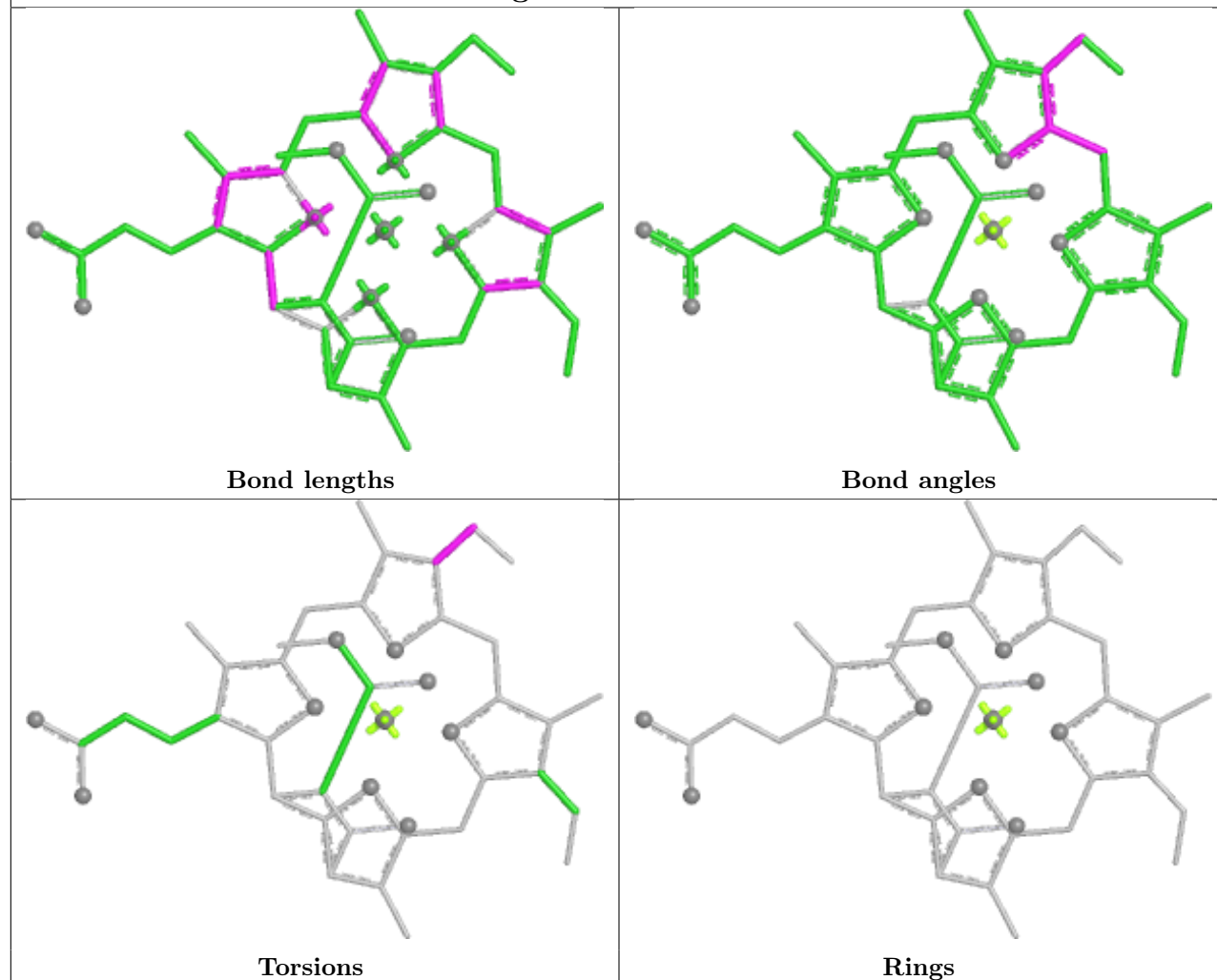


Rings

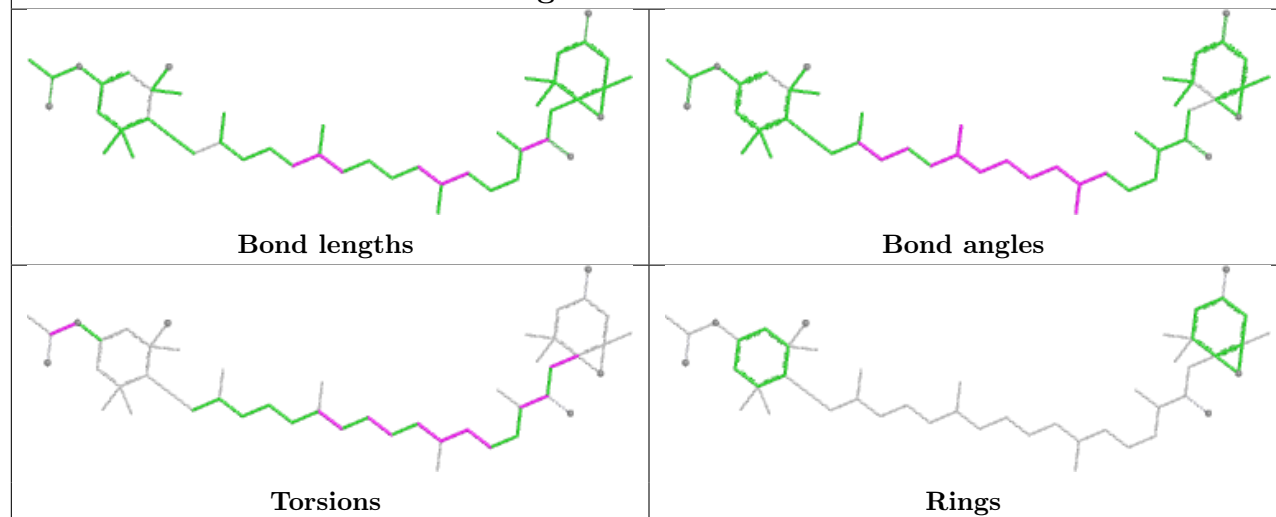


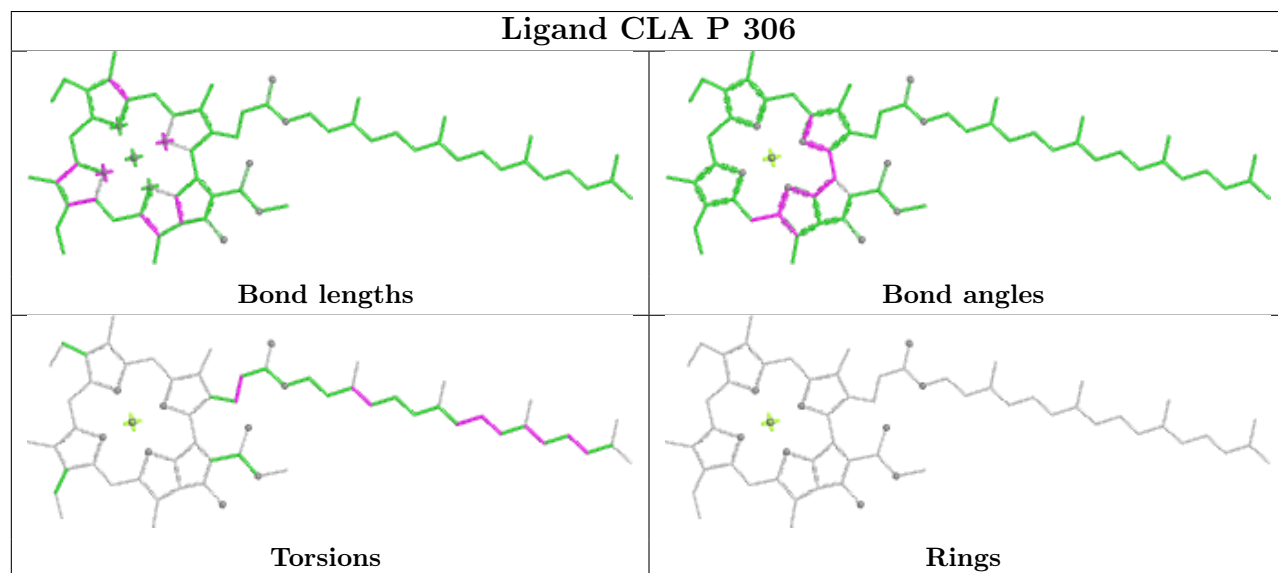
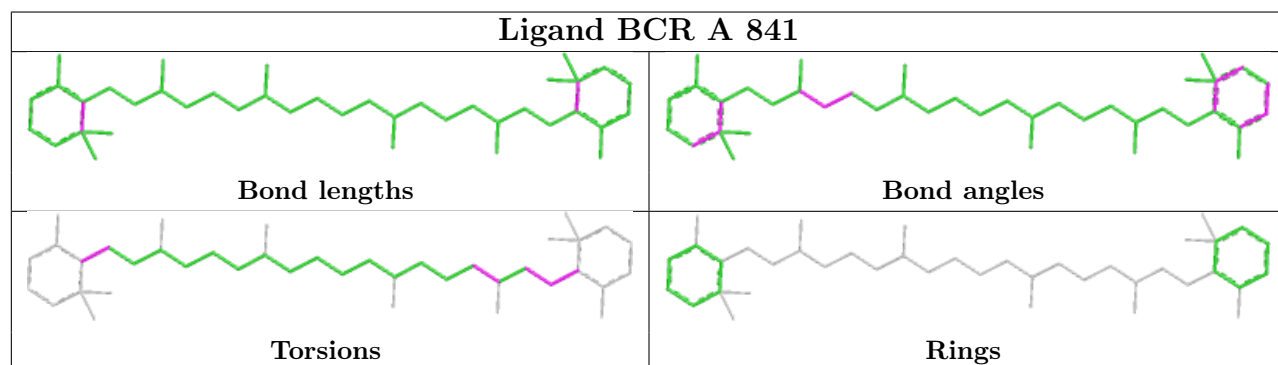
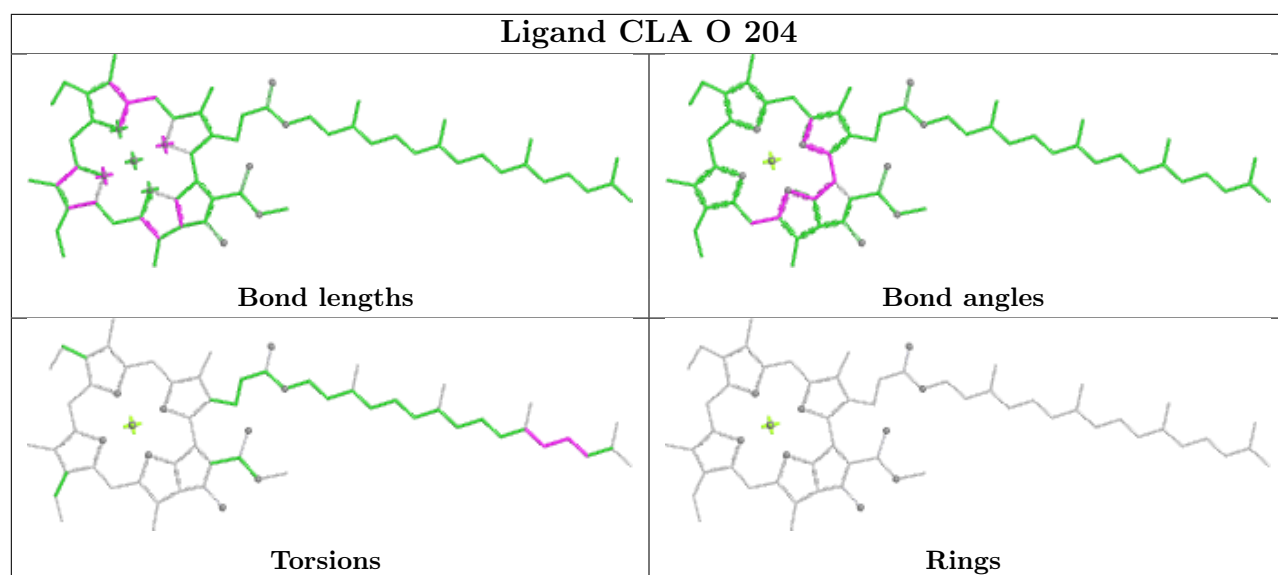


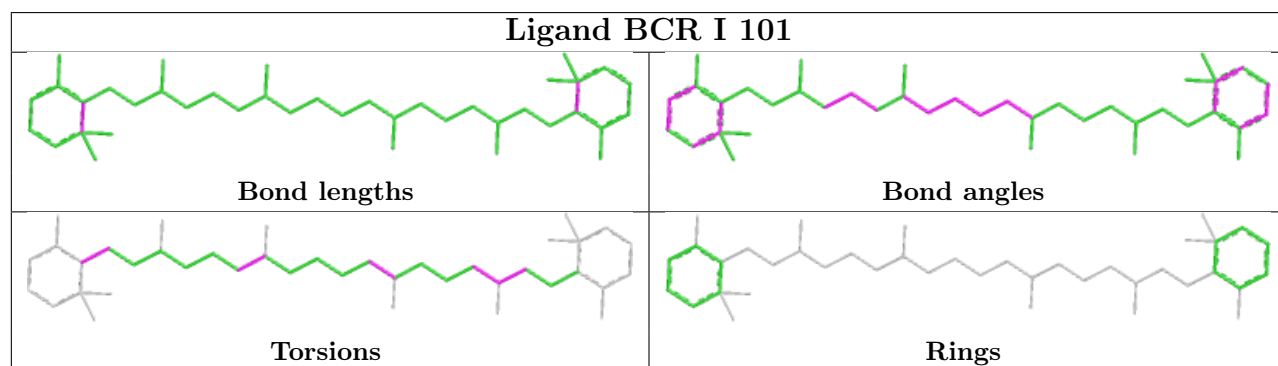
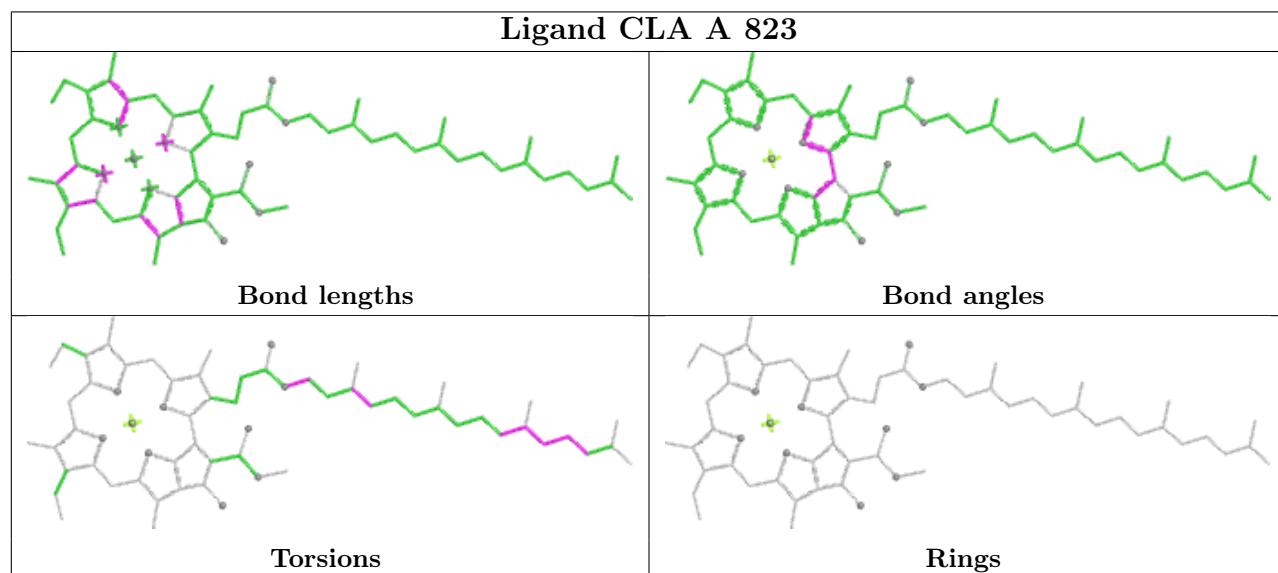
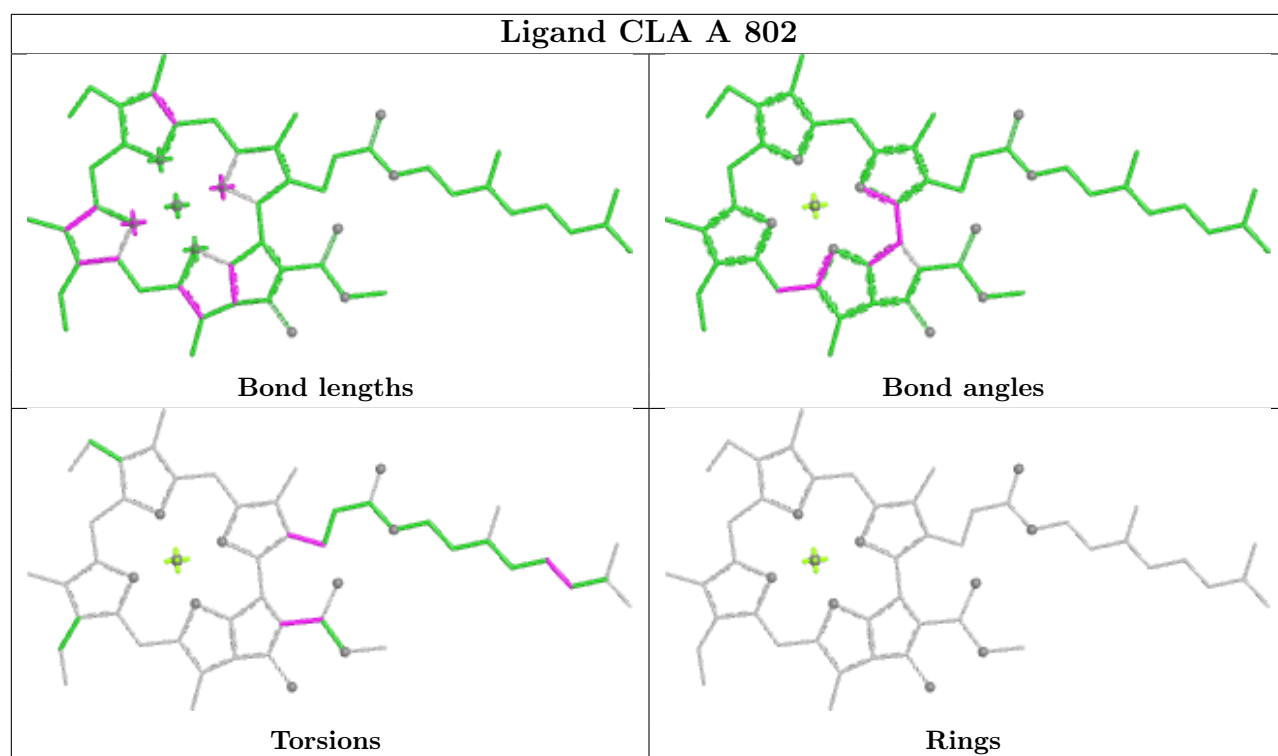
Ligand KC1 S 208

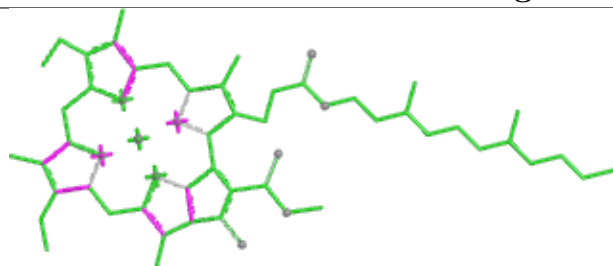


Ligand A86 U 202

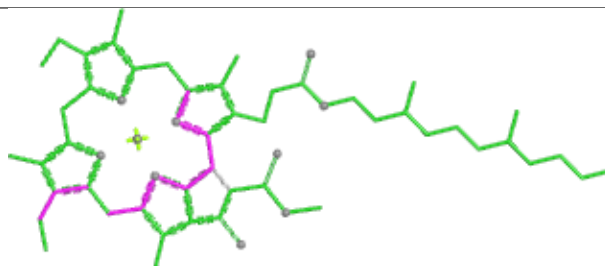




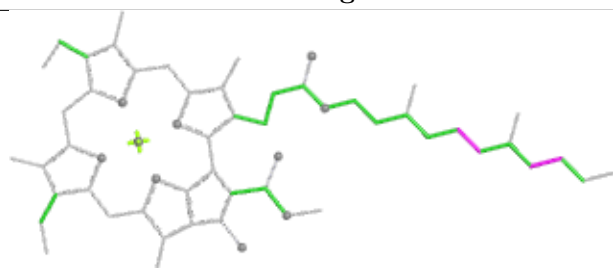


Ligand CLA B 830

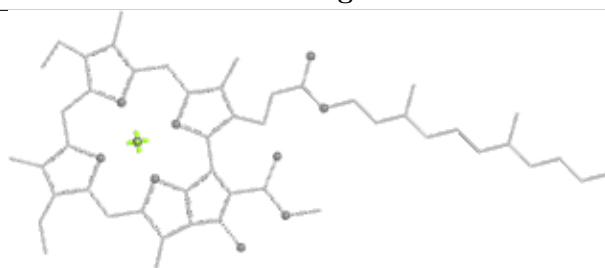
Bond lengths



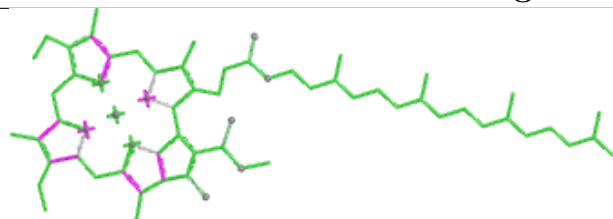
Bond angles



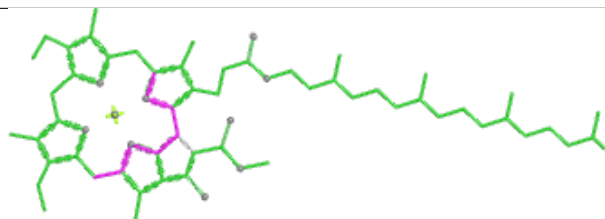
Torsions



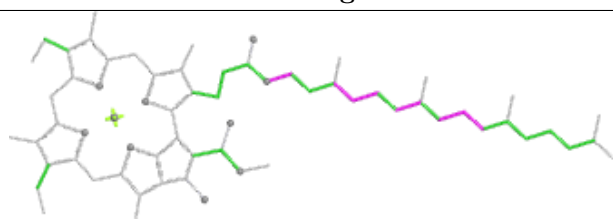
Rings

Ligand CLA B 809

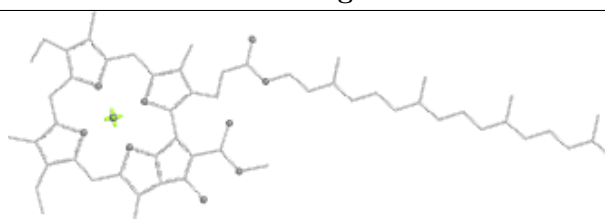
Bond lengths



Bond angles

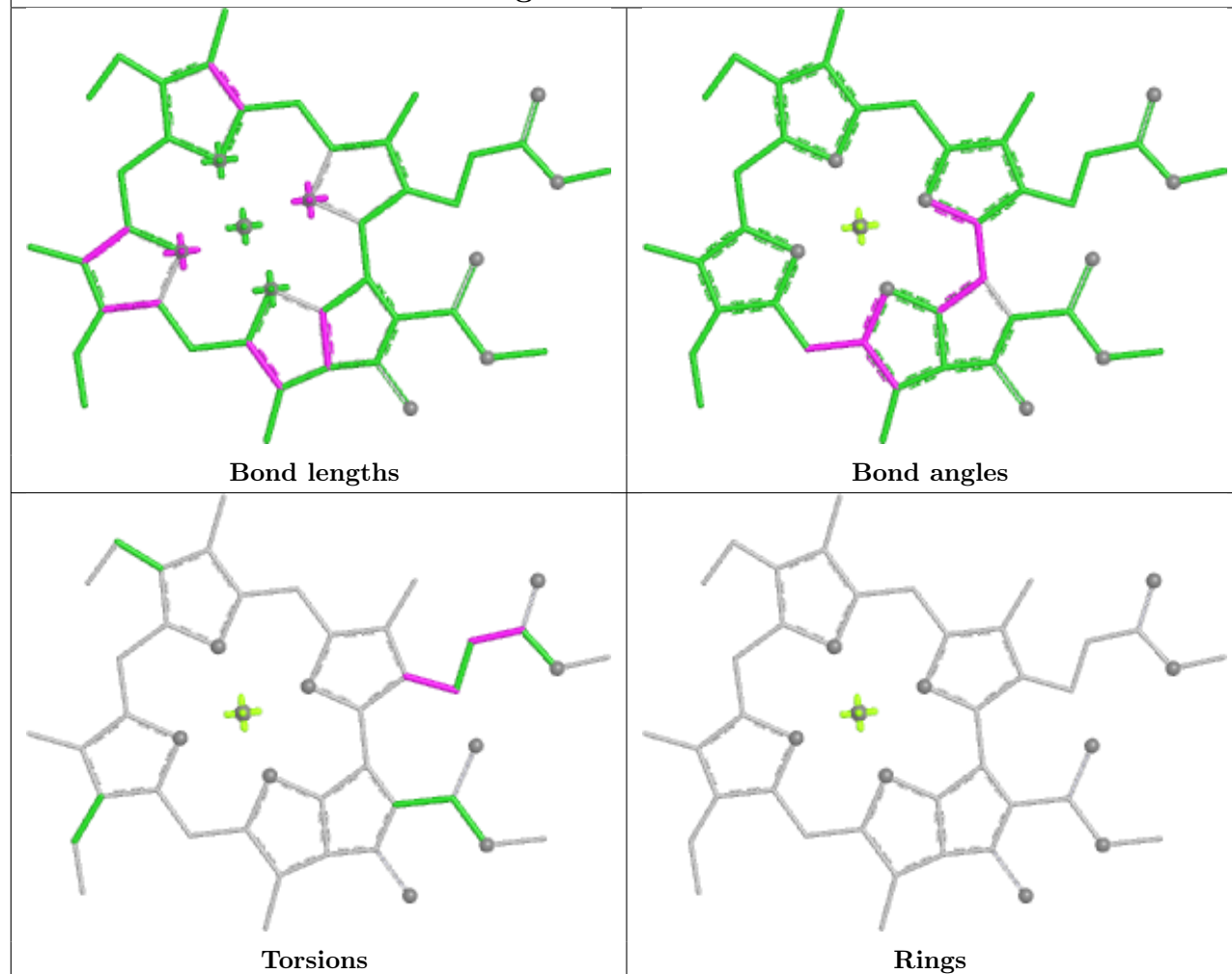


Torsions

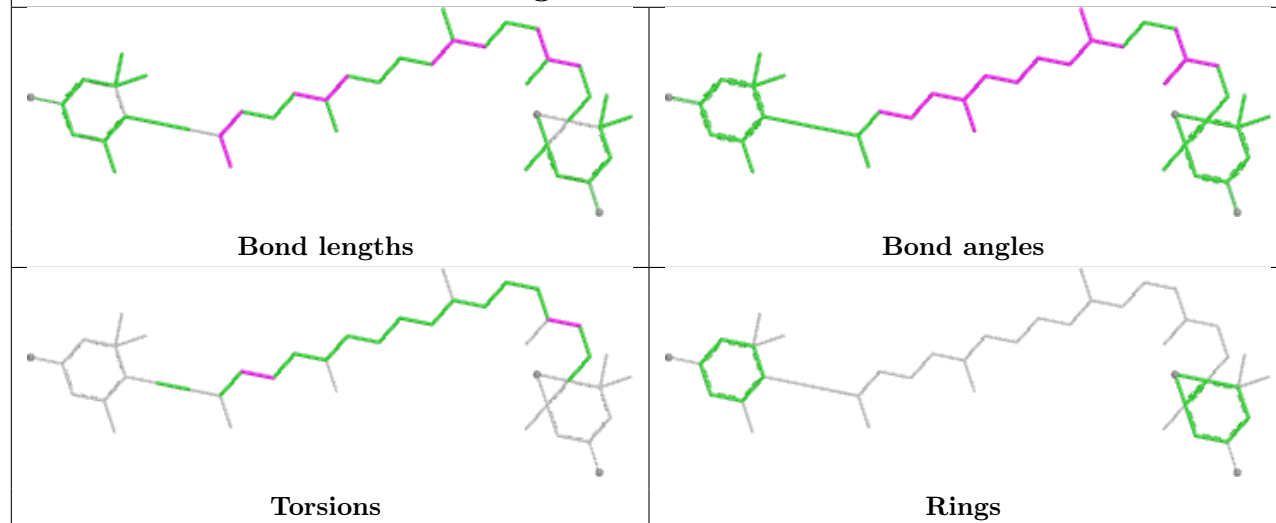


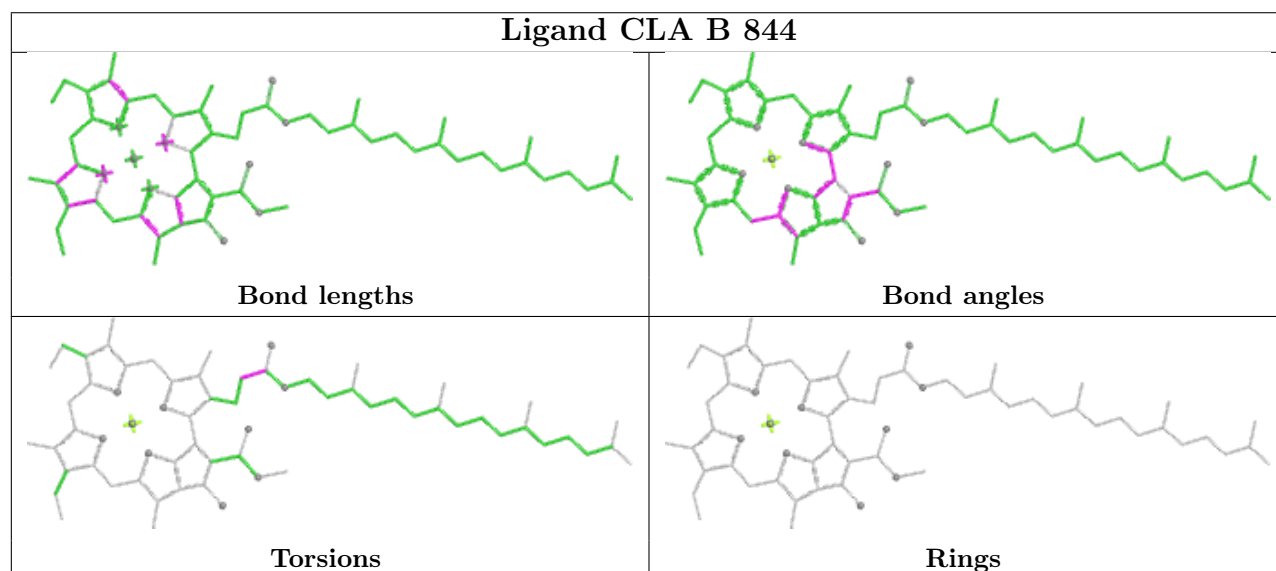
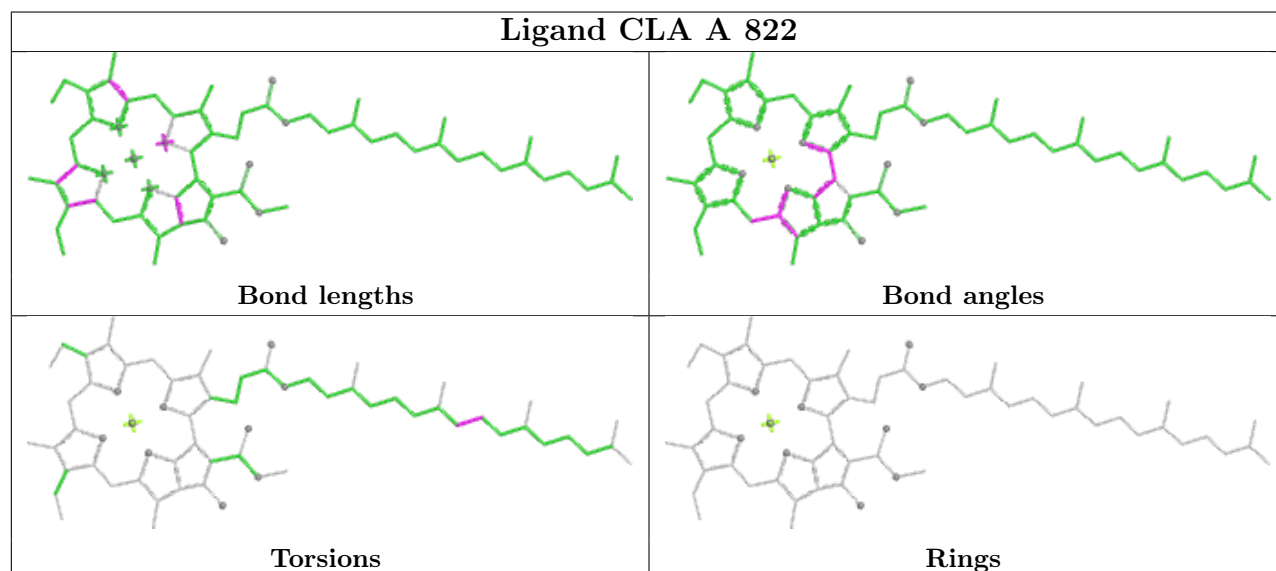
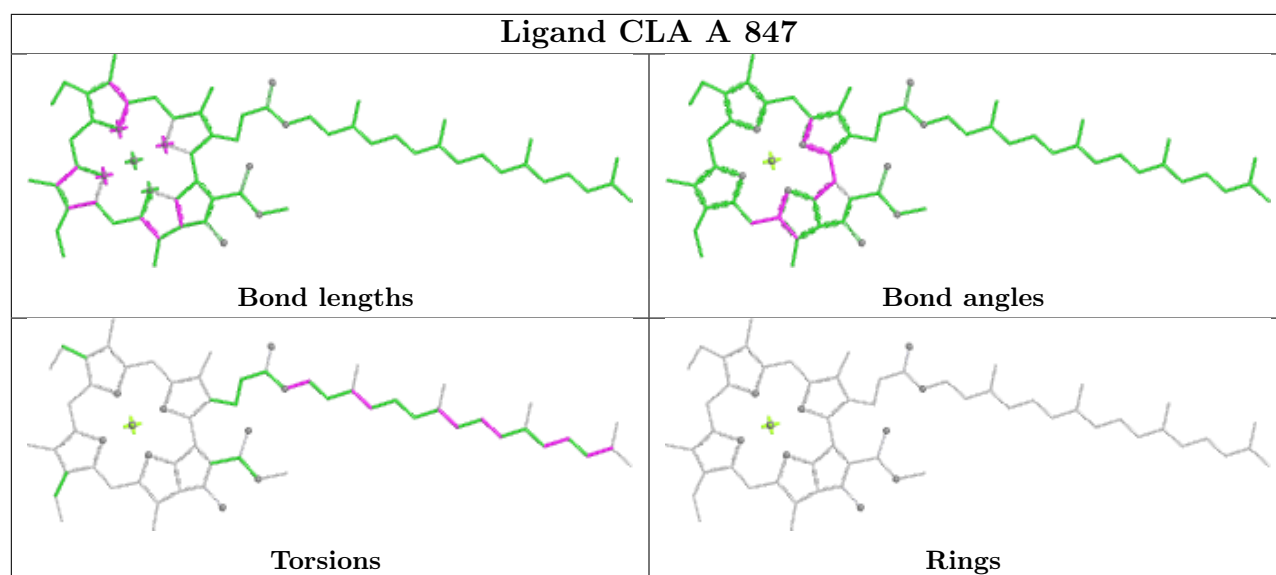
Rings

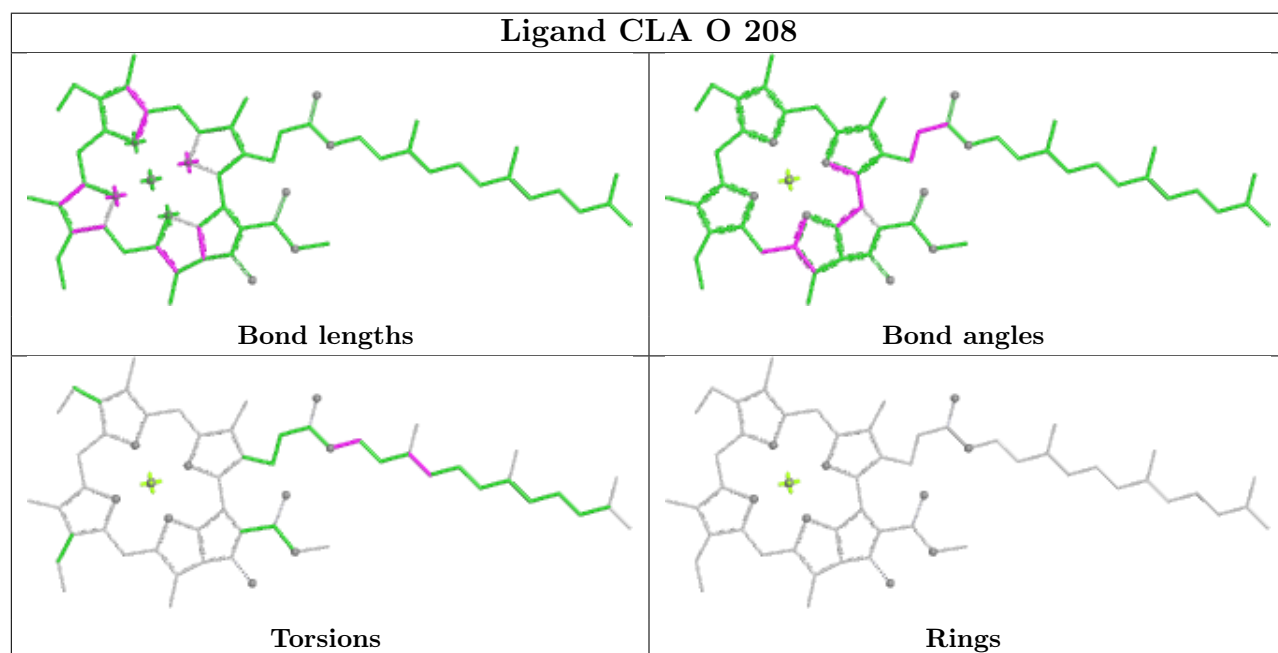
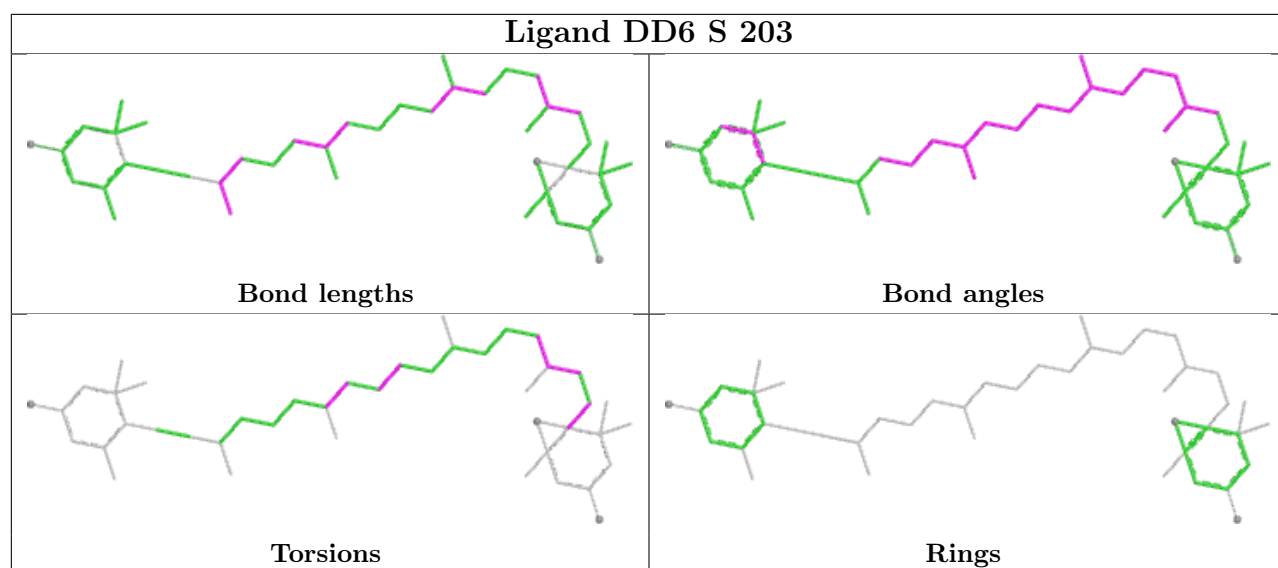
Ligand CLA F 803



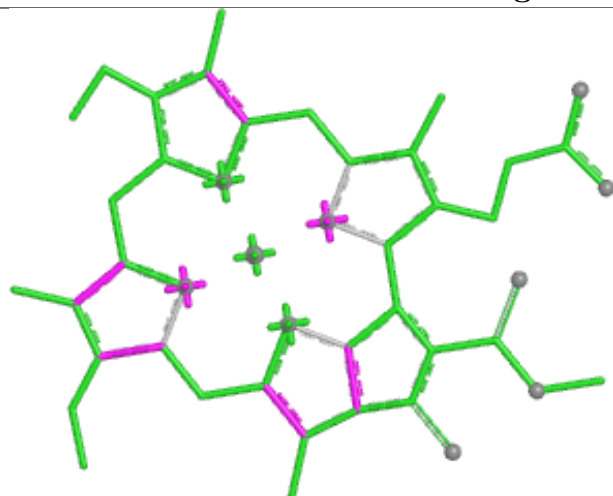
Ligand DD6 O 201



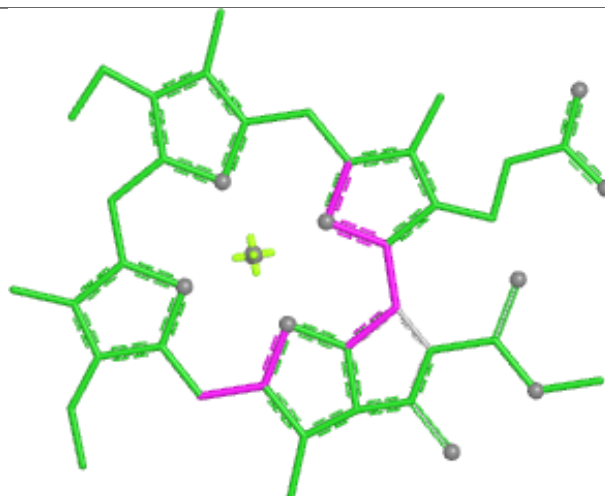




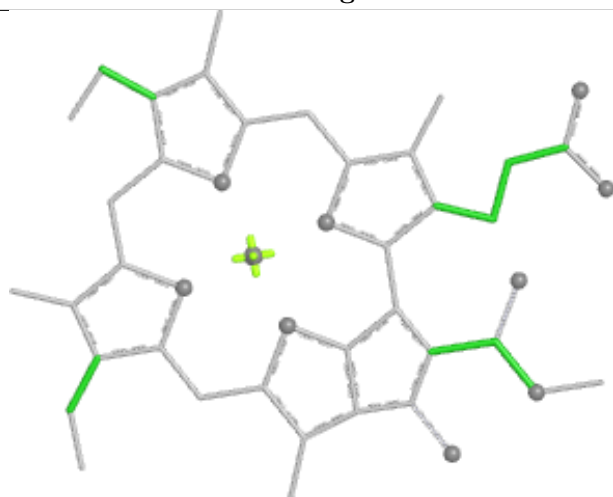
Ligand CLA A 817



Bond lengths



Bond angles

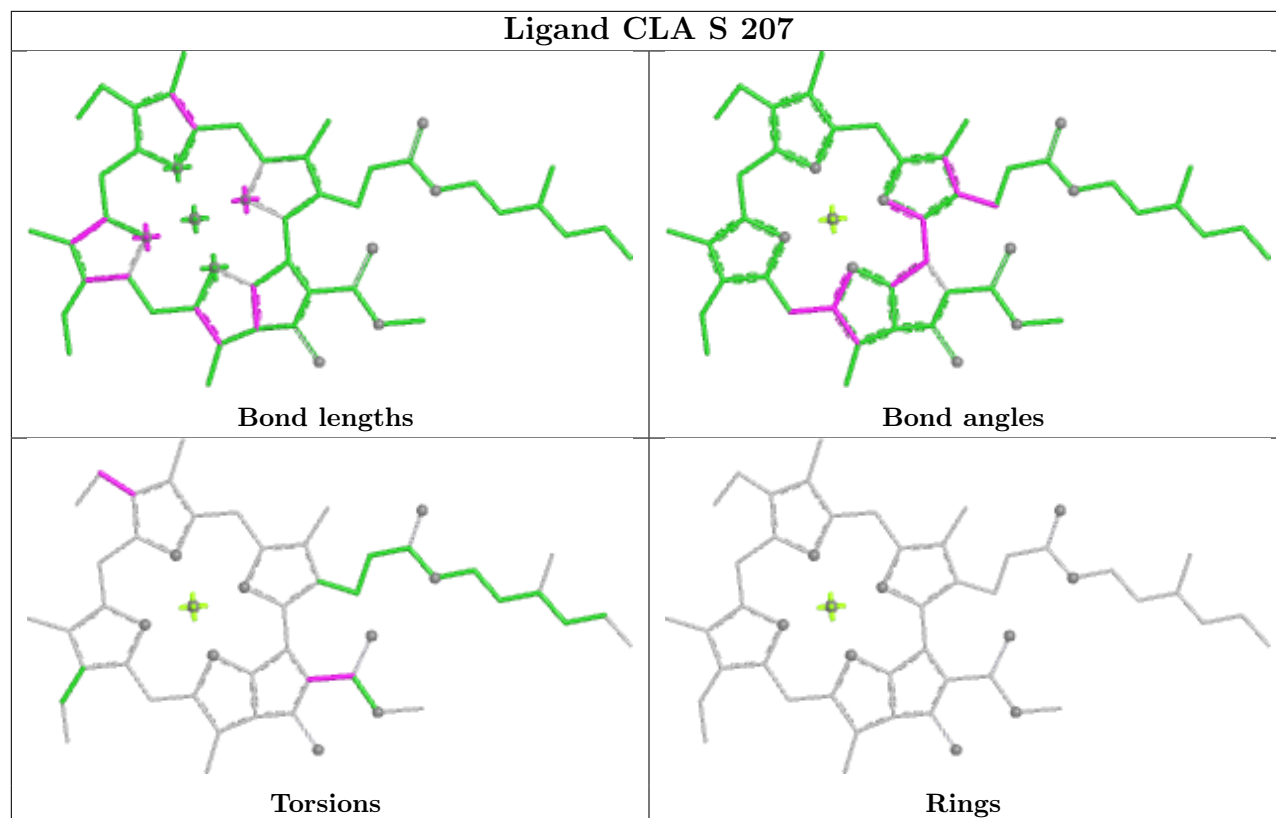


Torsions

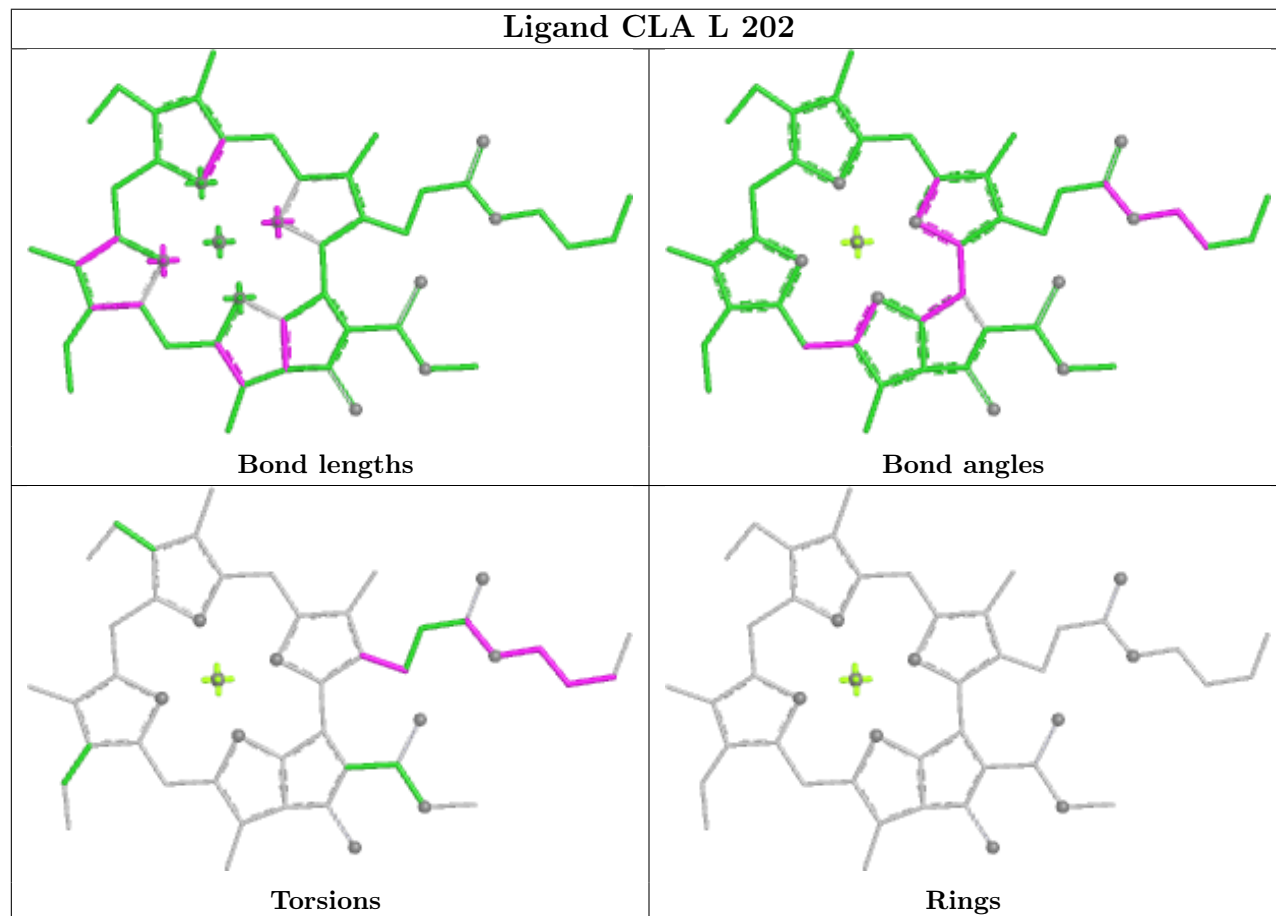


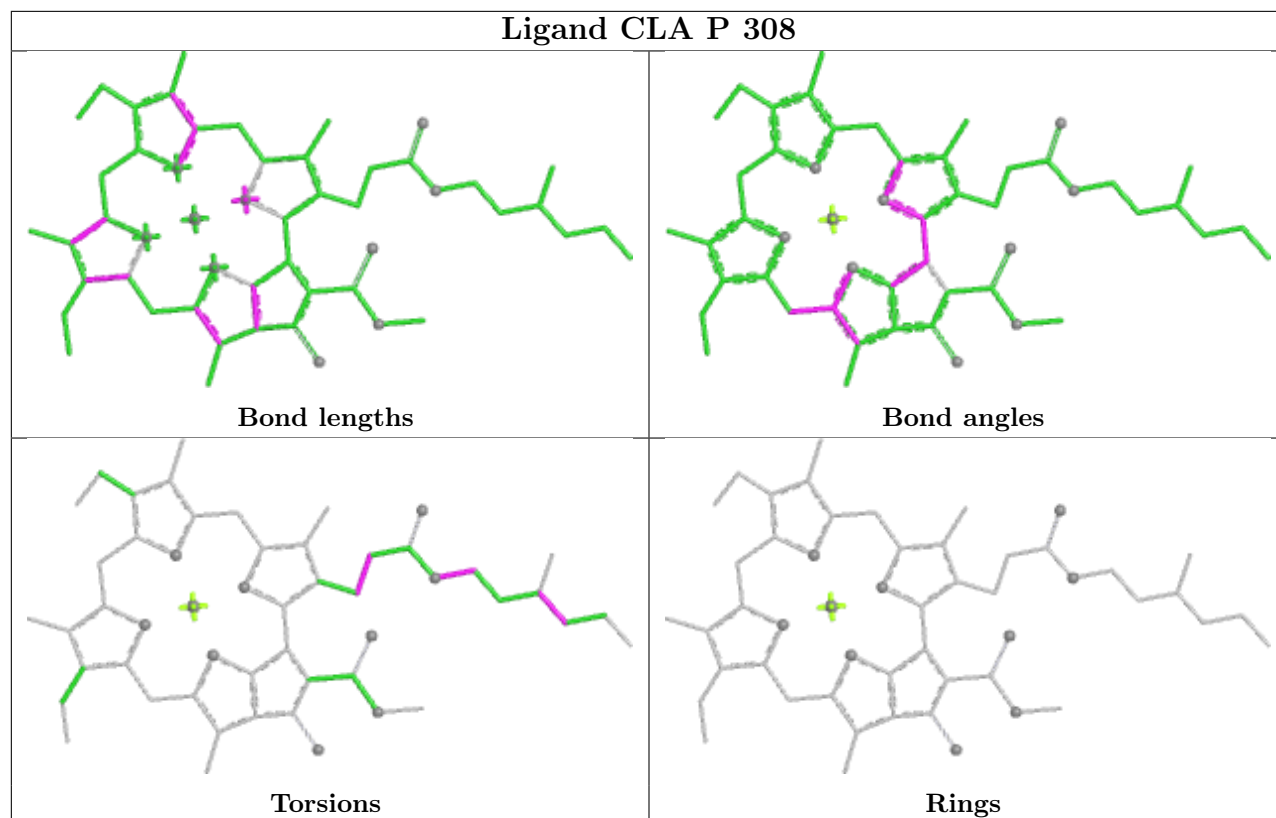
Rings

Ligand CLA S 207

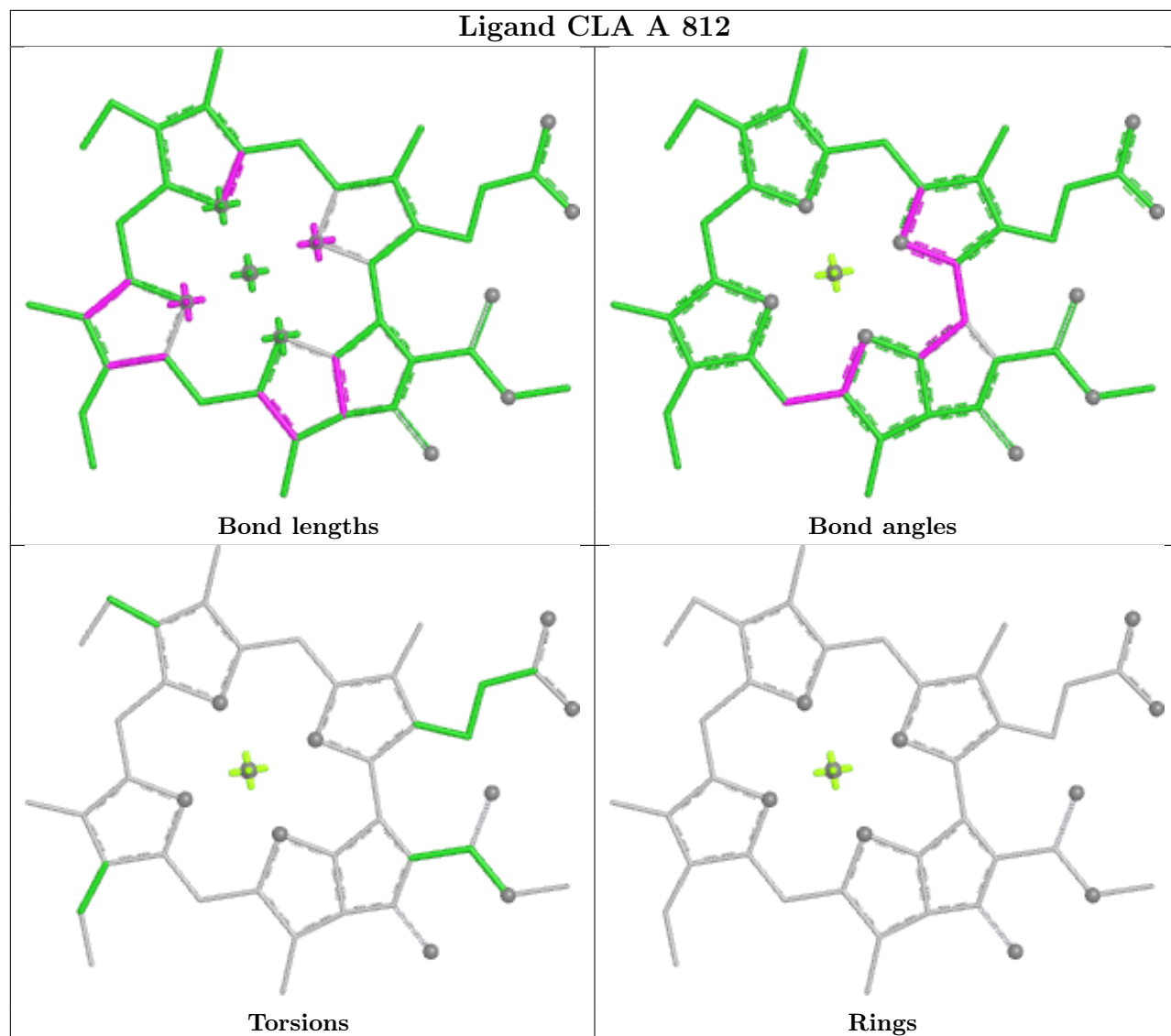


Ligand CLA L 202

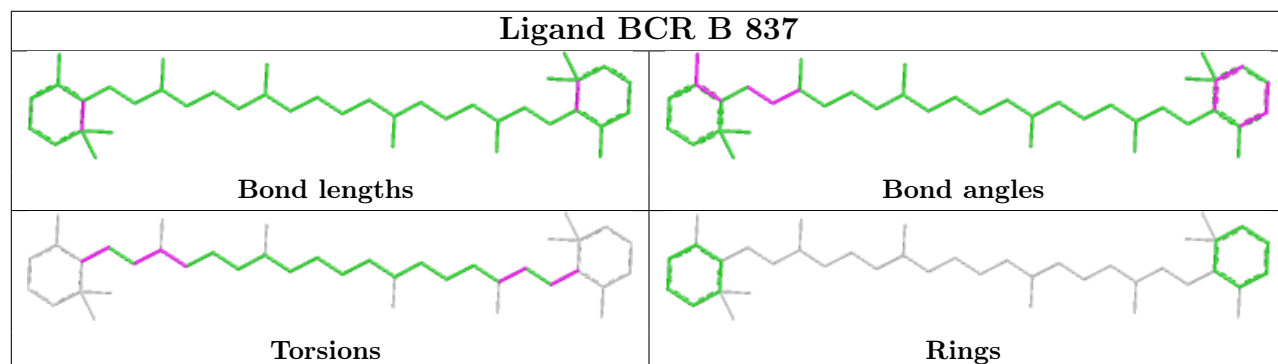


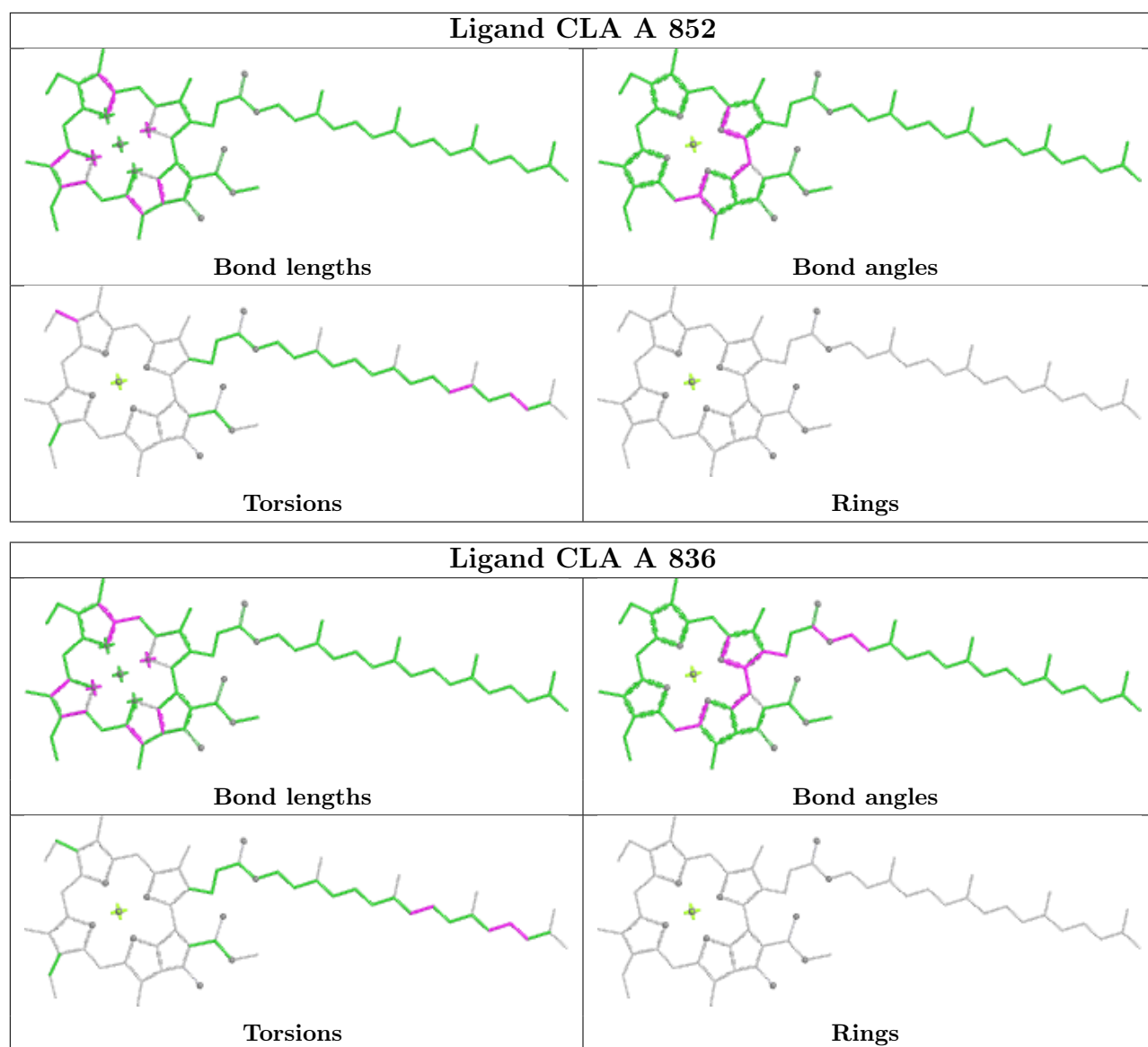


Ligand CLA A 812

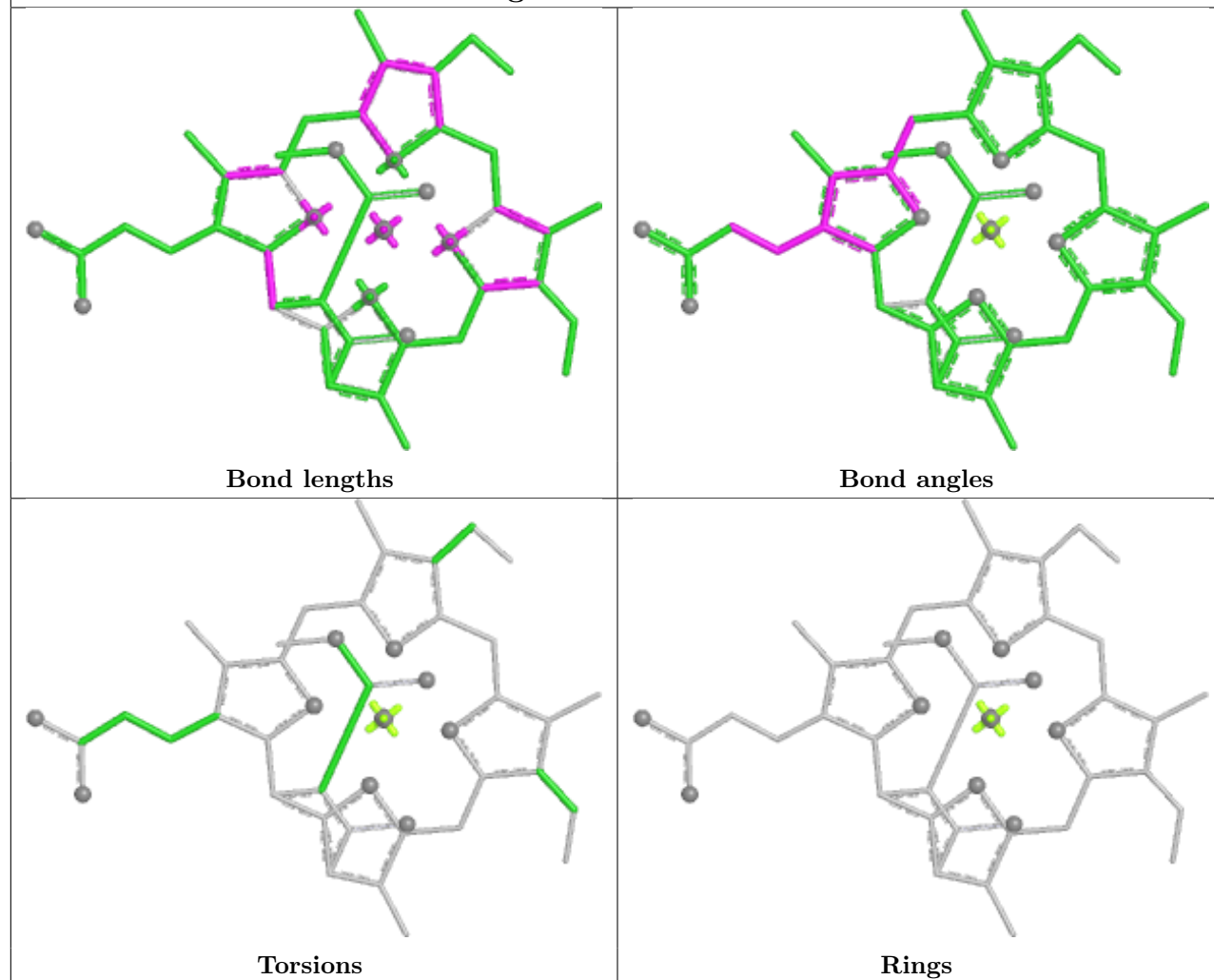


Ligand BCR B 837

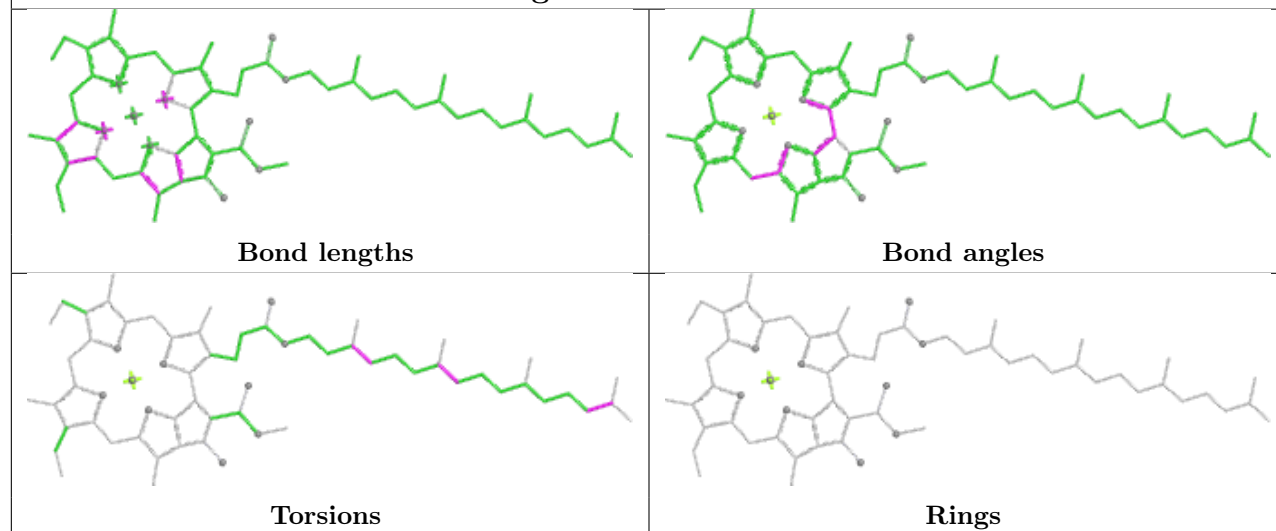


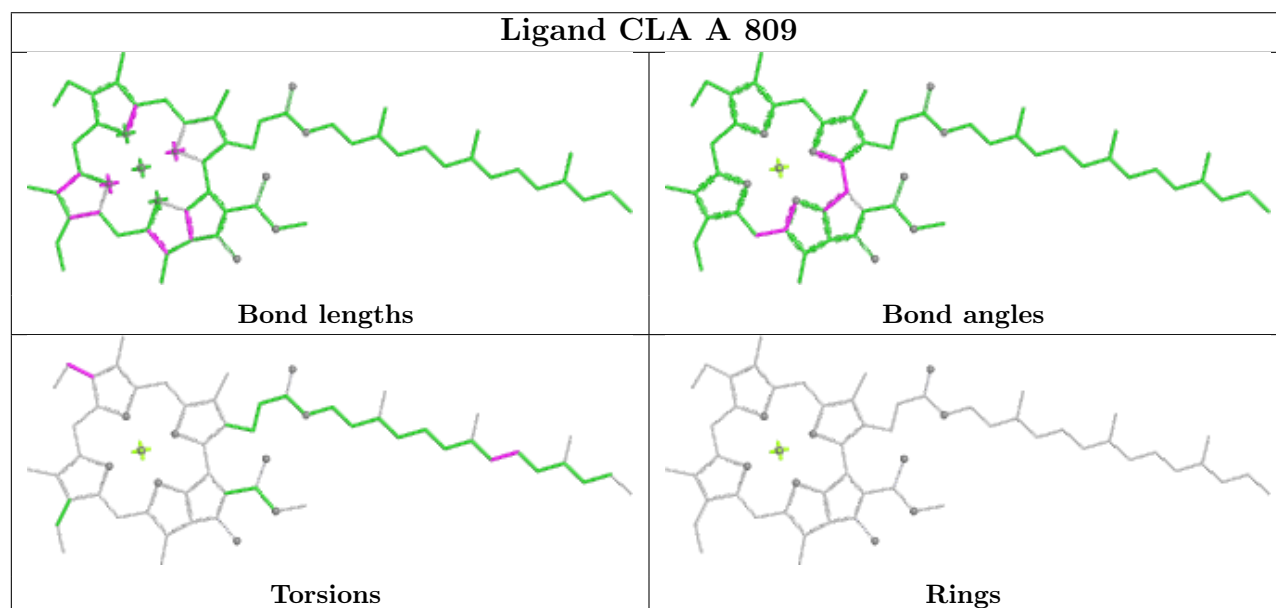
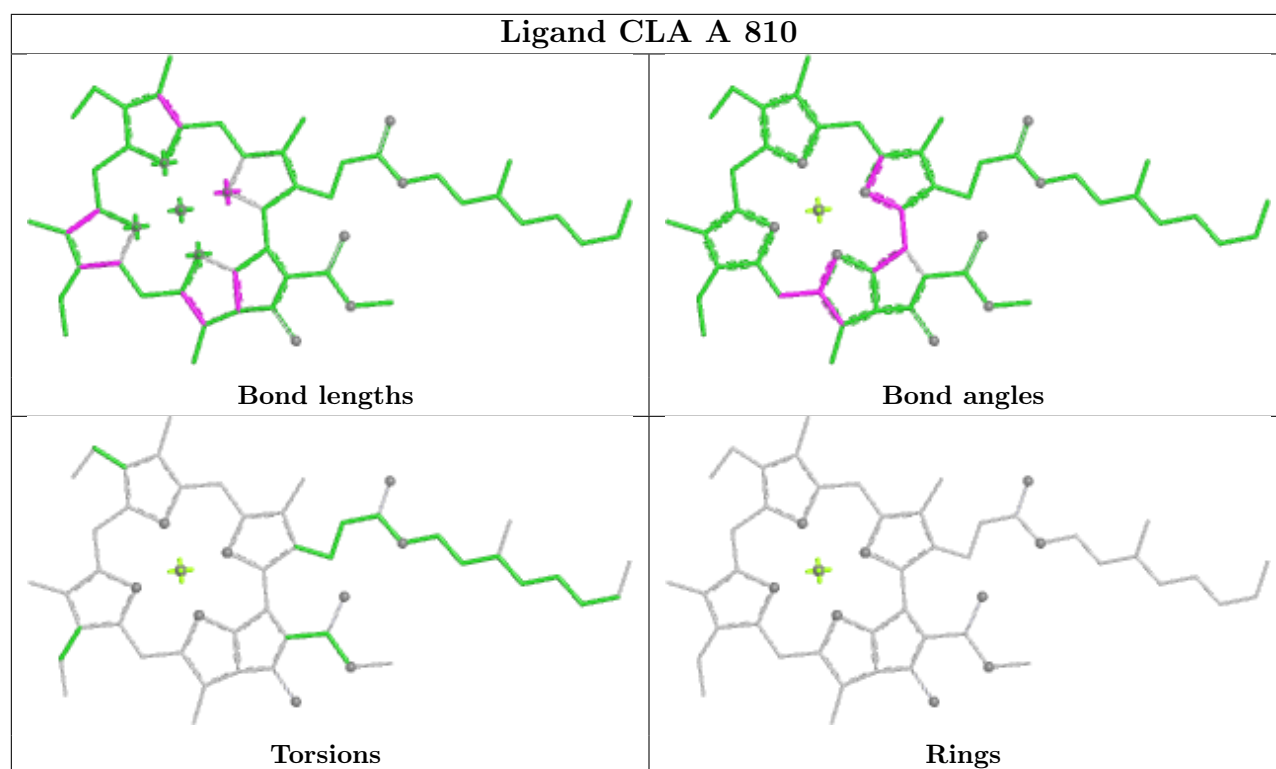


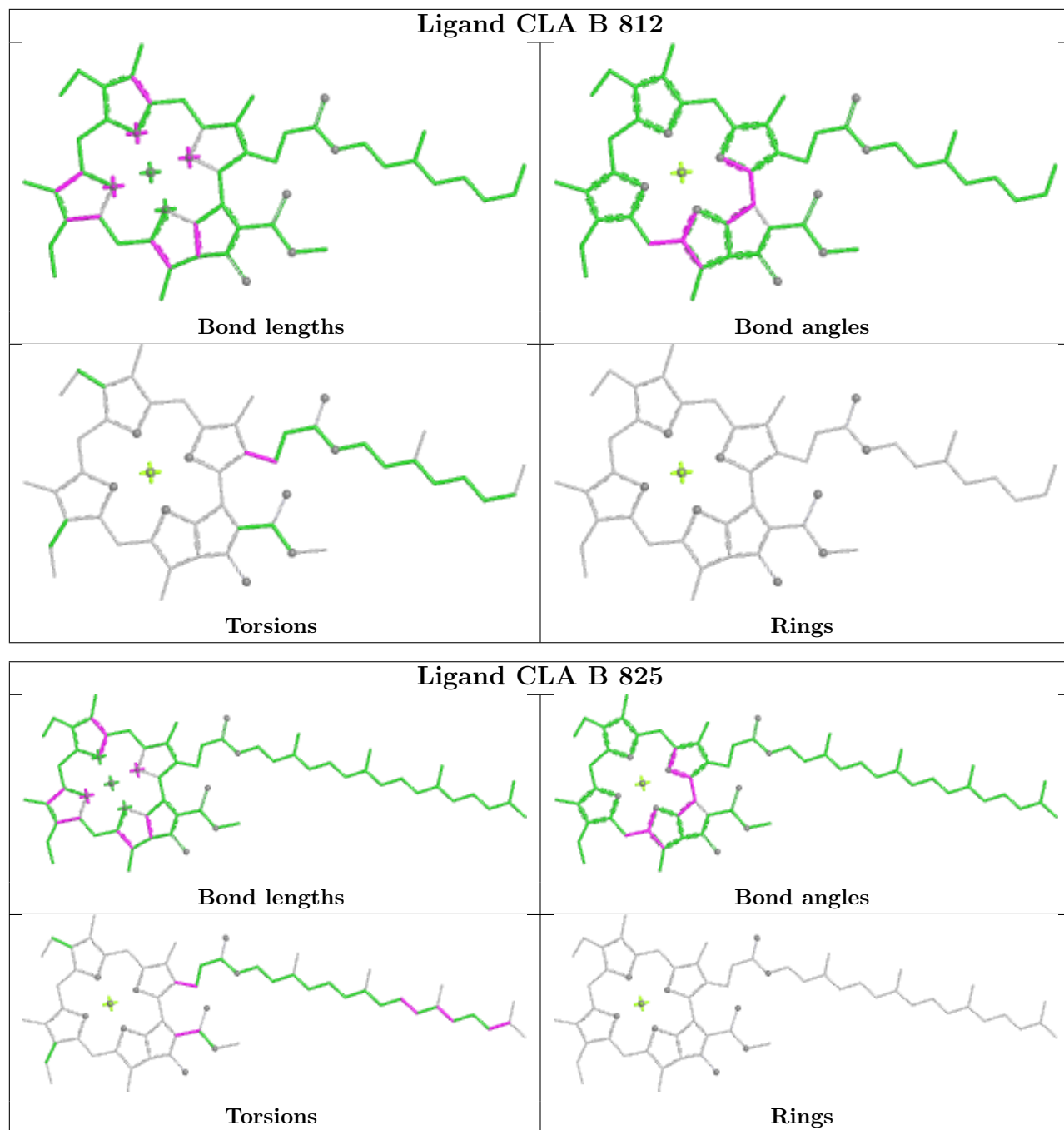
Ligand KC1 O 210

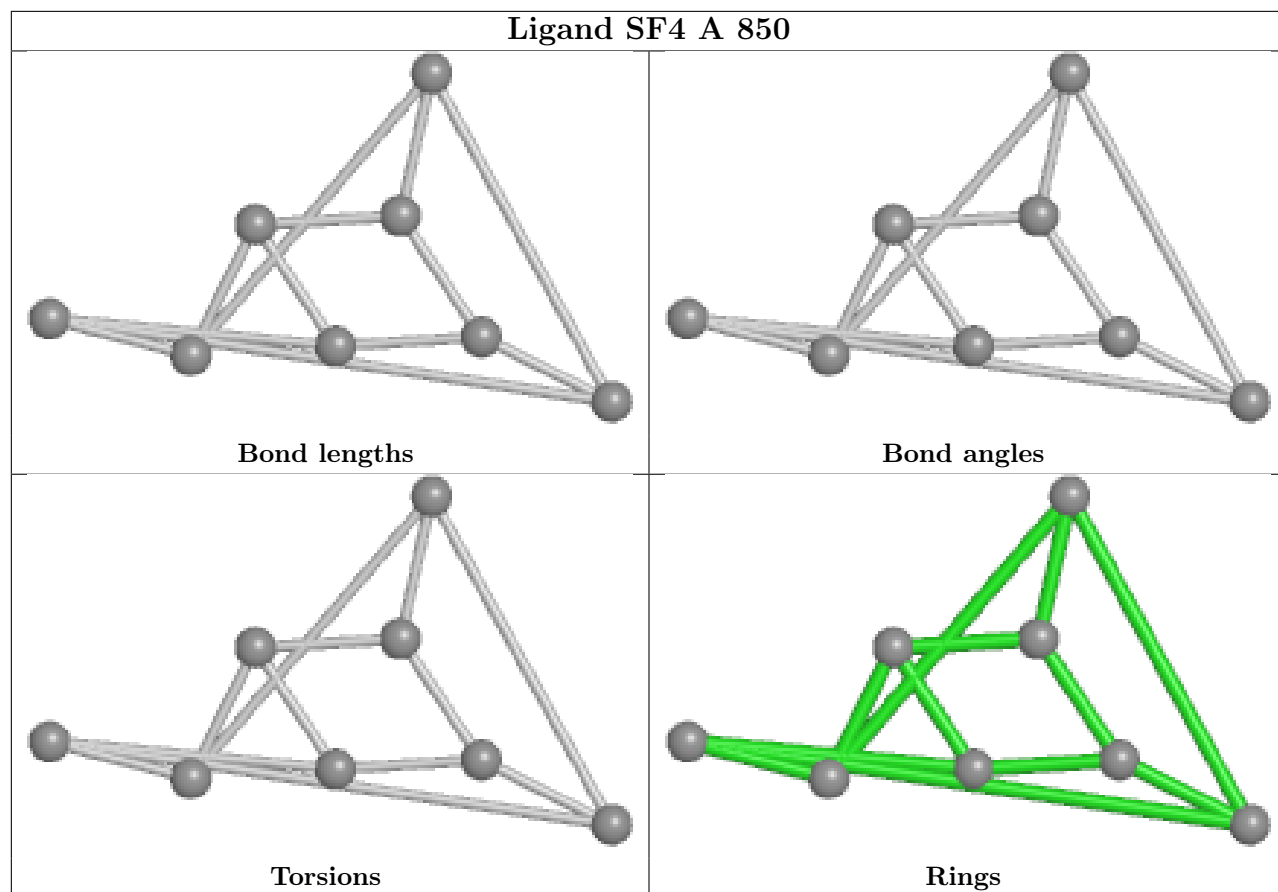


Ligand CLA A 849

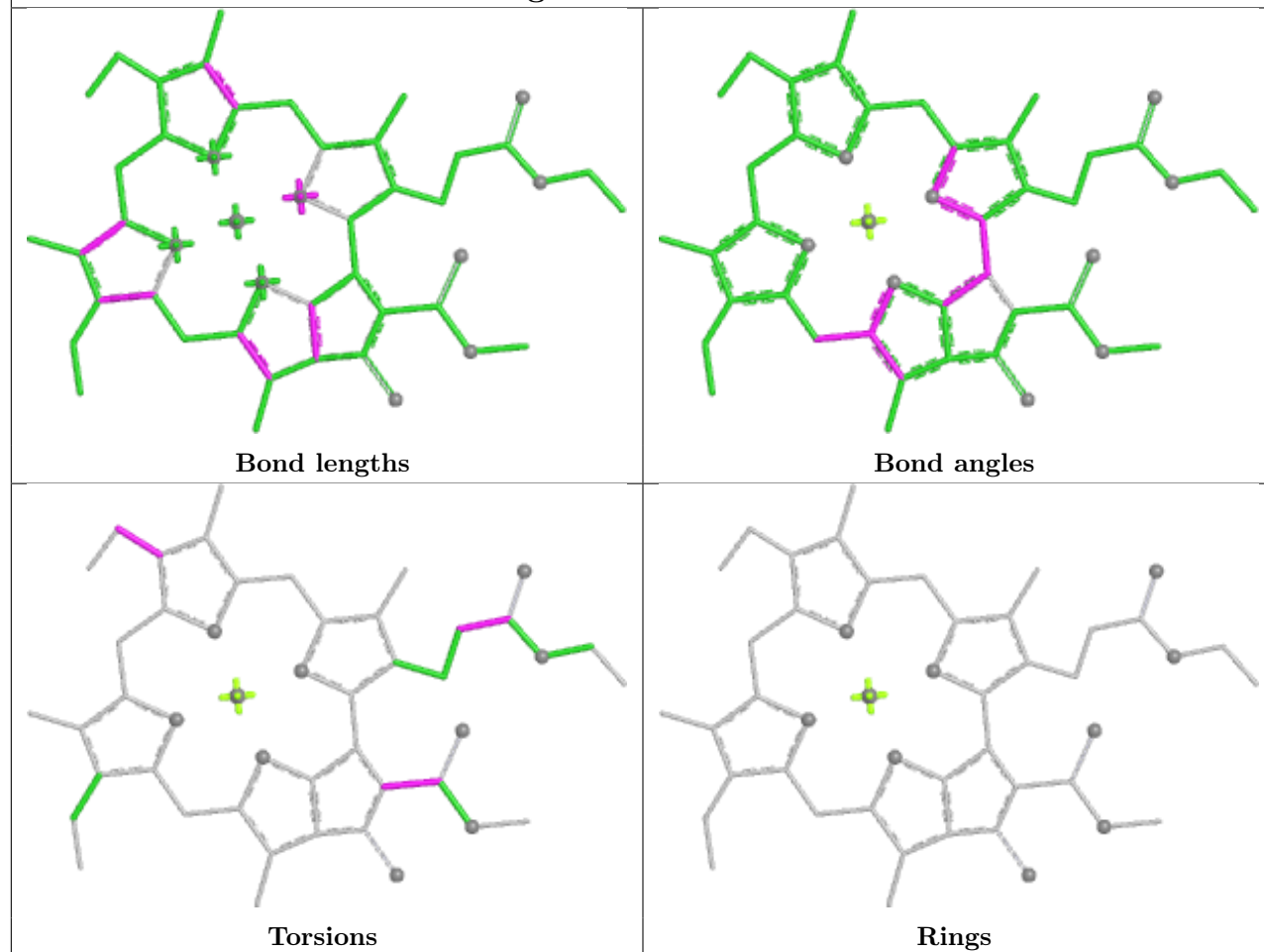




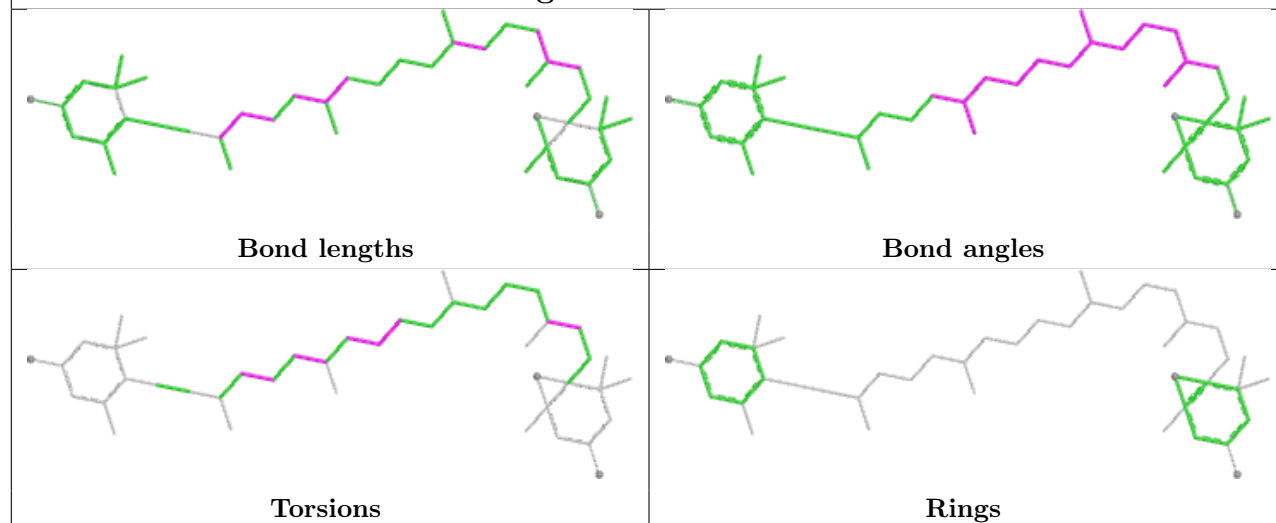


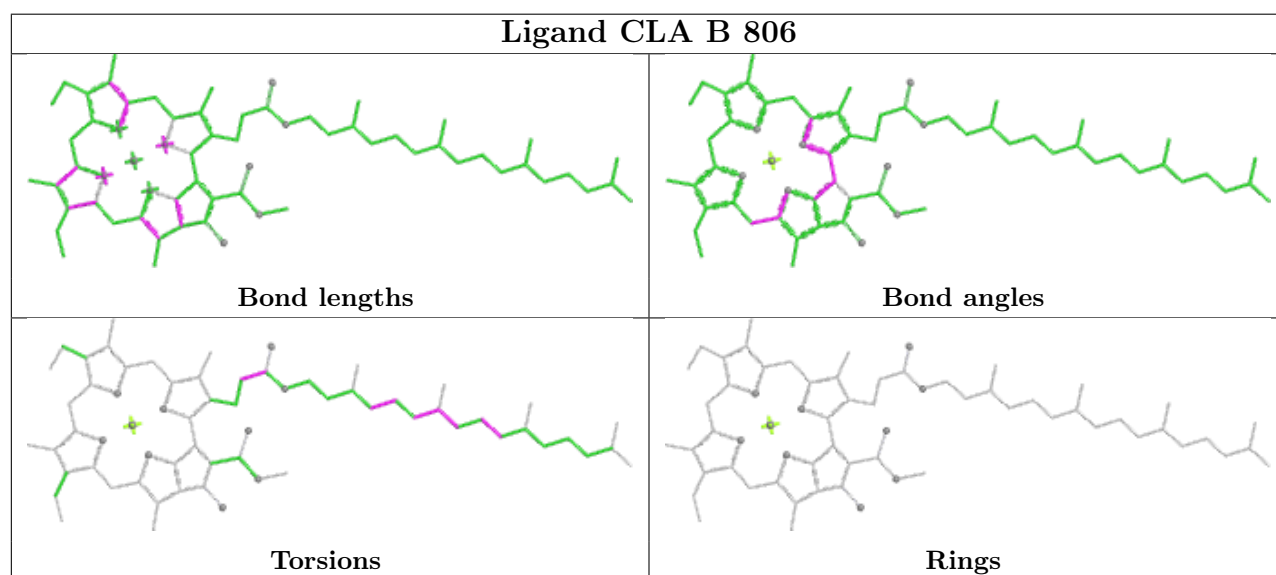
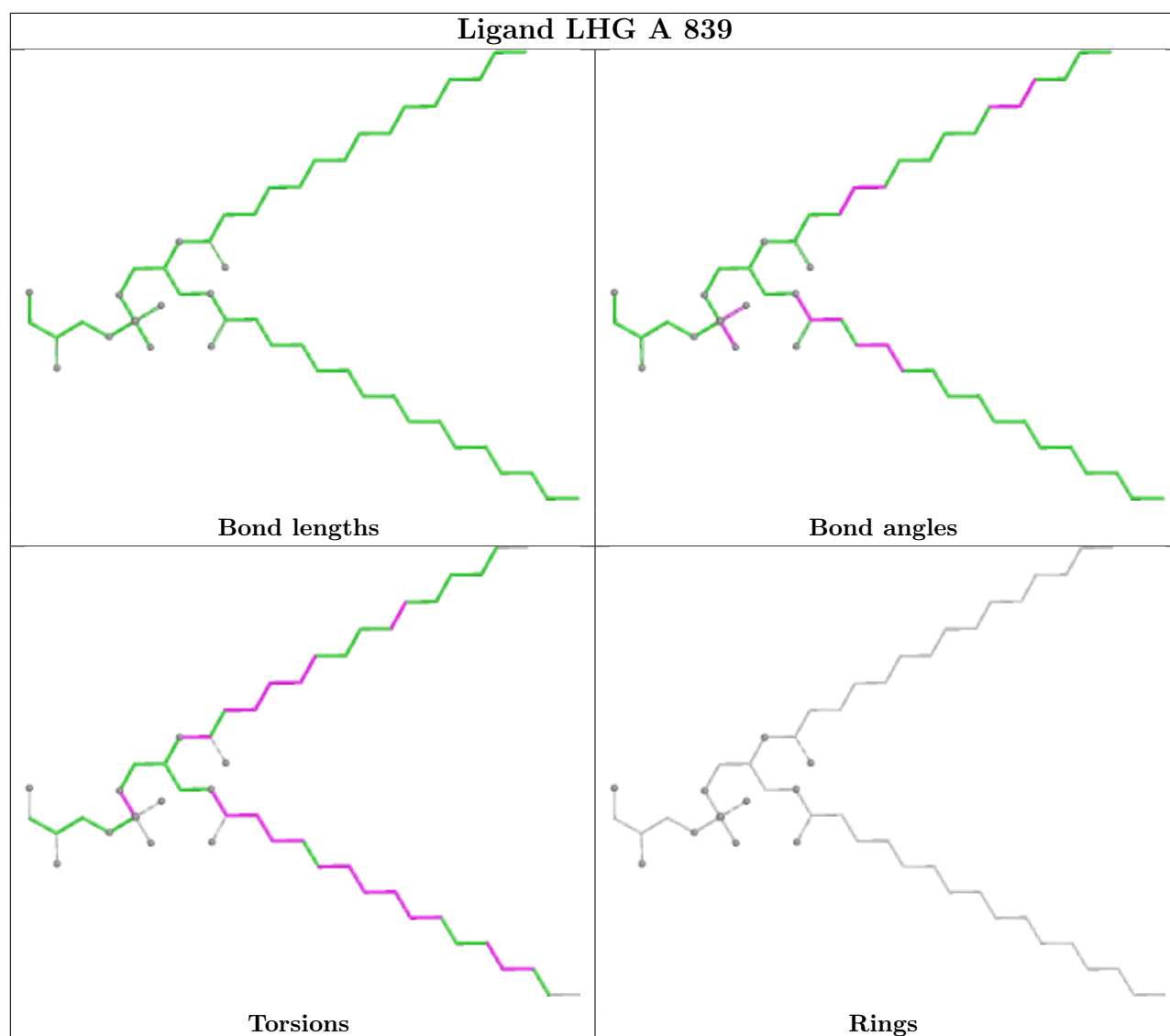


Ligand CLA T 212

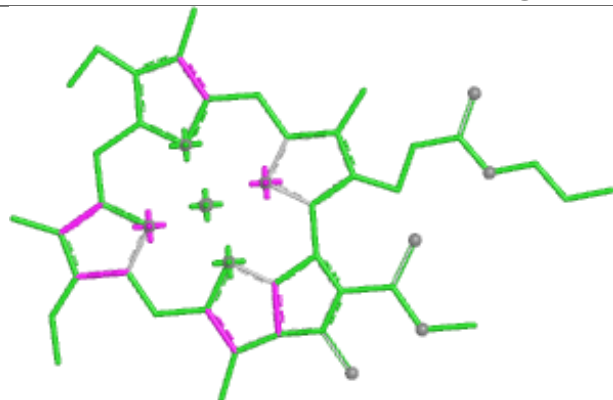


Ligand DD6 J 101

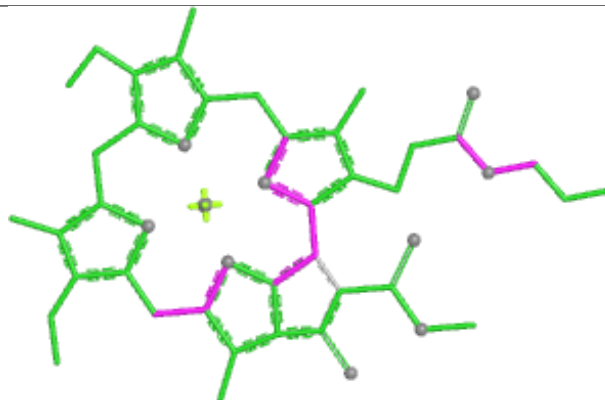




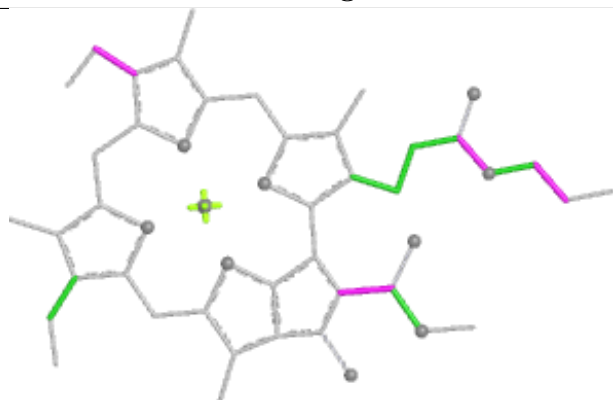
Ligand CLA F 802



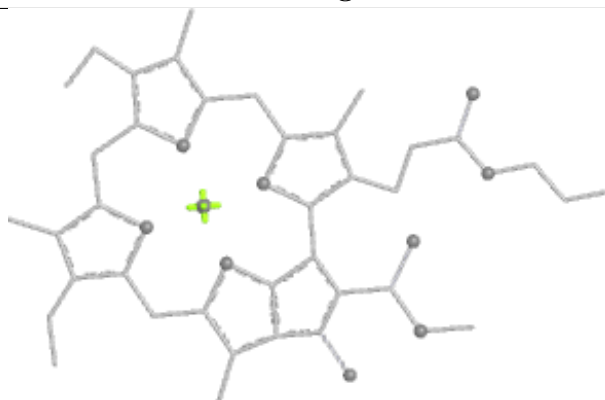
Bond lengths



Bond angles

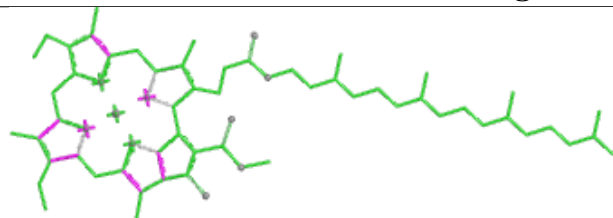


Torsions

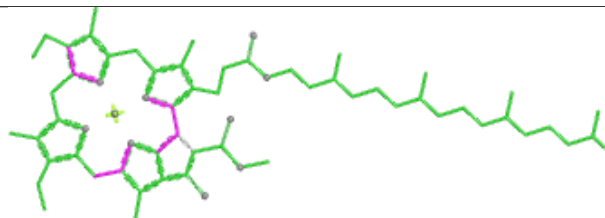


Rings

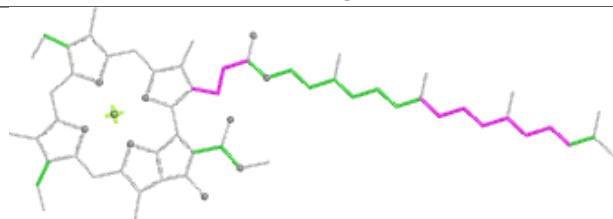
Ligand CLA A 815



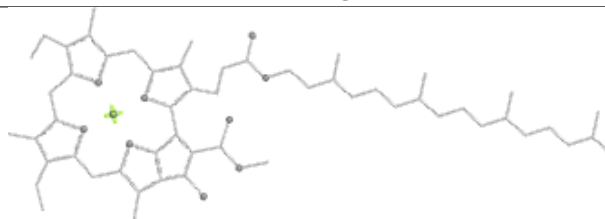
Bond lengths



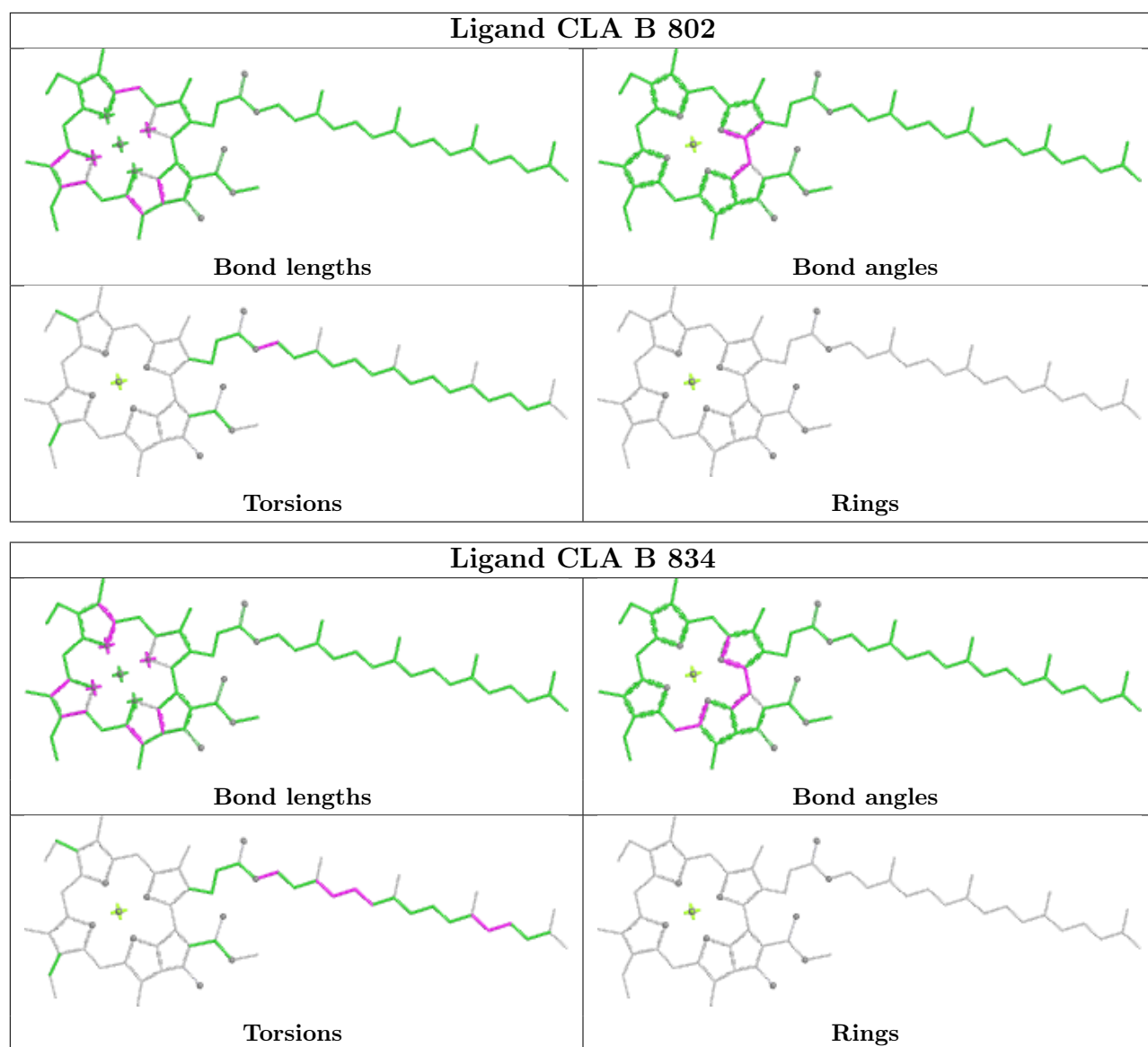
Bond angles



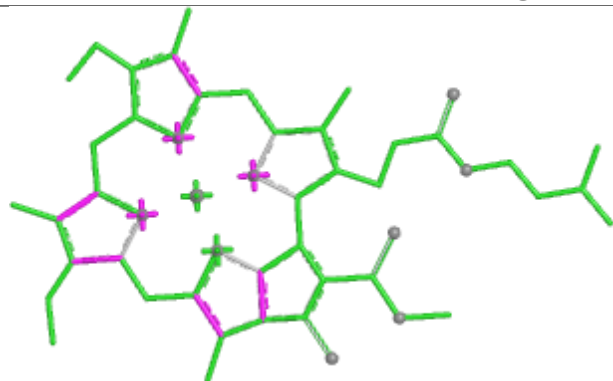
Torsions



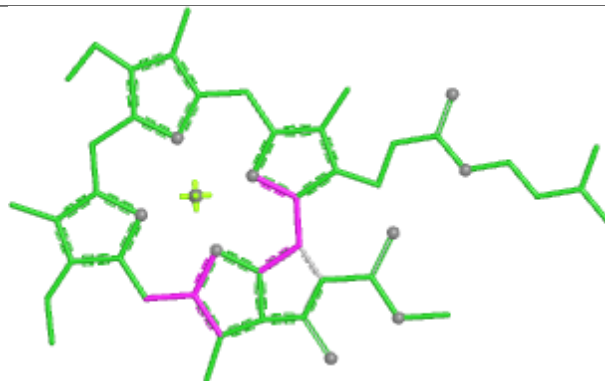
Rings



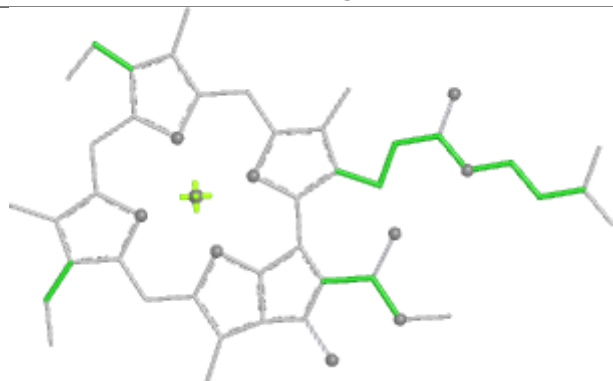
Ligand CLA B 826



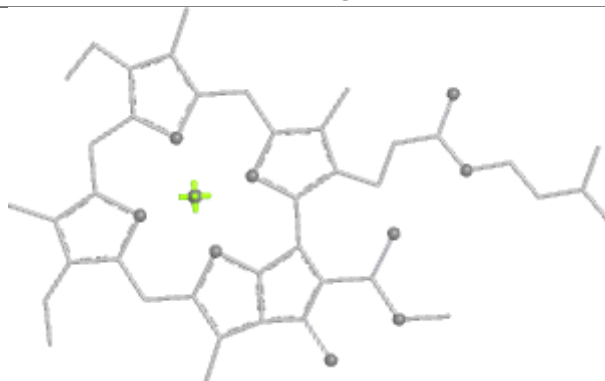
Bond lengths



Bond angles

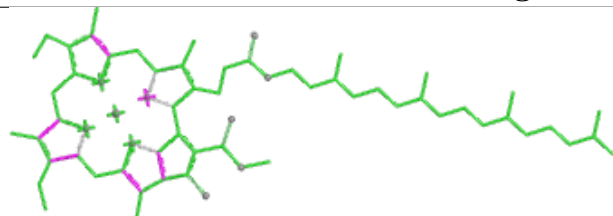


Torsions

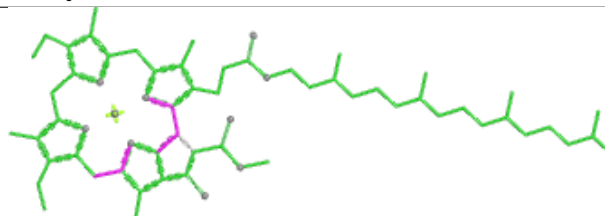


Rings

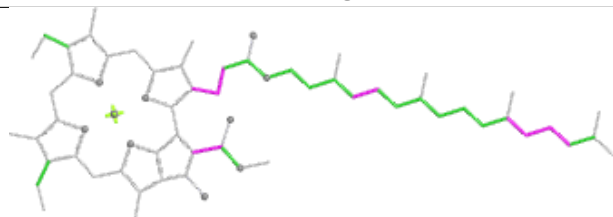
Ligand CLA Q 209



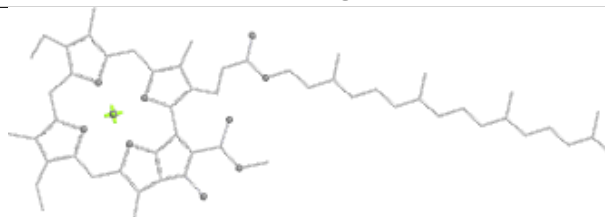
Bond lengths



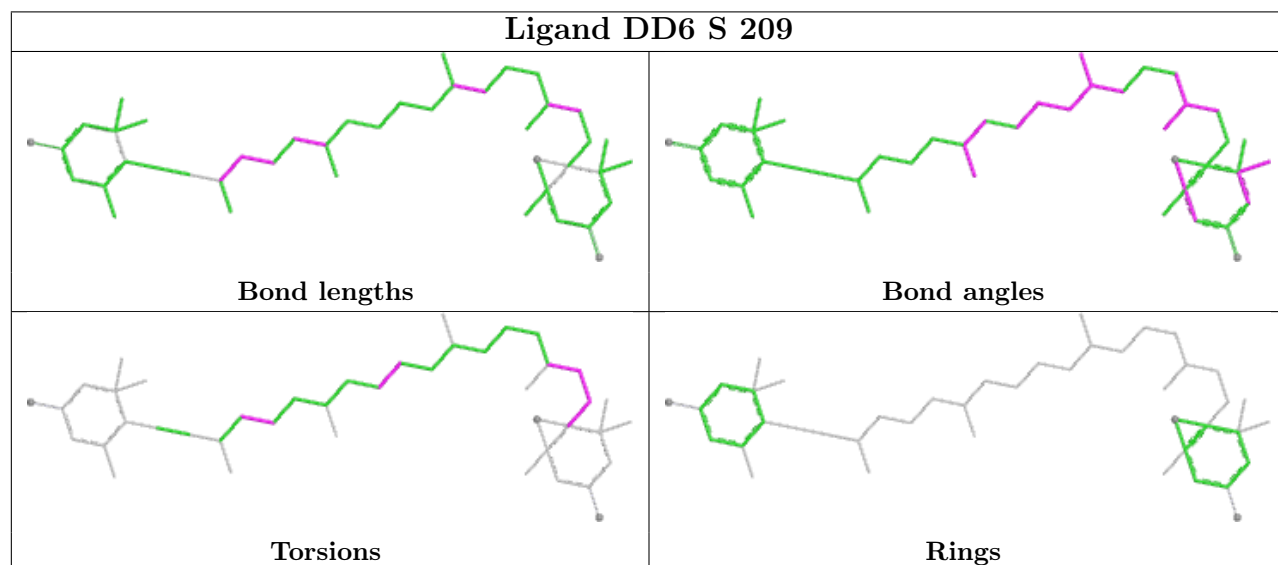
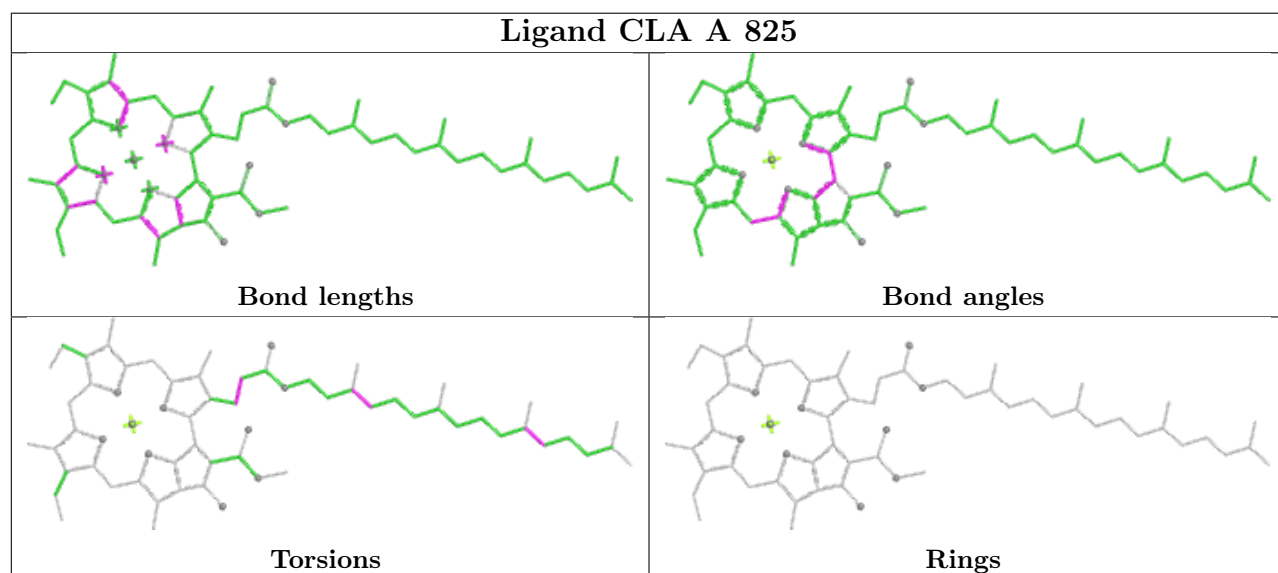
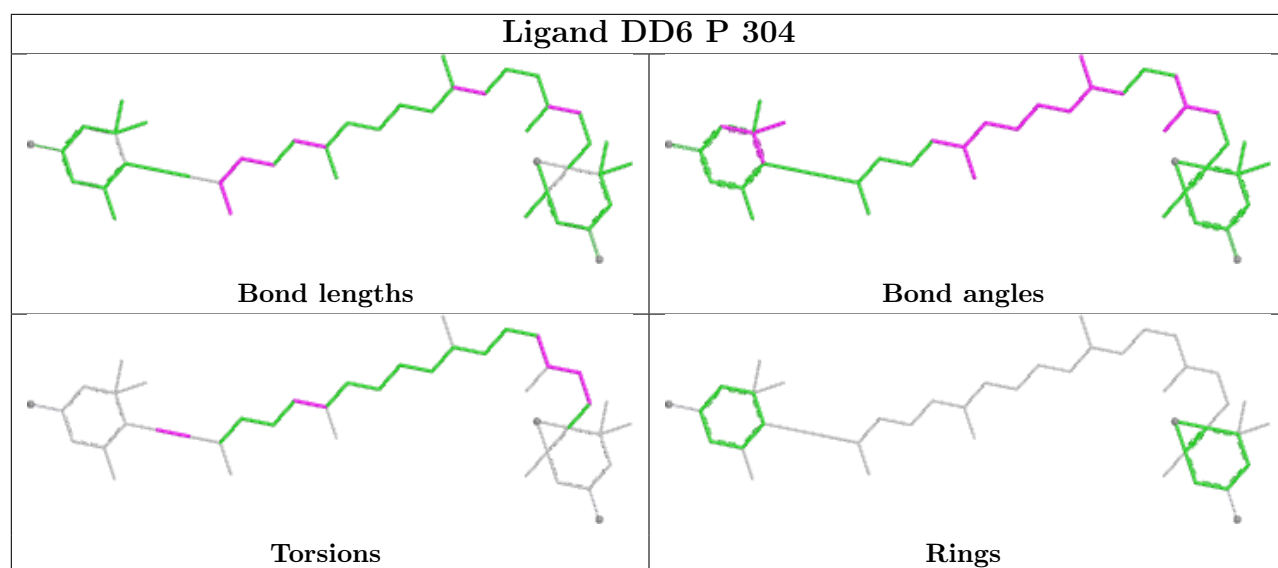
Bond angles



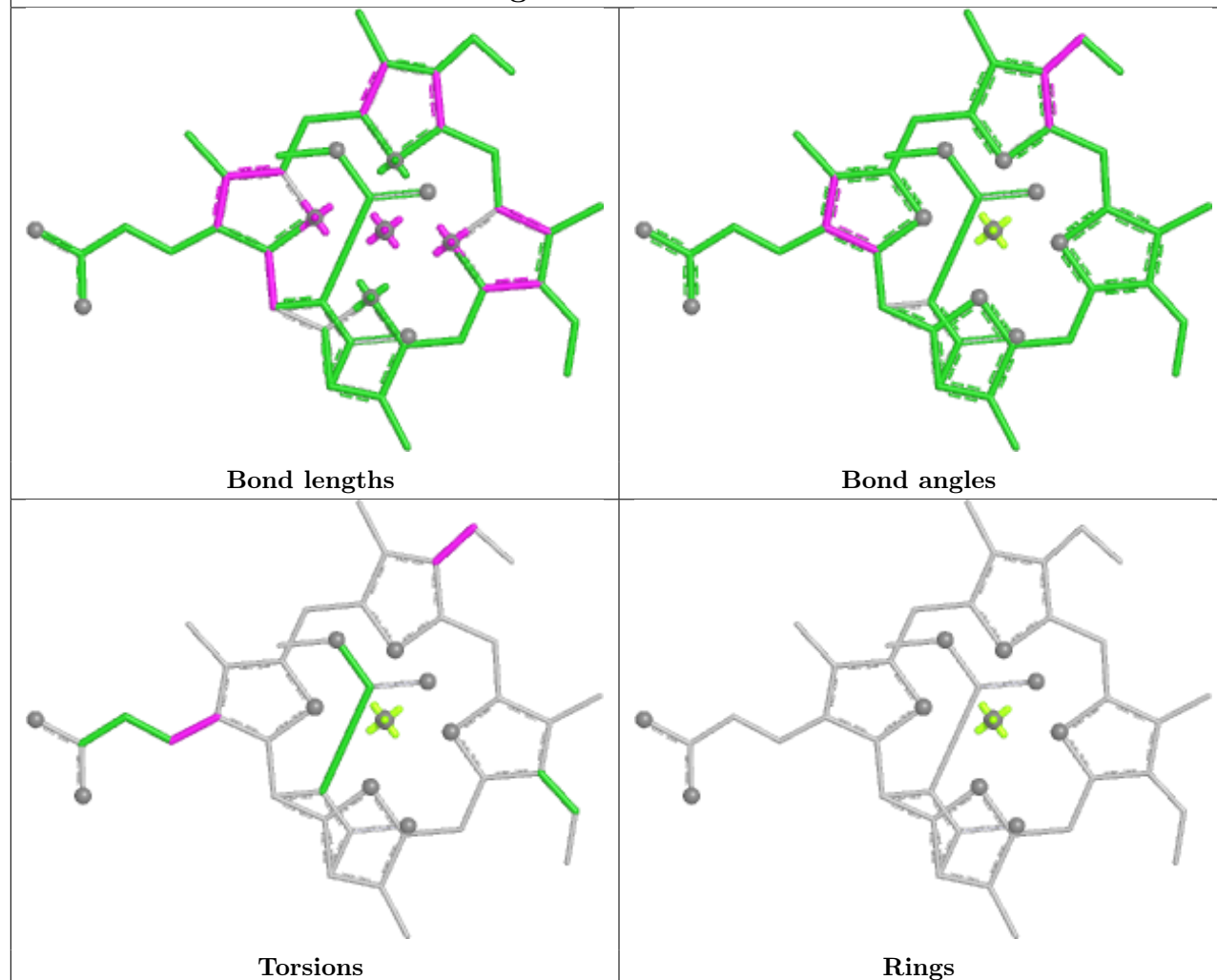
Torsions



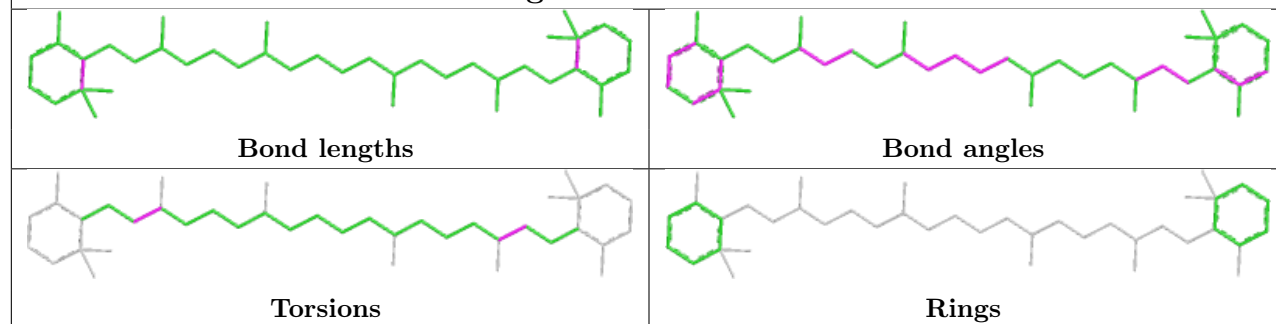
Rings



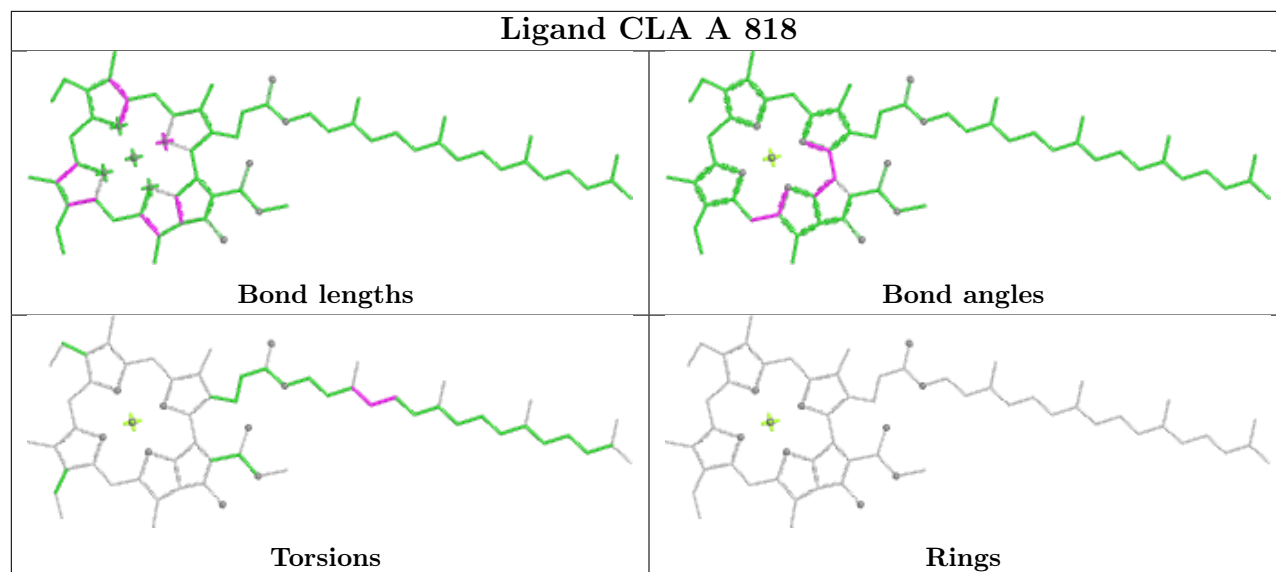
Ligand KC1 T 209



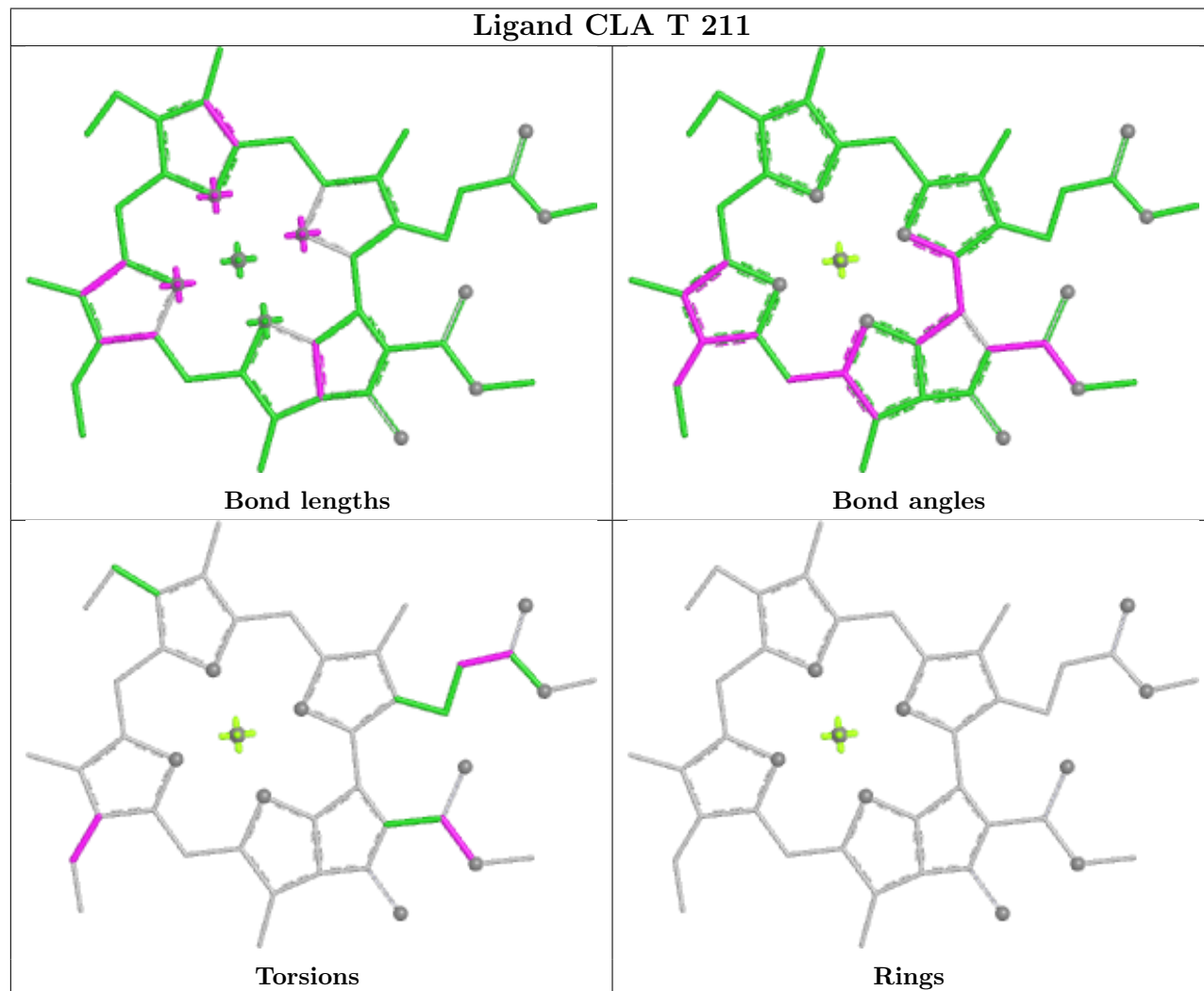
Ligand BCR M 101

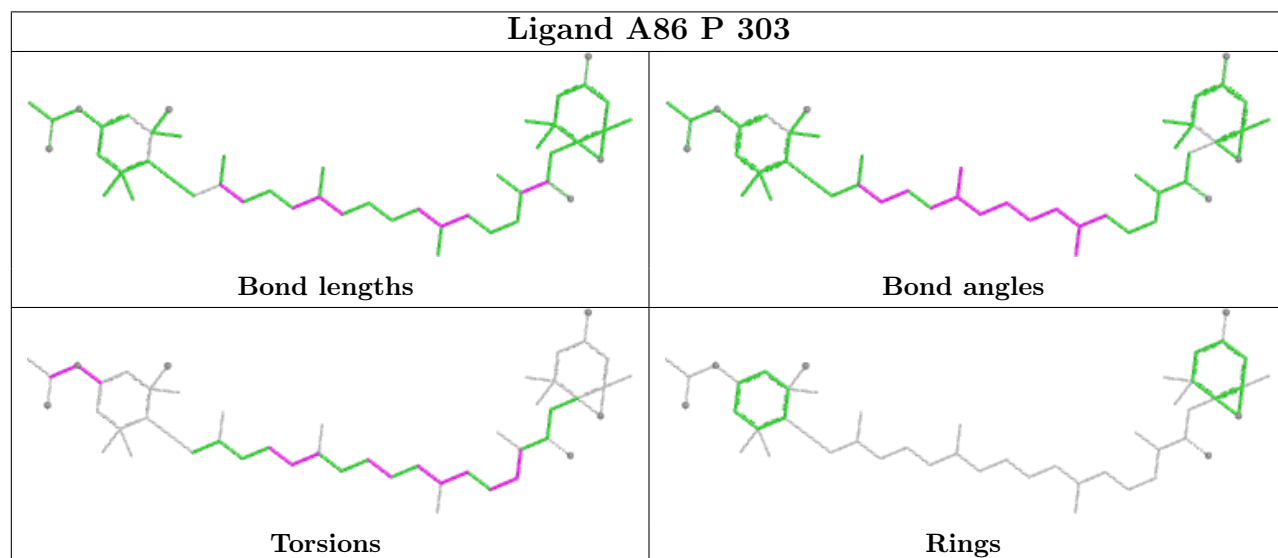
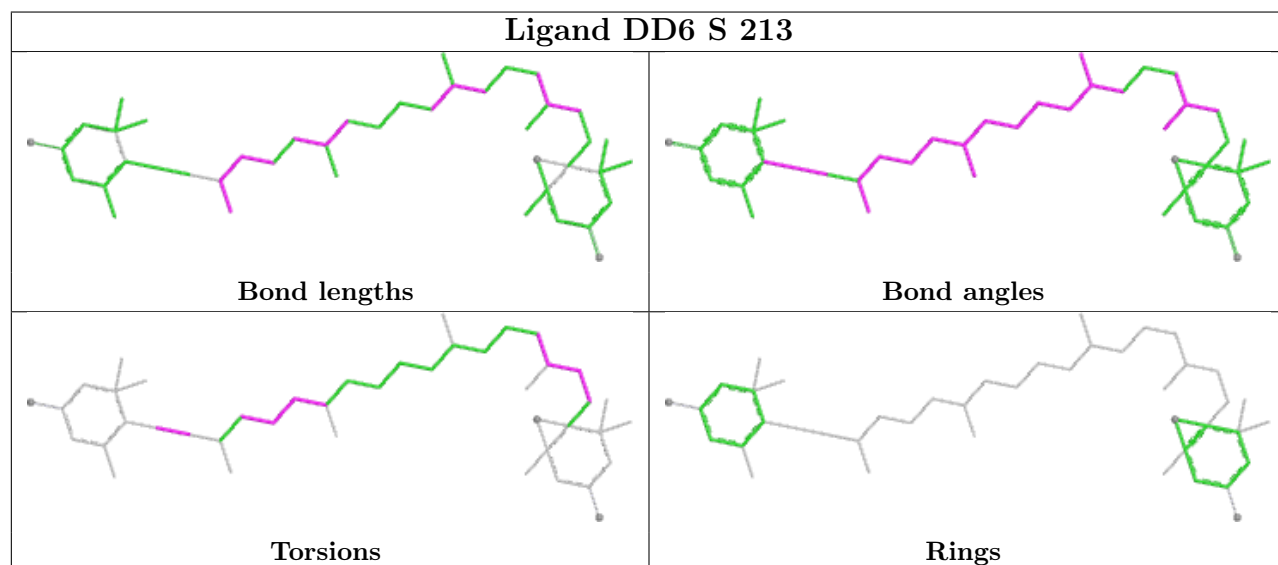
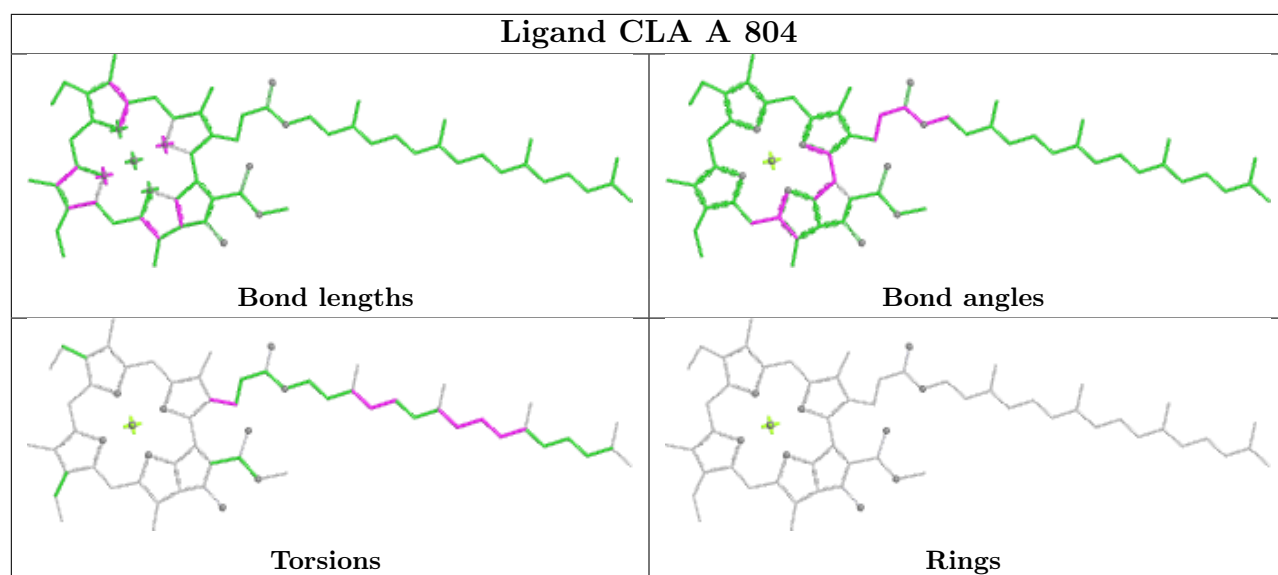


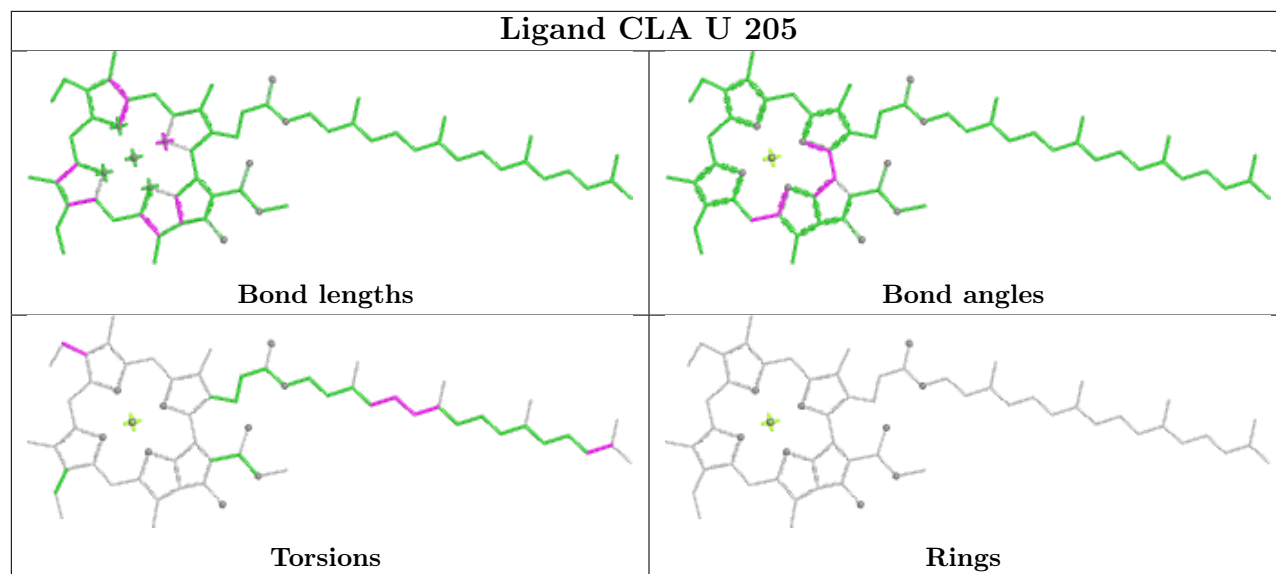
Ligand CLA A 818



Ligand CLA T 211







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.

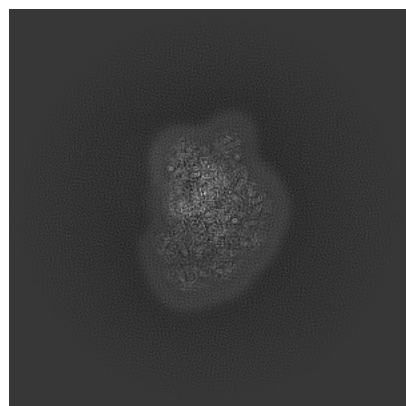
6 Map visualisation [i](#)

This section contains visualisations of the EMDB entry EMD-64378. These allow visual inspection of the internal detail of the map and identification of artifacts.

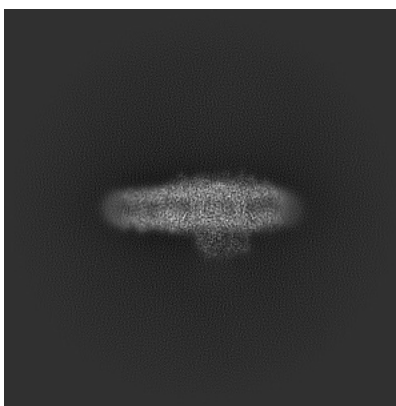
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections [i](#)

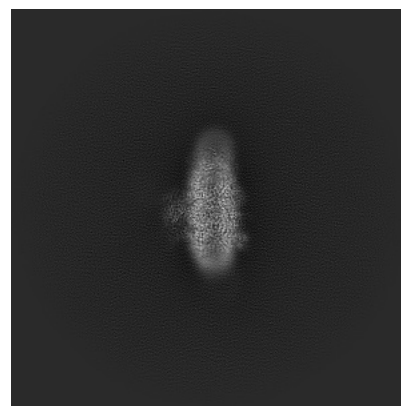
6.1.1 Primary map



X

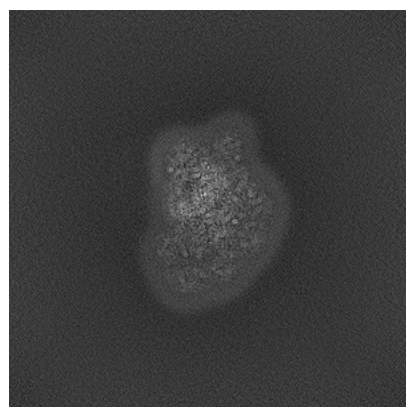


Y

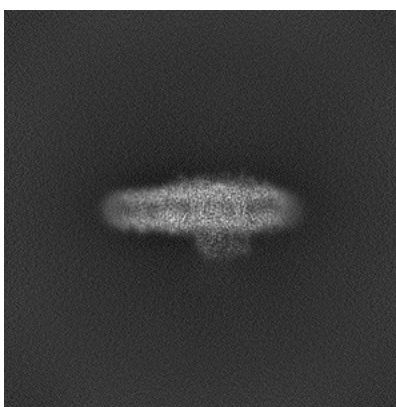


Z

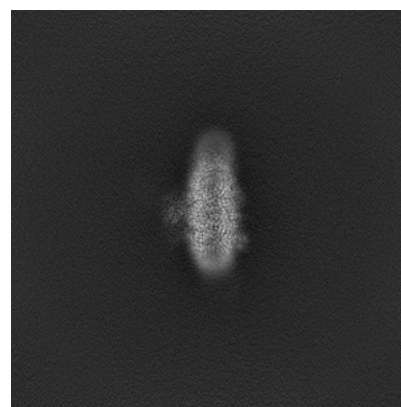
6.1.2 Raw map



X



Y

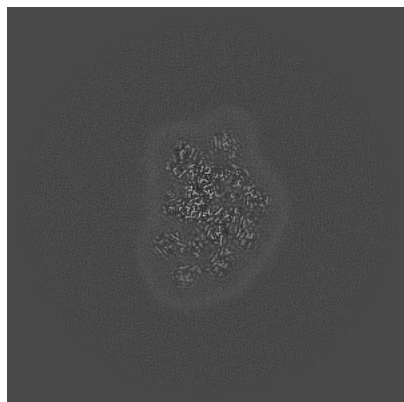


Z

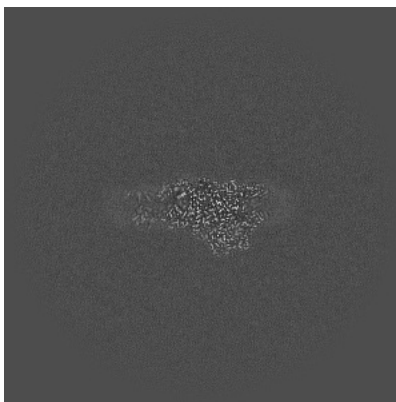
The images above show the map projected in three orthogonal directions.

6.2 Central slices [i](#)

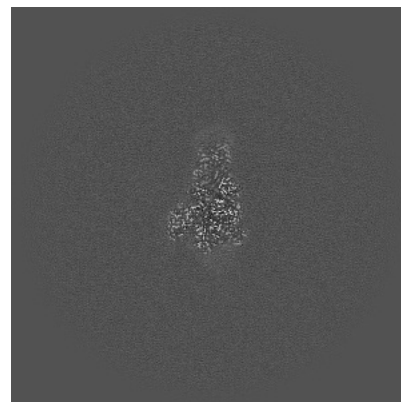
6.2.1 Primary map



X Index: 300

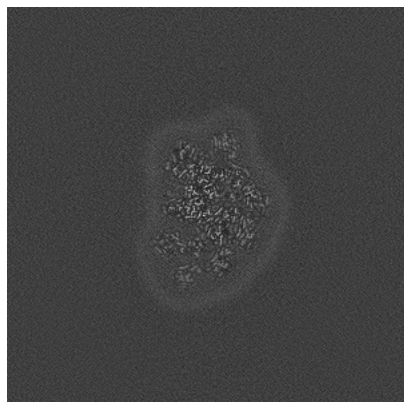


Y Index: 300

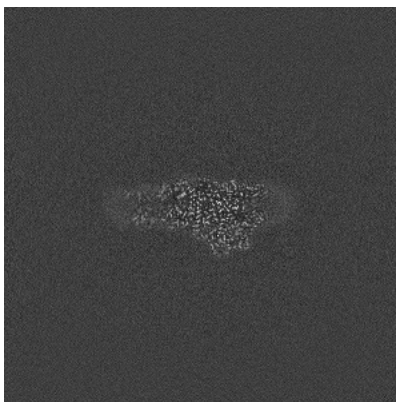


Z Index: 300

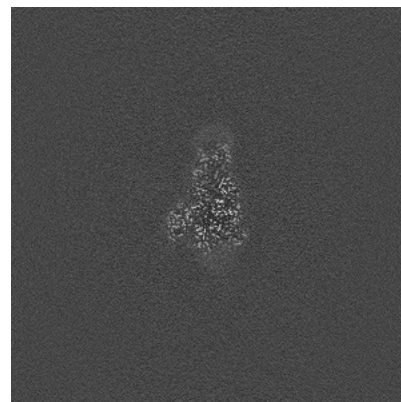
6.2.2 Raw map



X Index: 300



Y Index: 300

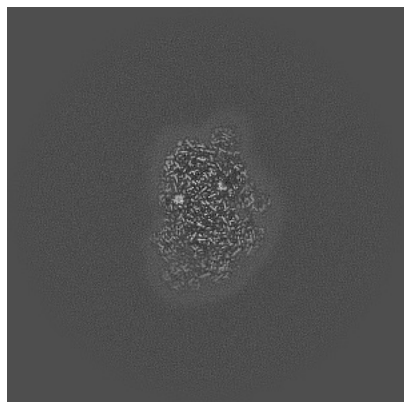


Z Index: 300

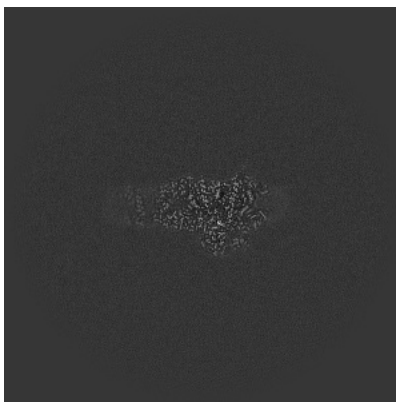
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices [i](#)

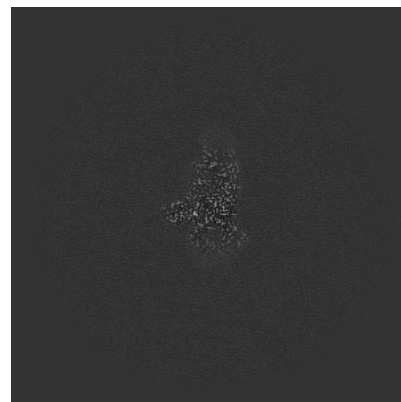
6.3.1 Primary map



X Index: 284

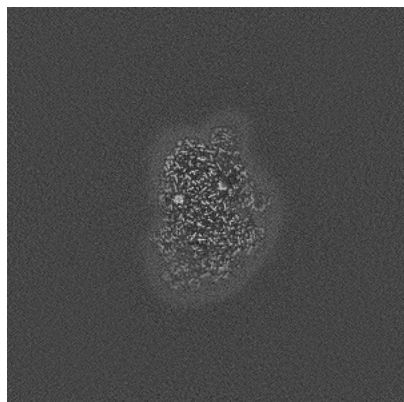


Y Index: 289

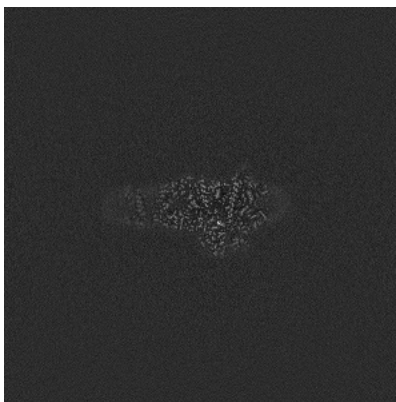


Z Index: 320

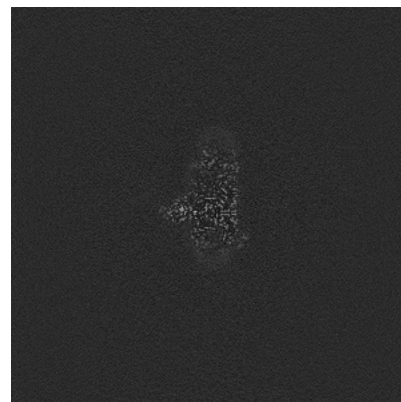
6.3.2 Raw map



X Index: 284



Y Index: 289

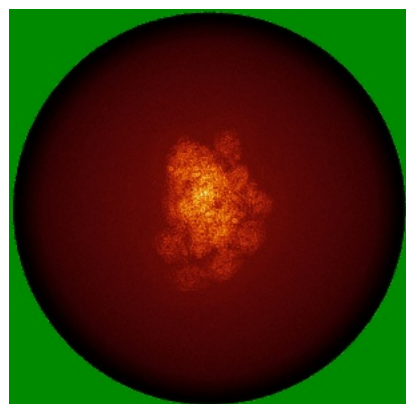


Z Index: 322

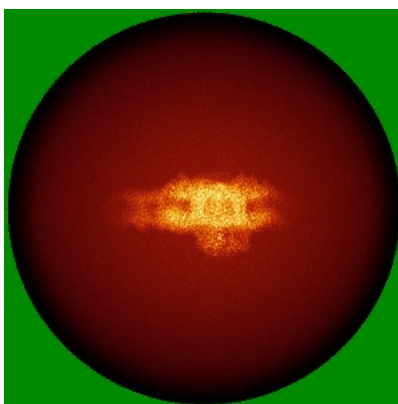
The images above show the largest variance slices of the map in three orthogonal directions.

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

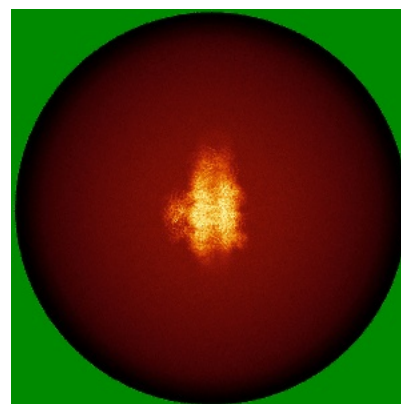
6.4.1 Primary map



X

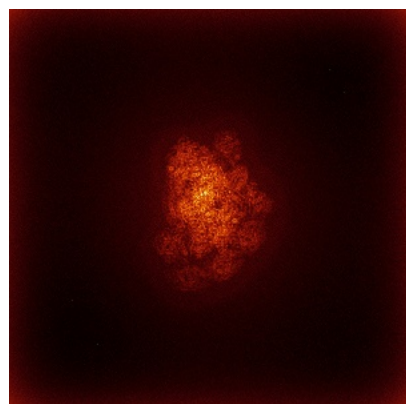


Y

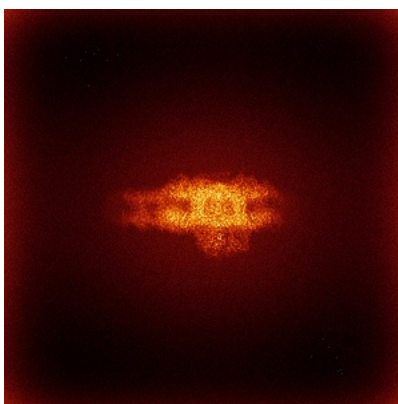


Z

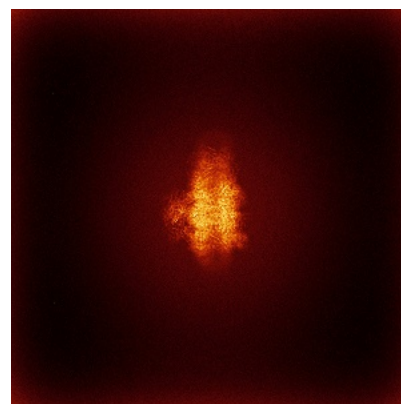
6.4.2 Raw map



X



Y

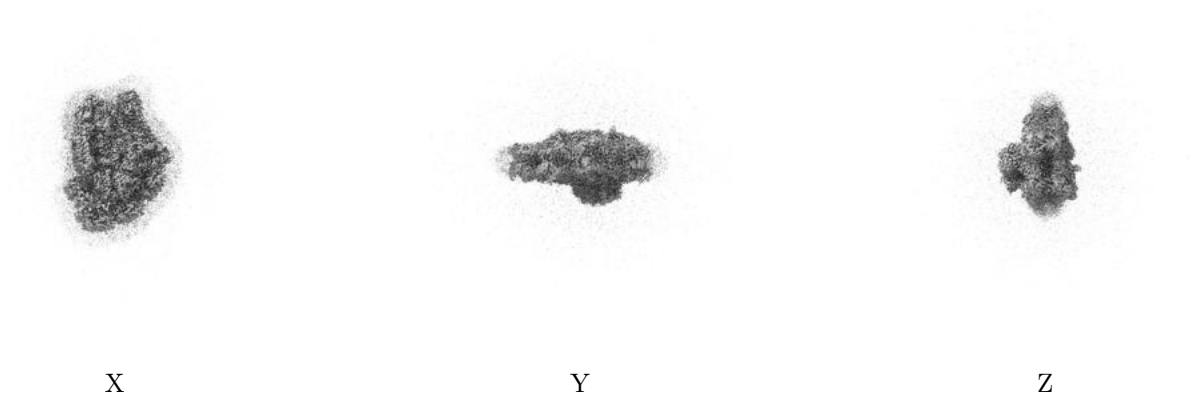


Z

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

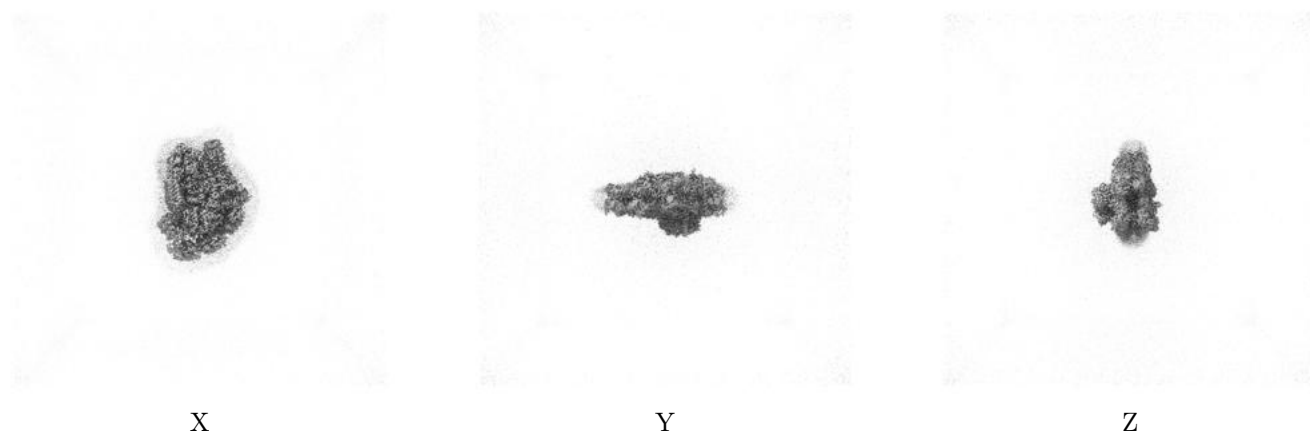
6.5 Orthogonal surface views [i](#)

6.5.1 Primary map



The images above show the 3D surface view of the map at the recommended contour level 0.08. These images, in conjunction with the slice images, may facilitate assessment of whether an appropriate contour level has been provided.

6.5.2 Raw map



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

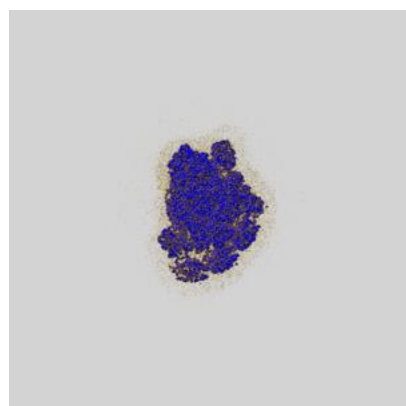
6.6 Mask visualisation [i](#)

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

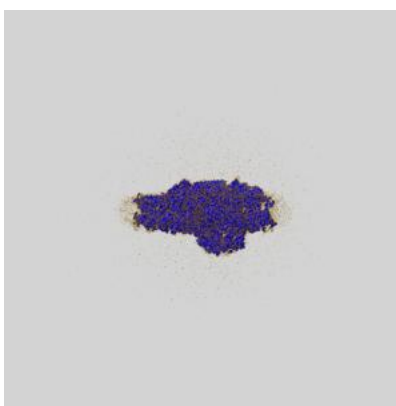
A mask typically either:

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

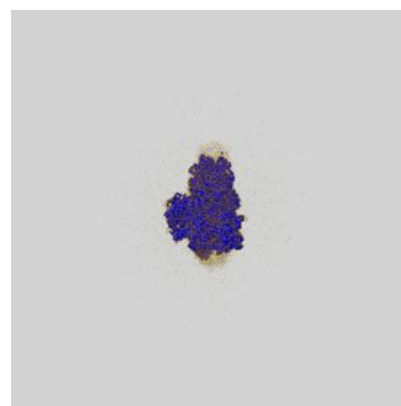
6.6.1 emd_64378_msk_1.map [i](#)



X



Y

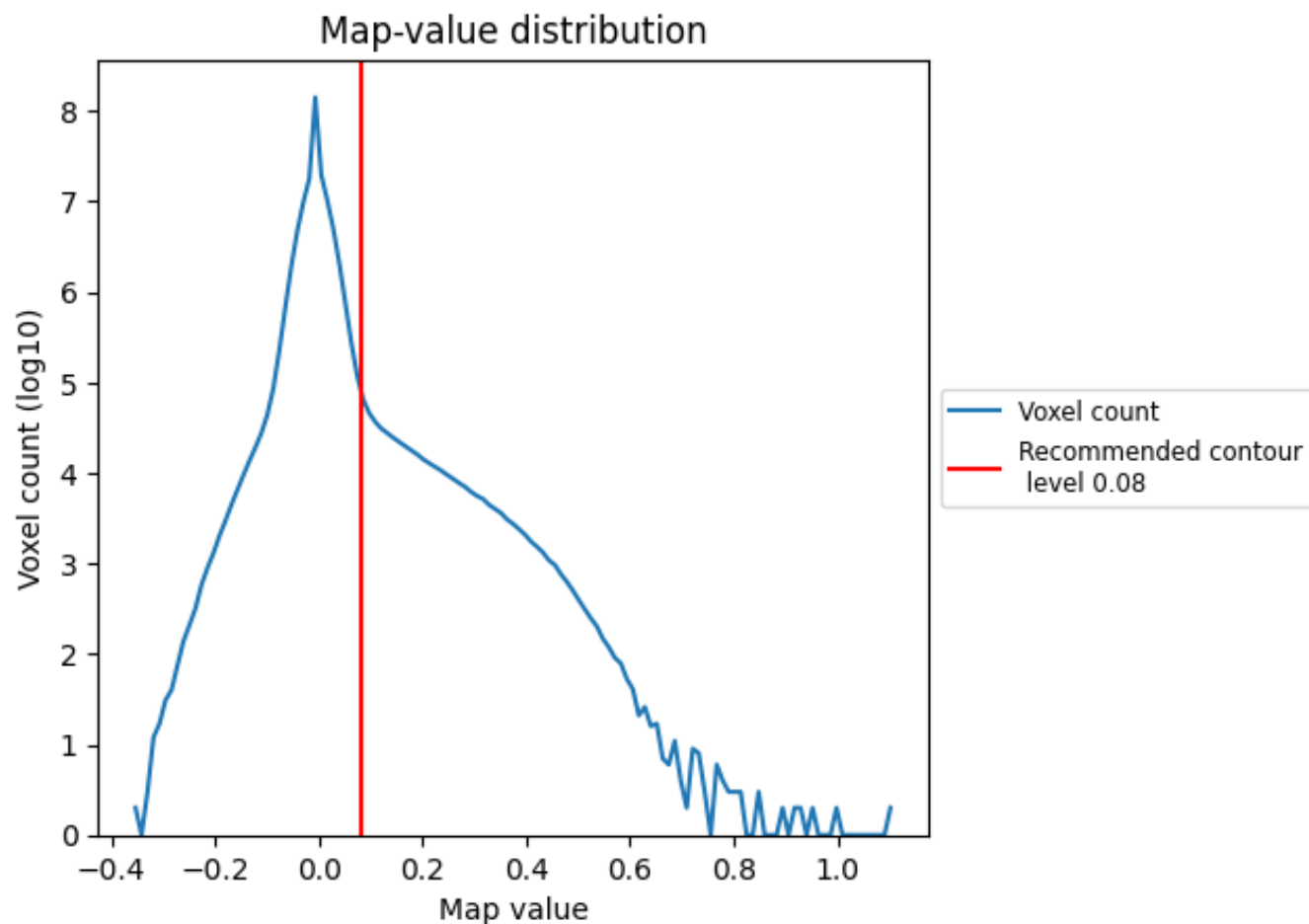


Z

7 Map analysis [i](#)

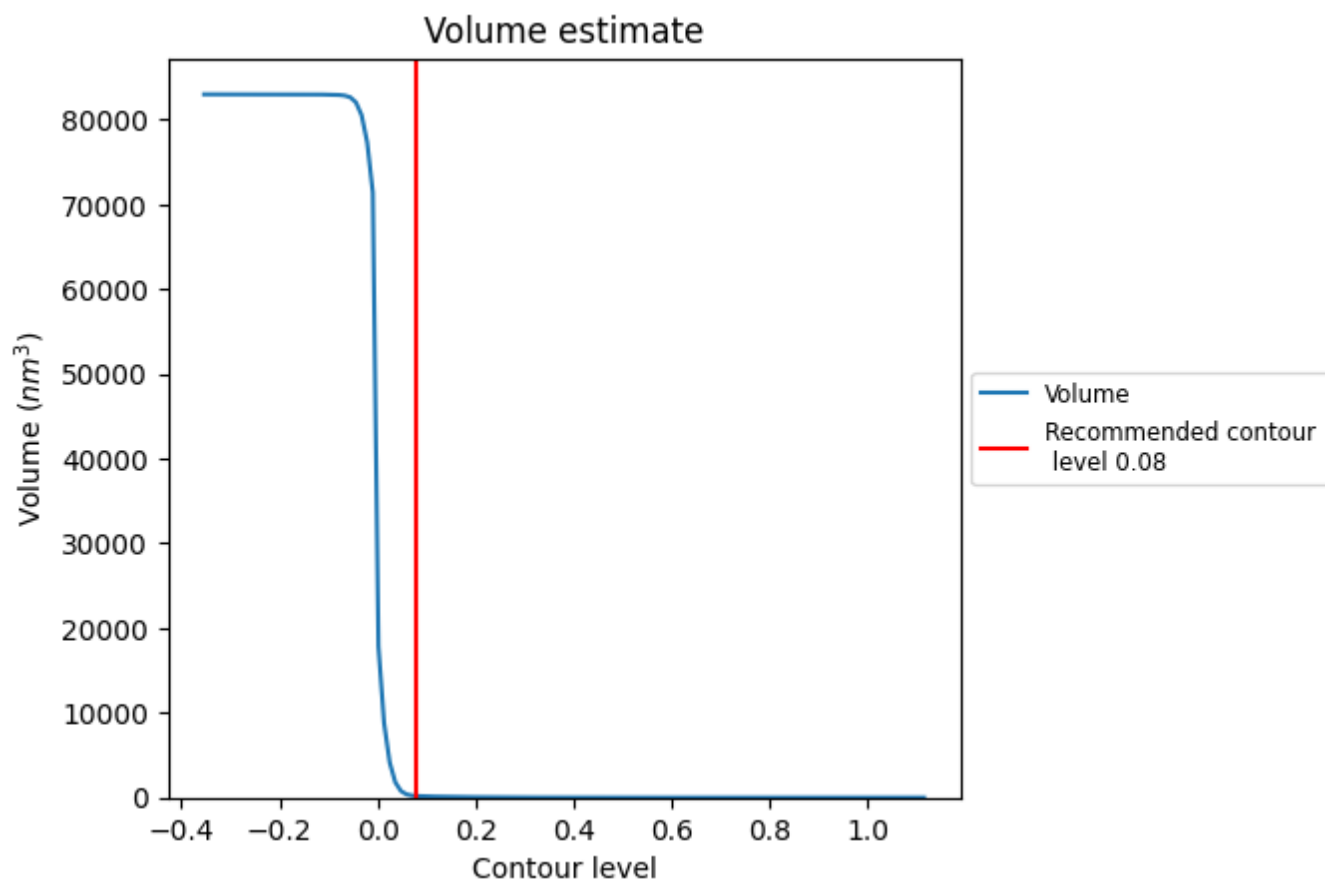
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution [i](#)



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

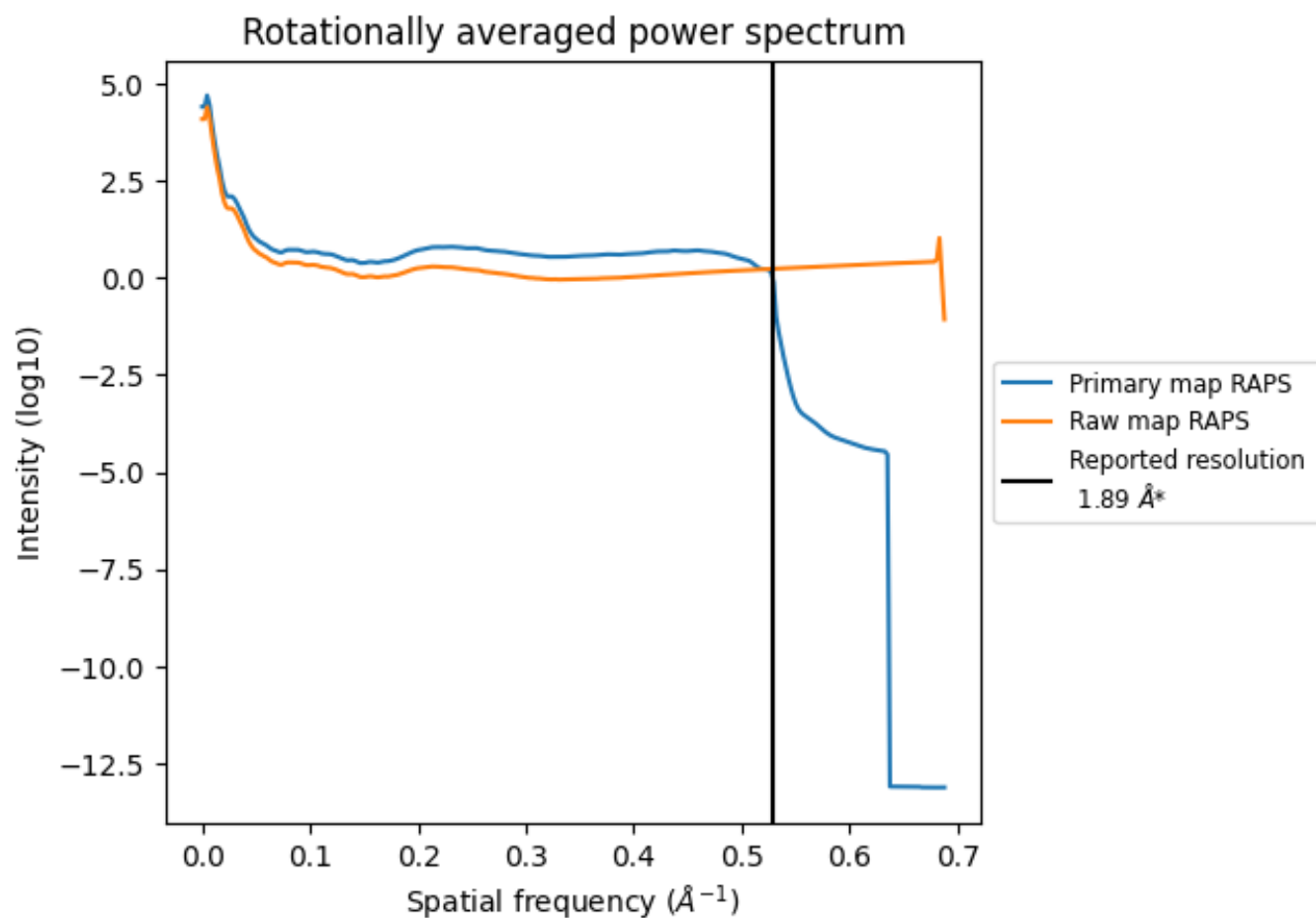
7.2 Volume estimate [i](#)



The volume at the recommended contour level is 189 nm^3 ; this corresponds to an approximate mass of 171 kDa.

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

7.3 Rotationally averaged power spectrum ⓘ

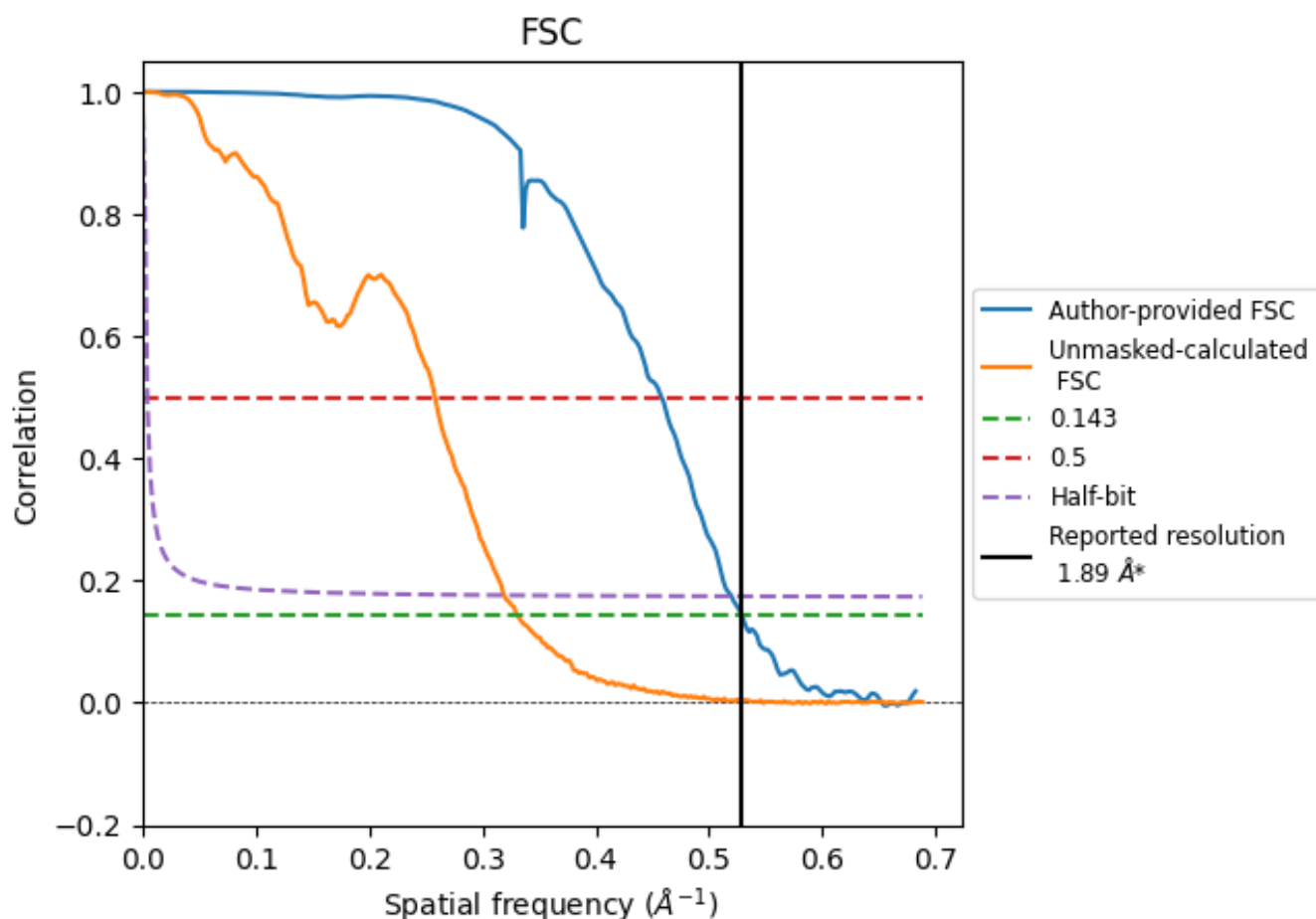


*Reported resolution corresponds to spatial frequency of 0.529 Å⁻¹

8 Fourier-Shell correlation [i](#)

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

8.1 FSC [i](#)



*Reported resolution corresponds to spatial frequency of 0.529 \AA^{-1}

8.2 Resolution estimates [i](#)

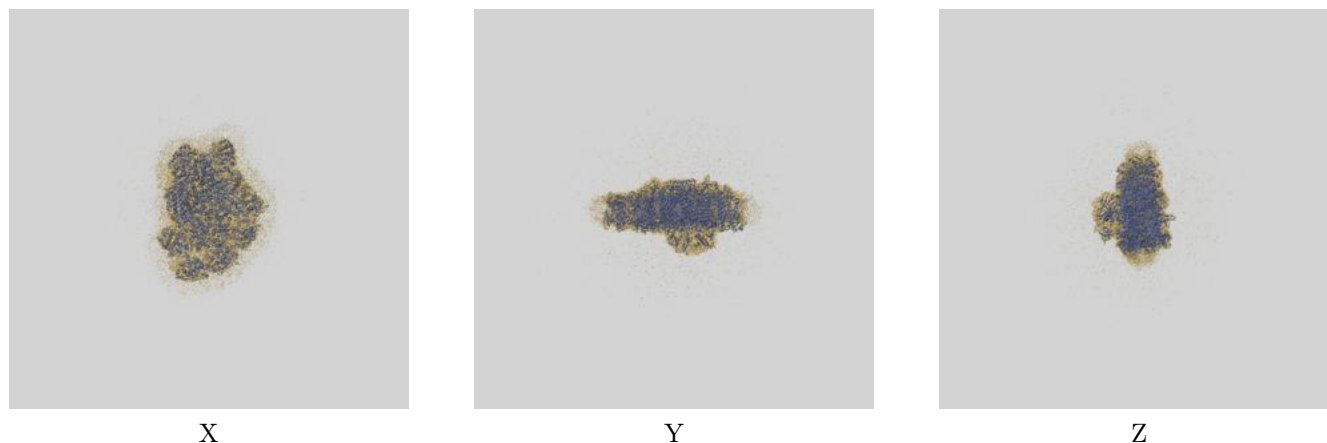
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	1.89	-	-
Author-provided FSC curve	1.89	2.18	1.92
Unmasked-calculated*	3.02	3.87	3.13

*Resolution estimate based on FSC curve calculated by comparison of deposited half-maps. The value from deposited half-maps intersecting FSC 0.143 CUT-OFF 3.02 differs from the reported value 1.89 by more than 10 %

9 Map-model fit [i](#)

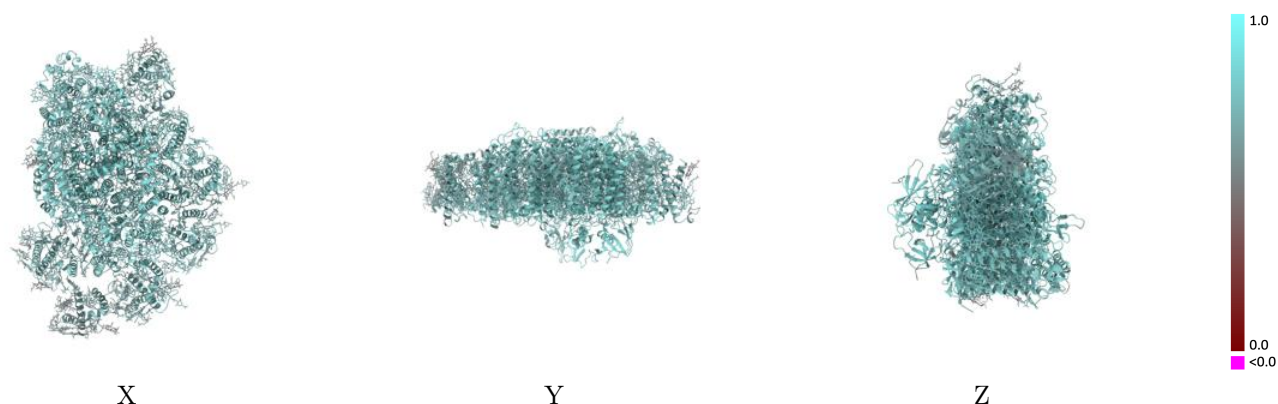
This section contains information regarding the fit between EMDB map EMD-64378 and PDB model 9UOF. Per-residue inclusion information can be found in section [3](#) on page [26](#).

9.1 Map-model overlay [i](#)



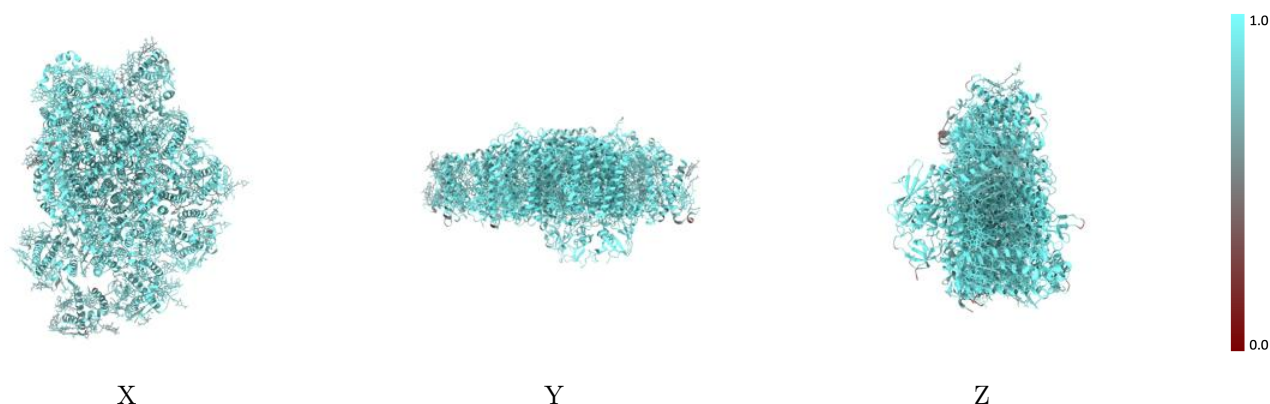
The images above show the 3D surface view of the map at the recommended contour level 0.08 at 50% transparency in yellow overlaid with a ribbon representation of the model coloured in blue. These images allow for the visual assessment of the quality of fit between the atomic model and the map.

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



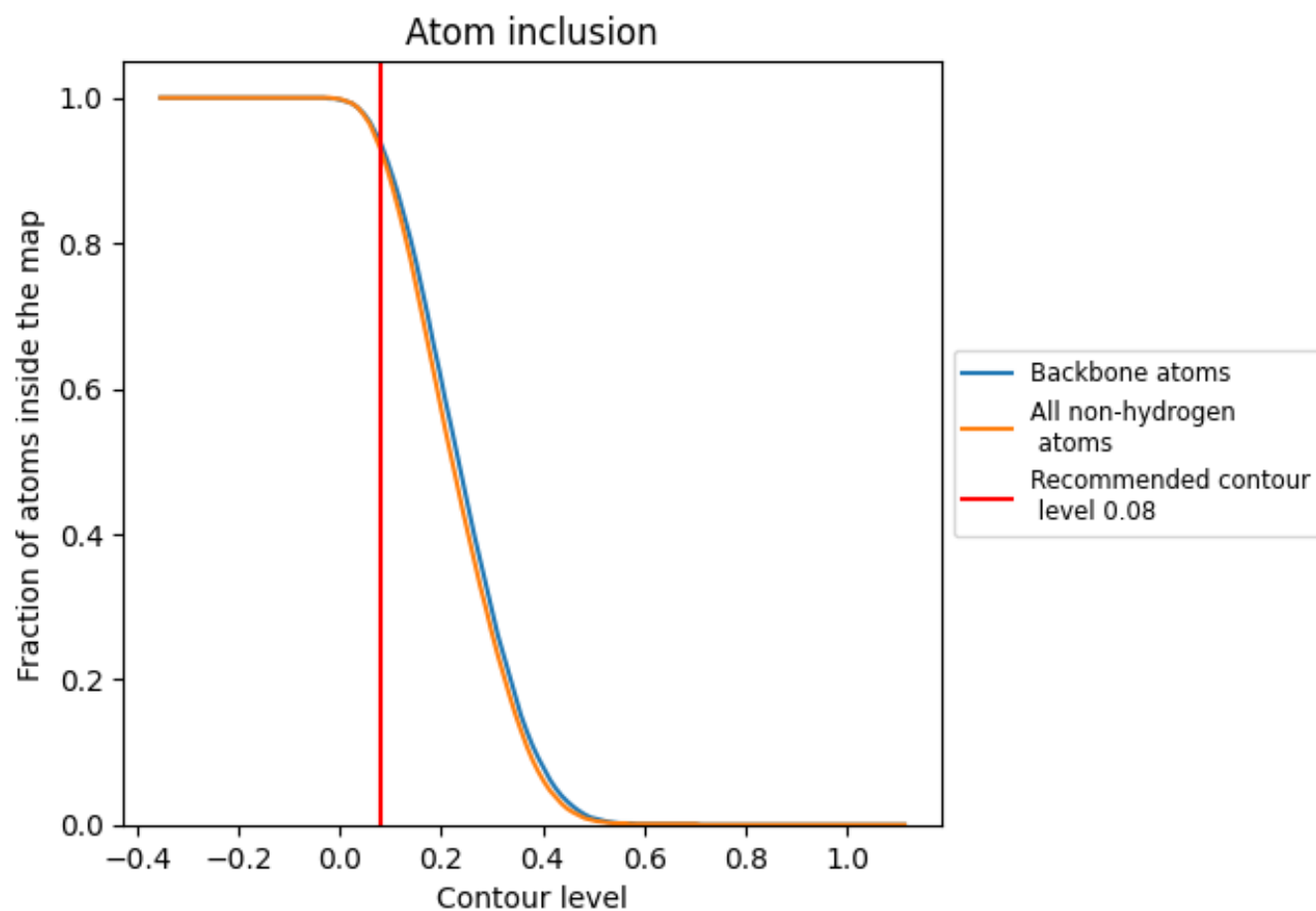
The images above show the model with each residue coloured according to its Q-score. This shows their resolvability in the map with higher Q-score values reflecting better resolvability. Please note: Q-score is calculating the resolvability of atoms, and thus high values are only expected at resolutions 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 (0.08).

9.4 Atom inclusion ⓘ



At the recommended contour level, 94% of all backbone atoms, 93% 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 (0.08) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	<div><div></div></div> 0.9300	<div><div></div></div> 0.7370
A	<div><div></div></div> 0.9580	<div><div></div></div> 0.7640
B	<div><div></div></div> 0.9770	<div><div></div></div> 0.7850
C	<div><div></div></div> 0.9920	<div><div></div></div> 0.8000
D	<div><div></div></div> 0.9440	<div><div></div></div> 0.7530
E	<div><div></div></div> 0.9060	<div><div></div></div> 0.7410
F	<div><div></div></div> 0.9400	<div><div></div></div> 0.7380
I	<div><div></div></div> 0.9740	<div><div></div></div> 0.7650
J	<div><div></div></div> 0.8270	<div><div></div></div> 0.6700
L	<div><div></div></div> 0.9540	<div><div></div></div> 0.7540
M	<div><div></div></div> 0.9530	<div><div></div></div> 0.7490
O	<div><div></div></div> 0.9110	<div><div></div></div> 0.7000
P	<div><div></div></div> 0.8790	<div><div></div></div> 0.6780
Q	<div><div></div></div> 0.8560	<div><div></div></div> 0.6690
R	<div><div></div></div> 0.9380	<div><div></div></div> 0.7320
S	<div><div></div></div> 0.9110	<div><div></div></div> 0.7100
T	<div><div></div></div> 0.7480	<div><div></div></div> 0.6010
U	<div><div></div></div> 0.8360	<div><div></div></div> 0.6690

1.0

0.0

<0.0