



wwPDB X-ray Structure Validation Summary Report ⓘ

Apr 21, 2024 – 08:09 am BST

PDB ID : 2BHW
Title : PEA LIGHT-HARVESTING COMPLEX II AT 2.5 ANGSTROM RESOLUTION
Authors : Standfuss, J.; Terwisscha van Scheltinga, A.C.; Lamborghini, M.; Kuehlbrandt, W.
Deposited on : 2005-01-19
Resolution : 2.50 Å(reported)

This is a wwPDB X-ray Structure 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/XrayValidationReportHelp>

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:

MolProbity : 4.02b-467
Mogul : 1.8.4, CSD as541be (2020)
Xtriage (Phenix) : 1.13
EDS : 2.36.2
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac : 5.8.0158
CCP4 : 7.0.044 (Gargrove)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.36.2

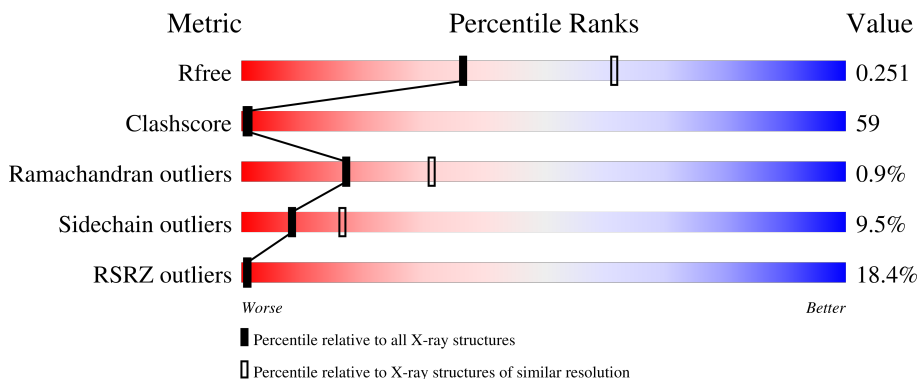
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 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)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	4661 (2.50-2.50)
Clashscore	141614	5346 (2.50-2.50)
Ramachandran outliers	138981	5231 (2.50-2.50)
Sidechain outliers	138945	5233 (2.50-2.50)
RSRZ outliers	127900	4559 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower 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 electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	232	
1	B	232	
1	C	232	

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
2	LUX	A	501	X	-	X	-
2	LUX	A	502	X	-	X	-
2	LUX	B	501	X	-	X	-
2	LUX	B	502	X	-	-	-
2	LUX	C	501	X	-	X	-
2	LUX	C	502	X	-	-	-
4	XAT	A	504	X	-	-	-
4	XAT	B	504	X	-	-	-
4	XAT	C	504	X	-	-	-
5	CLA	A	601	X	-	X	-
5	CLA	A	602	X	-	-	-
5	CLA	A	603	X	-	-	-
5	CLA	A	604	X	-	X	-
5	CLA	A	605	X	-	X	-
5	CLA	A	606	X	-	-	-
5	CLA	A	607	X	-	-	-
5	CLA	A	608	X	-	-	-
5	CLA	B	601	X	-	X	-
5	CLA	B	602	X	-	-	-
5	CLA	B	603	X	-	X	-
5	CLA	B	604	X	-	X	-
5	CLA	B	605	X	-	X	-
5	CLA	B	606	X	-	-	-
5	CLA	B	607	X	-	-	-
5	CLA	B	608	X	-	X	-
5	CLA	C	601	X	-	X	-
5	CLA	C	602	X	-	-	-
5	CLA	C	603	X	-	-	-
5	CLA	C	604	X	-	-	-
5	CLA	C	605	X	-	X	-
5	CLA	C	606	X	-	-	-
5	CLA	C	607	X	-	-	-
5	CLA	C	608	X	-	-	-
6	CHL	A	609	X	-	X	-
6	CHL	A	610	X	-	X	-
6	CHL	A	611	X	-	X	-
6	CHL	A	612	X	-	-	-
6	CHL	A	613	X	-	-	-
6	CHL	A	614	X	-	-	-
6	CHL	B	609	X	-	X	-
6	CHL	B	610	X	-	X	-

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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
6	CHL	B	611	X	-	X	-
6	CHL	B	612	X	-	X	-
6	CHL	B	613	X	-	-	-
6	CHL	B	614	X	-	-	-
6	CHL	C	609	X	-	X	-
6	CHL	C	610	X	-	X	-
6	CHL	C	611	X	-	X	-
6	CHL	C	612	X	-	X	-
6	CHL	C	613	X	-	-	-
6	CHL	C	614	X	-	-	-
8	DGD	A	802	X	-	-	X
8	DGD	B	802	X	-	-	X
8	DGD	C	802	X	-	-	X

2 Entry composition i

There are 9 unique types of molecules in this entry. The entry contains 8373 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, 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.

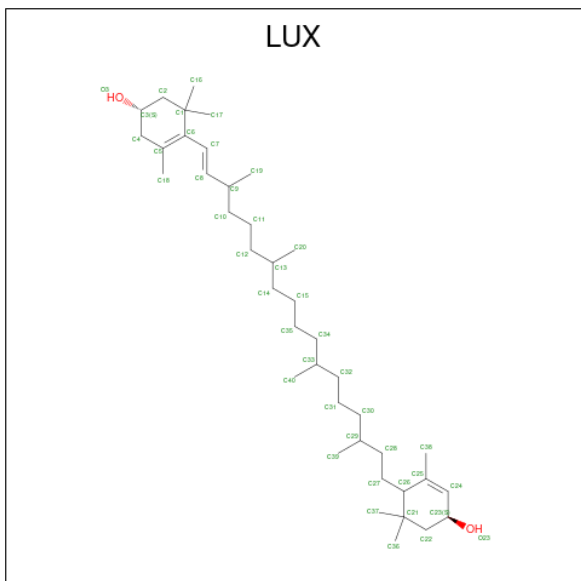
- Molecule 1 is a protein called CHLOROPHYLL A-B BINDING PROTEIN AB80.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	223	Total 1683	C 1090	N 274	O 315	S 4	0	0	0
1	B	223	Total 1683	C 1090	N 274	O 315	S 4	0	0	0
1	C	223	Total 1683	C 1090	N 274	O 315	S 4	0	0	0

There are 3 discrepancies between the modelled and reference sequences:

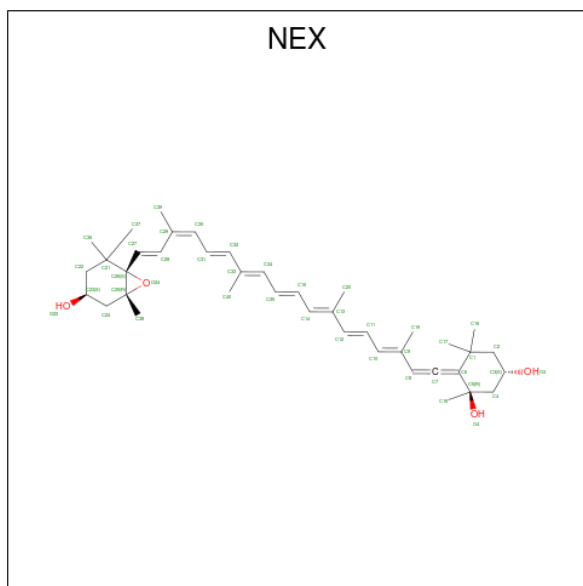
Chain	Residue	Modelled	Actual	Comment	Reference
A	79	SER	CYS	conflict	UNP P07371
B	79	SER	CYS	conflict	UNP P07371
C	79	SER	CYS	conflict	UNP P07371

- Molecule 2 is (3R,3'R,6'S,9R,9'R,13R,13'S)-4',5'-DIDEHYDRO-5',6',7',8',9,9',10,10',11,11',12,12',13,13',14,14',15,15'-OCTADECALHYDRO-BETA,BETA-CAROTENE-3,3'-DIOL (three-letter code: LUX) (formula: C₄₀H₇₂O₂).



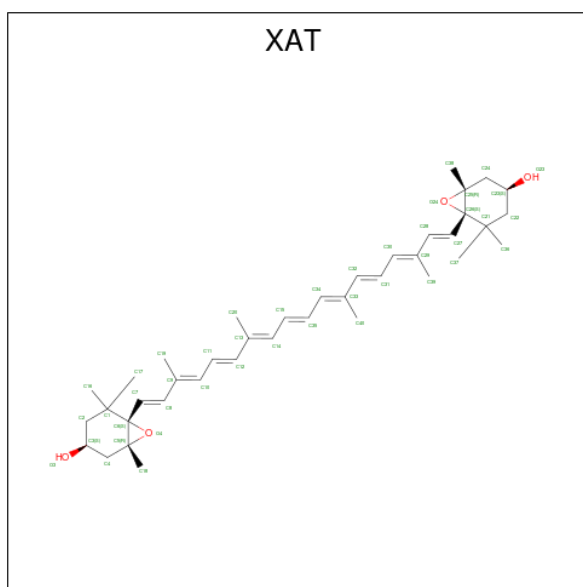
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			42	40	2		
2	A	1	Total	C	O	0	0
			42	40	2		
2	B	1	Total	C	O	0	0
			42	40	2		
2	B	1	Total	C	O	0	0
			42	40	2		
2	C	1	Total	C	O	0	0
			42	40	2		
2	C	1	Total	C	O	0	0
			42	40	2		

- Molecule 3 is (1R,3R)-6-[(3E,5E,7E,9E,11E,13E,15E,17E)-18-[(1S,4R,6R)-4-HYDROXY-2,2,6-TRIMETHYL-7-OXABICYCLO[4.1.0]HEPT-1-YL]-3,7,12,16-TETRAMETHYLOCTADEC-1,3,5,7,9,11,13,15,17-NONAENYLIDENE]-1,5,5-TRIMETHYLCYCLOHEXANE-1,3-DIOL (three-letter code: NEX) (formula: C₄₀H₅₆O₄).



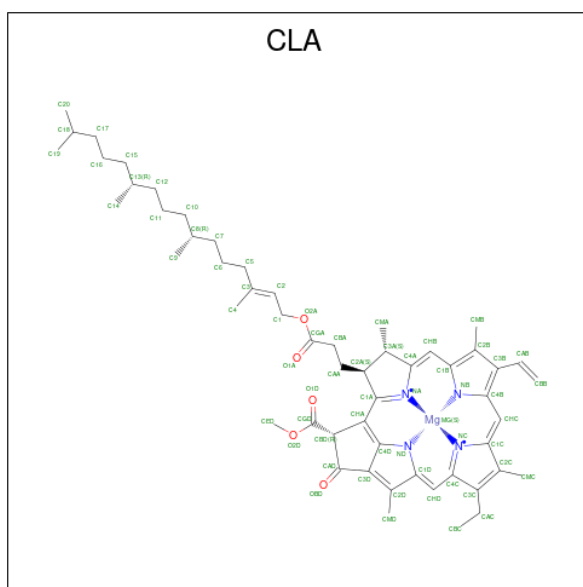
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total	C	O	0	0
			44	40	4		
3	B	1	Total	C	O	0	0
			44	40	4		
3	C	1	Total	C	O	0	0
			44	40	4		

- Molecule 4 is (3S,5R,6S,3'S,5'R,6'S)-5,6,5',6'-DIEPOXY-5,6,5',6'-TETRAHYDRO-BETA, BETA-CAROTENE-3,3'-DIOL (three-letter code: XAT) (formula: C₄₀H₅₆O₄).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
4	A	1	Total	C	O	0	0
			44	40	4		
4	B	1	Total	C	O	0	0
			44	40	4		
4	C	1	Total	C	O	0	0
			44	40	4		

- Molecule 5 is CHLOROPHYLL A (three-letter code: CLA) (formula: $C_{55}H_{72}MgN_4O_5$).



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

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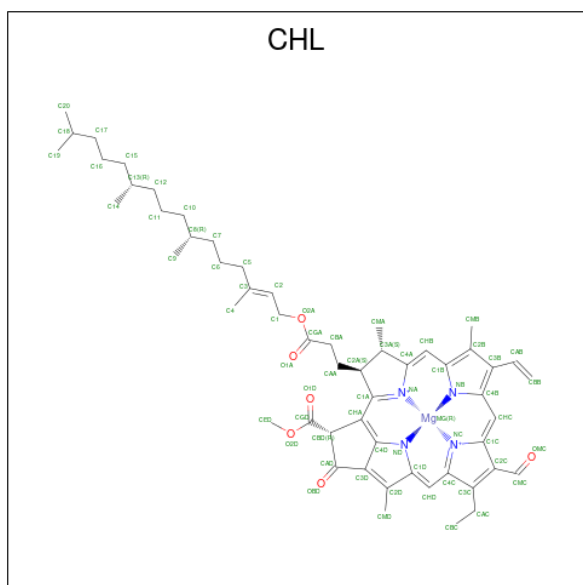
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
5	A	1	Total	C	Mg	N	O	0	0
			60	50	1	4	5		
5	A	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	A	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	A	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	A	1	Total	C	Mg	N	O	0	0
			57	47	1	4	5		
5	A	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	A	1	Total	C	Mg	N	O	0	0
			48	38	1	4	5		
5	B	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	B	1	Total	C	Mg	N	O	0	0
			60	50	1	4	5		
5	B	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	B	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	B	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	B	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	B	1	Total	C	Mg	N	O	0	0
			48	38	1	4	5		
5	C	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	C	1	Total	C	Mg	N	O	0	0
			60	50	1	4	5		
5	C	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	C	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	C	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	C	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	C	1	Total	C	Mg	N	O	0	0
			57	47	1	4	5		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
5	C	1	Total	C	Mg	N	O	0	0
			65	55	1	4	5		
5	C	1	Total	C	Mg	N	O	0	0
			48	38	1	4	5		

- Molecule 6 is CHLOROPHYLL B (three-letter code: CHL) (formula: $C_{55}H_{70}MgN_4O_6$).



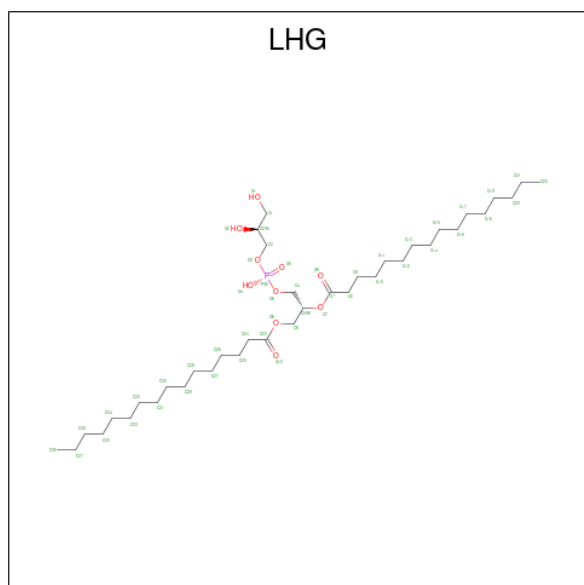
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
6	A	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	A	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	A	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	A	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	A	1	Total	C	Mg	N	O	0	0
			46	35	1	4	6		
6	A	1	Total	C	Mg	N	O	0	0
			42	33	1	4	4		
6	B	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	B	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	B	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
6	B	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	B	1	Total	C	Mg	N	O	0	0
			46	35	1	4	6		
6	B	1	Total	C	Mg	N	O	0	0
			42	33	1	4	4		
6	C	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	C	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	C	1	Total	C	Mg	N	O	0	0
			66	55	1	4	6		
6	C	1	Total	C	Mg	N	O	0	0
			46	35	1	4	6		
6	C	1	Total	C	Mg	N	O	0	0
			42	33	1	4	4		

- Molecule 7 is 1,2-DIPALMITOYL-PHOSPHATIDYL-GLYCEROLE (three-letter code: LHG) (formula: $C_{38}H_{76}O_{10}P$).



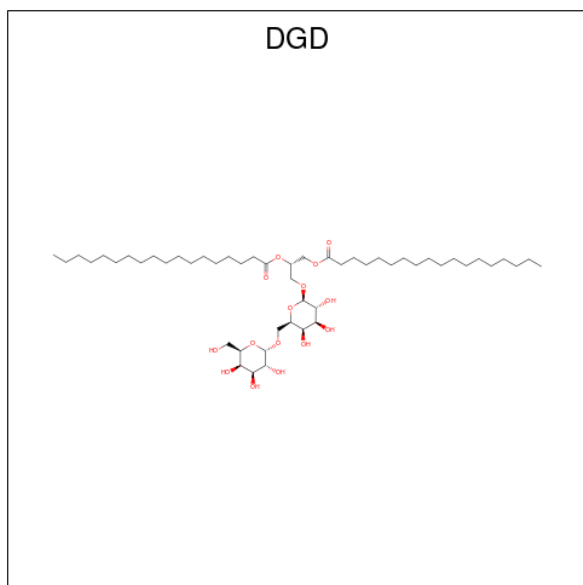
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
7	A	1	Total	C	O	P	0	0
			49	38	10	1		
7	B	1	Total	C	O	P	0	0
			49	38	10	1		

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Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	P		
7	C	1	49	38	10	1	0	0

- Molecule 8 is DIGALACTOSYL DIACYL GLYCEROL (DGDG) (three-letter code: DGD) (formula: $C_{51}H_{96}O_{15}$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	O		
8	A	1	38	25	13	0	0
8	B	1	38	25	13	0	0
8	C	1	38	25	13	0	0

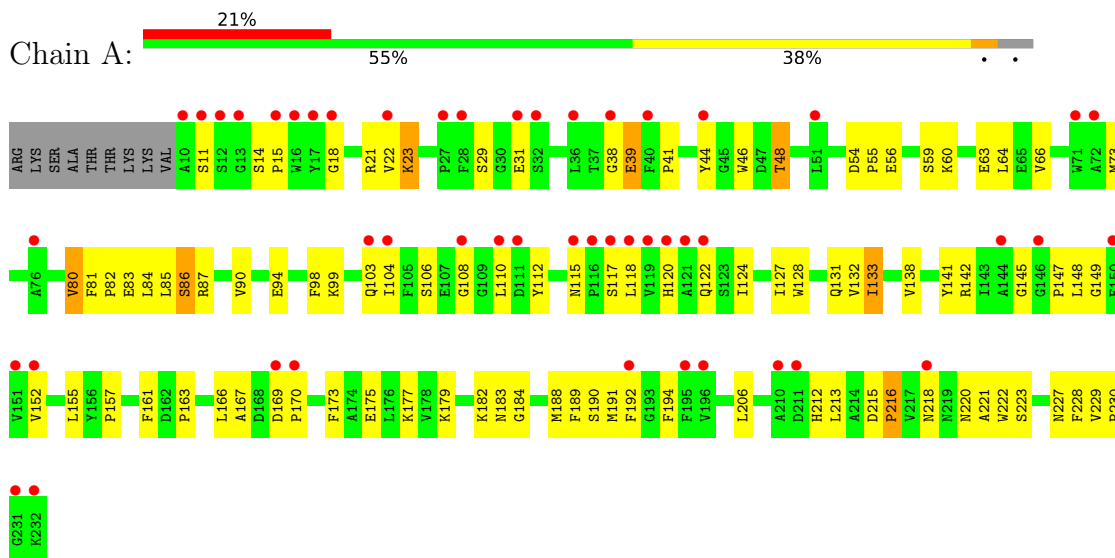
- Molecule 9 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
9	A	9	9	9	0	0
9	B	7	7	7	0	0
9	C	5	5	5	0	0

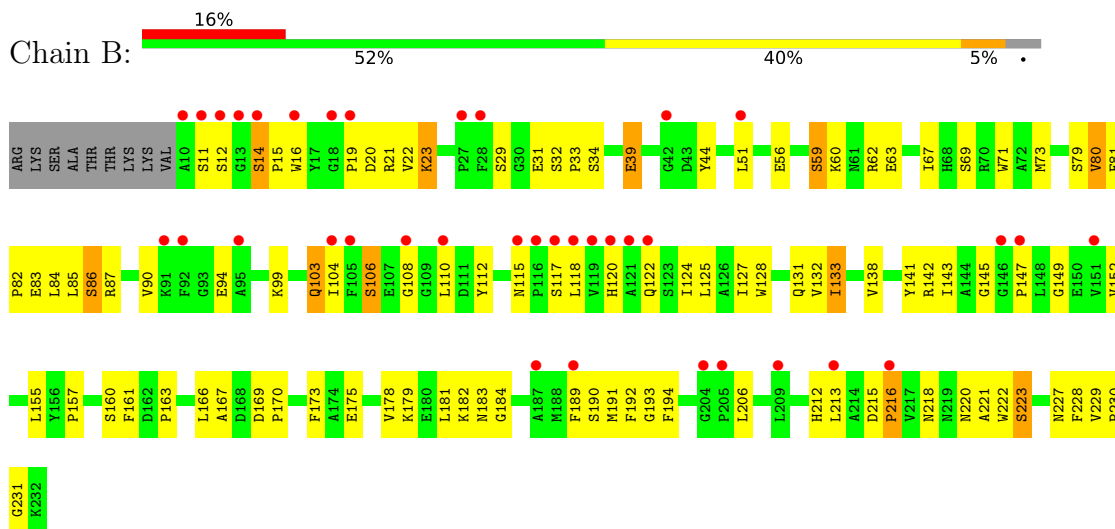
3 Residue-property plots [i](#)

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 electron 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 dot above a residue indicates a poor fit to the electron density ($RSRZ > 2$). 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: CHLOROPHYLL A-B BINDING PROTEIN AB80

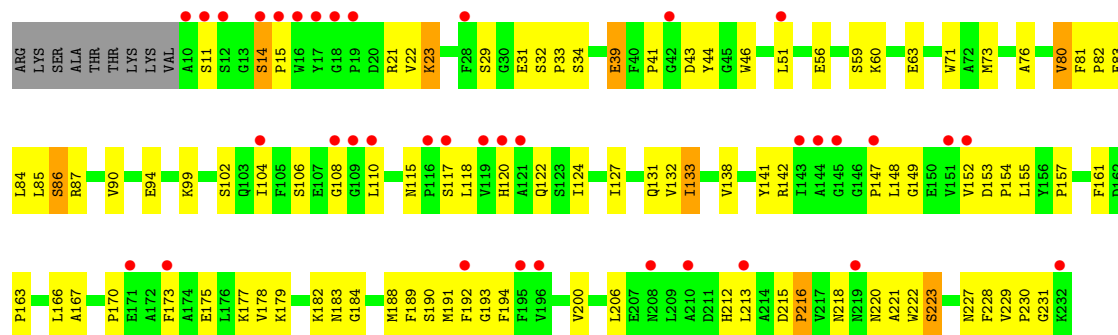


- Molecule 1: CHLOROPHYLL A-B BINDING PROTEIN AB80



- Molecule 1: CHLOROPHYLL A-B BINDING PROTEIN AB80





4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	211.40Å 128.00Å 62.00Å 90.00° 101.80° 90.00°	Depositor
Resolution (Å)	50.00 – 2.50 48.22 – 2.50	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-2.50) 85.7 (48.22-2.50)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.24 (at 2.51Å)	Xtrriage
Refinement program	TNT 5F	Depositor
R, R_{free}	0.220 , 0.241 0.232 , 0.251	Depositor DCC
R_{free} test set	990 reflections (2.07%)	wwPDB-VP
Wilson B-factor (Å ²)	92.5	Xtrriage
Anisotropy	0.183	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 103.8	EDS
L-test for twinning ²	$\langle L \rangle = 0.47$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	8373	wwPDB-VP
Average B, all atoms (Å ²)	91.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.96% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality [i](#)

5.1 Standard geometry [i](#)

Bond lengths and bond angles in the following residue types are not validated in this section: DGD, XAT, CLA, LUX, NEX, LHG, CHL

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	0.50	0/1735	0.66	0/2363
1	B	0.53	0/1735	0.69	0/2363
1	C	0.52	0/1735	0.66	0/2363
All	All	0.51	0/5205	0.67	0/7089

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts [i](#)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1683	0	1601	127	0
1	B	1683	0	1601	138	0
1	C	1683	0	1601	135	0
2	A	84	0	134	50	0
2	B	84	0	134	49	0
2	C	84	0	134	50	0
3	A	44	0	56	4	0
3	B	44	0	56	4	0
3	C	44	0	56	4	0
4	A	44	0	56	18	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	B	44	0	56	16	0
4	C	44	0	56	16	0
5	A	490	0	508	133	0
5	B	490	0	508	141	0
5	C	490	0	508	131	0
6	A	352	0	338	114	0
6	B	352	0	338	112	0
6	C	352	0	338	114	0
7	A	49	0	74	12	0
7	B	49	0	74	13	0
7	C	49	0	74	13	0
8	A	38	0	40	4	0
8	B	38	0	40	2	0
8	C	38	0	40	4	0
9	A	9	0	0	1	0
9	B	7	0	0	1	0
9	C	5	0	0	1	0
All	All	8373	0	8421	982	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 59.

The worst 5 of 982 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)
5:A:604:CLA:HMB1	5:A:604:CLA:HBB1	1.22	1.18
5:B:601:CLA:HBB1	5:B:601:CLA:HMB1	1.21	1.18
5:B:604:CLA:HBB1	5:B:604:CLA:HMB1	1.22	1.16
6:B:613:CHL:HBB1	6:B:613:CHL:HMB1	1.27	1.16
6:C:613:CHL:HBB1	6:C:613:CHL:HMB1	1.28	1.16

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

In the following table, the Percentiles column shows the percent Ramachandran outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	221/232 (95%)	210 (95%)	10 (4%)	1 (0%)	29	48
1	B	221/232 (95%)	210 (95%)	9 (4%)	2 (1%)	17	31
1	C	221/232 (95%)	208 (94%)	10 (4%)	3 (1%)	11	20
All	All	663/696 (95%)	628 (95%)	29 (4%)	6 (1%)	17	31

5 of 6 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	14	SER
1	C	14	SER
1	A	216	PRO
1	B	216	PRO
1	C	216	PRO

5.3.2 Protein sidechains [i](#)

In the following table, the Percentiles column shows the percent sidechain outliers of the chain as a percentile score with respect to all X-ray entries followed by that with respect to entries of similar resolution.

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	169/181 (93%)	154 (91%)	15 (9%)	9	19
1	B	169/181 (93%)	150 (89%)	19 (11%)	6	11
1	C	169/181 (93%)	155 (92%)	14 (8%)	11	22
All	All	507/543 (93%)	459 (90%)	48 (10%)	8	17

5 of 48 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	110	LEU
1	C	23	LYS
1	B	133	ILE
1	B	175	GLU
1	C	34	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	208	ASN
1	B	218	ASN
1	C	218	ASN
1	C	122	GLN
1	A	218	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 monosaccharides in this entry.

5.6 Ligand geometry [i](#)

60 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\# Z > 2$	Counts	RMSZ	$\# Z > 2$
4	XAT	C	504	-	39,47,47	2.20	13 (33%)	54,74,74	2.64	21 (38%)
5	CLA	C	604	1	65,73,73	1.23	8 (12%)	76,113,113	1.69	11 (14%)
3	NEX	C	503	-	38,46,46	2.25	8 (21%)	50,70,70	2.54	14 (28%)
6	CHL	C	609	1	66,74,74	1.49	10 (15%)	73,114,114	1.63	10 (13%)
2	LUX	C	501	-	43,43,43	4.65	25 (58%)	53,60,60	3.60	31 (58%)
8	DGD	C	802	-	39,39,67	1.41	5 (12%)	51,51,81	1.97	9 (17%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	CLA	B	607	7	65,73,73	1.56	10 (15%)	76,113,113	1.81	11 (14%)
2	LUX	B	501	-	43,43,43	4.71	25 (58%)	53,60,60	3.60	31 (58%)
6	CHL	B	610	9	66,74,74	1.59	10 (15%)	73,114,114	1.58	12 (16%)
2	LUX	B	502	-	43,43,43	4.65	25 (58%)	53,60,60	3.78	31 (58%)
5	CLA	A	604	1	65,73,73	1.36	8 (12%)	76,113,113	1.61	11 (14%)
4	XAT	A	504	-	39,47,47	2.21	15 (38%)	54,74,74	2.85	20 (37%)
5	CLA	B	608	1	48,56,73	1.56	8 (16%)	55,92,113	1.82	13 (23%)
5	CLA	B	603	1	65,73,73	1.28	9 (13%)	76,113,113	1.75	11 (14%)
3	NEX	B	503	-	38,46,46	2.40	10 (26%)	50,70,70	2.53	12 (24%)
5	CLA	C	602	1	60,68,73	1.50	11 (18%)	70,107,113	1.65	9 (12%)
6	CHL	C	612	1	66,74,74	1.58	12 (18%)	73,114,114	1.82	12 (16%)
6	CHL	A	613	-	46,54,74	1.72	8 (17%)	49,90,114	1.78	9 (18%)
7	LHG	C	801	5	48,48,48	1.68	6 (12%)	51,54,54	1.33	4 (7%)
5	CLA	C	605	1	65,73,73	1.33	10 (15%)	76,113,113	1.71	9 (11%)
5	CLA	C	607	7	65,73,73	1.43	11 (16%)	76,113,113	1.76	11 (14%)
4	XAT	B	504	-	39,47,47	2.20	14 (35%)	54,74,74	2.85	21 (38%)
6	CHL	A	611	9	66,74,74	1.46	12 (18%)	73,114,114	1.84	14 (19%)
5	CLA	A	608	1	48,56,73	1.71	10 (20%)	55,92,113	2.05	13 (23%)
5	CLA	B	604	1	65,73,73	1.43	9 (13%)	76,113,113	1.65	14 (18%)
6	CHL	A	614	1	42,50,74	1.77	8 (19%)	44,85,114	2.15	11 (25%)
2	LUX	C	502	-	43,43,43	4.52	26 (60%)	53,60,60	3.75	32 (60%)
5	CLA	A	602	1	60,68,73	1.39	9 (15%)	70,107,113	1.63	9 (12%)
8	DGD	B	802	-	39,39,67	1.32	4 (10%)	51,51,81	1.93	10 (19%)
5	CLA	C	603	1	65,73,73	1.20	7 (10%)	76,113,113	1.73	11 (14%)
6	CHL	B	614	1	42,50,74	1.86	9 (21%)	44,85,114	2.29	10 (22%)
5	CLA	A	601	1	65,73,73	1.34	10 (15%)	76,113,113	1.83	14 (18%)
5	CLA	A	603	1	65,73,73	1.23	8 (12%)	76,113,113	1.71	11 (14%)
6	CHL	C	614	1	42,50,74	1.87	9 (21%)	44,85,114	2.22	10 (22%)
5	CLA	A	606	-	57,65,73	1.27	6 (10%)	66,103,113	1.86	12 (18%)
3	NEX	A	503	-	38,46,46	2.32	11 (28%)	50,70,70	2.39	11 (22%)
6	CHL	A	609	1	66,74,74	1.44	9 (13%)	73,114,114	1.62	10 (13%)
2	LUX	A	502	-	43,43,43	4.59	26 (60%)	53,60,60	3.70	32 (60%)
5	CLA	B	601	1	65,73,73	1.33	13 (20%)	76,113,113	1.80	16 (21%)
6	CHL	A	610	9	66,74,74	1.56	12 (18%)	73,114,114	1.51	9 (12%)
6	CHL	A	612	1	66,74,74	1.50	11 (16%)	73,114,114	1.84	11 (15%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
5	CLA	B	606	-	57,65,73	1.30	9 (15%)	66,103,113	1.80	13 (19%)
6	CHL	C	611	9	66,74,74	1.47	12 (18%)	73,114,114	1.80	13 (17%)
5	CLA	B	605	1	65,73,73	1.39	11 (16%)	76,113,113	1.66	8 (10%)
6	CHL	B	609	1	66,74,74	1.52	10 (15%)	73,114,114	1.61	12 (16%)
6	CHL	C	613	-	46,54,74	1.64	9 (19%)	49,90,114	1.90	9 (18%)
8	DGD	A	802	-	39,39,67	1.32	4 (10%)	51,51,81	1.95	9 (17%)
2	LUX	A	501	-	43,43,43	4.65	26 (60%)	53,60,60	3.53	30 (56%)
6	CHL	B	612	1	66,74,74	1.47	12 (18%)	73,114,114	1.68	10 (13%)
5	CLA	C	601	1	65,73,73	1.37	9 (13%)	76,113,113	1.90	12 (15%)
5	CLA	C	606	-	57,65,73	1.33	7 (12%)	66,103,113	1.83	13 (19%)
5	CLA	C	608	1	48,56,73	1.45	7 (14%)	55,92,113	2.12	13 (23%)
6	CHL	C	610	9	66,74,74	1.56	12 (18%)	73,114,114	1.59	11 (15%)
6	CHL	B	613	-	46,54,74	1.61	10 (21%)	49,90,114	1.82	10 (20%)
5	CLA	B	602	1	60,68,73	1.53	12 (20%)	70,107,113	1.73	10 (14%)
7	LHG	A	801	5	48,48,48	1.71	6 (12%)	51,54,54	1.35	4 (7%)
6	CHL	B	611	9	66,74,74	1.49	11 (16%)	73,114,114	1.86	12 (16%)
5	CLA	A	605	1	65,73,73	1.32	9 (13%)	76,113,113	1.72	11 (14%)
7	LHG	B	801	5	48,48,48	1.66	6 (12%)	51,54,54	1.37	4 (7%)
5	CLA	A	607	7	65,73,73	1.50	11 (16%)	76,113,113	1.78	10 (13%)

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
4	XAT	C	504	-	2/2/12/26	2/31/93/93	0/4/4/4
5	CLA	C	604	1	2/2/15/20	15/37/115/115	-
8	DGD	C	802	-	1/1/11/13	16/24/64/95	0/2/2/2
6	CHL	C	609	1	4/4/20/26	15/39/137/137	-
2	LUX	C	501	-	5/5/12/15	11/29/67/67	0/2/2/2
3	NEX	C	503	-	-	2/27/83/83	0/3/3/3
5	CLA	B	607	7	1/1/15/20	16/37/115/115	-
2	LUX	B	501	-	5/5/12/15	11/29/67/67	0/2/2/2
6	CHL	B	610	9	4/4/20/26	14/39/137/137	-
2	LUX	B	502	-	5/5/12/15	11/29/67/67	0/2/2/2

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	CLA	A	604	1	2/2/15/20	12/37/115/115	-
4	XAT	A	504	-	2/2/12/26	0/31/93/93	0/4/4/4
5	CLA	B	608	1	1/1/11/20	7/17/95/115	-
5	CLA	B	603	1	2/2/15/20	12/37/115/115	-
5	CLA	C	602	1	1/1/14/20	13/31/109/115	-
6	CHL	C	612	1	3/3/20/26	16/39/137/137	-
3	NEX	B	503	-	-	2/27/83/83	0/3/3/3
6	CHL	A	613	-	3/3/16/26	0/15/113/137	-
7	LHG	C	801	5	-	22/53/53/53	-
5	CLA	C	605	1	2/2/15/20	11/37/115/115	-
5	CLA	C	607	7	1/1/15/20	15/37/115/115	-
4	XAT	B	504	-	2/2/12/26	2/31/93/93	0/4/4/4
6	CHL	A	611	9	4/4/20/26	16/39/137/137	-
5	CLA	A	608	1	1/1/11/20	6/17/95/115	-
5	CLA	B	604	1	2/2/15/20	12/37/115/115	-
6	CHL	A	614	1	3/3/15/26	0/10/108/137	-
2	LUX	C	502	-	5/5/12/15	10/29/67/67	0/2/2/2
5	CLA	A	602	1	1/1/14/20	13/31/109/115	-
8	DGD	B	802	-	1/1/11/13	17/24/64/95	0/2/2/2
5	CLA	C	603	1	2/2/15/20	12/37/115/115	-
6	CHL	B	614	1	3/3/15/26	0/10/108/137	-
5	CLA	A	601	1	2/2/15/20	12/37/115/115	-
5	CLA	A	603	1	2/2/15/20	13/37/115/115	-
6	CHL	C	614	1	3/3/15/26	1/10/108/137	-
5	CLA	A	606	-	2/2/13/20	9/28/106/115	-
6	CHL	A	609	1	4/4/20/26	18/39/137/137	-
3	NEX	A	503	-	-	2/27/83/83	0/3/3/3
2	LUX	A	502	-	5/5/12/15	10/29/67/67	0/2/2/2
5	CLA	B	601	1	2/2/15/20	12/37/115/115	-
6	CHL	A	610	9	4/4/20/26	14/39/137/137	-
6	CHL	A	612	1	3/3/20/26	16/39/137/137	-
5	CLA	B	606	-	2/2/13/20	10/28/106/115	-
6	CHL	C	611	9	4/4/20/26	16/39/137/137	-
5	CLA	B	605	1	2/2/15/20	11/37/115/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
6	CHL	B	609	1	4/4/20/26	16/39/137/137	-
6	CHL	C	613	-	3/3/16/26	0/15/113/137	-
8	DGD	A	802	-	1/1/11/13	16/24/64/95	0/2/2/2
2	LUX	A	501	-	5/5/12/15	11/29/67/67	0/2/2/2
6	CHL	B	612	1	3/3/20/26	14/39/137/137	-
5	CLA	C	601	1	2/2/15/20	13/37/115/115	-
5	CLA	C	606	-	2/2/13/20	10/28/106/115	-
5	CLA	C	608	1	1/1/11/20	6/17/95/115	-
6	CHL	C	610	9	4/4/20/26	14/39/137/137	-
6	CHL	B	613	-	3/3/16/26	1/15/113/137	-
5	CLA	B	602	1	1/1/14/20	13/31/109/115	-
7	LHG	A	801	5	-	22/53/53/53	-
6	CHL	B	611	9	4/4/20/26	17/39/137/137	-
5	CLA	A	605	1	2/2/15/20	11/37/115/115	-
7	LHG	B	801	5	-	22/53/53/53	-
5	CLA	A	607	7	1/1/15/20	16/37/115/115	-

The worst 5 of 663 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	501	LUX	C24-C25	19.56	1.57	1.33
2	B	502	LUX	C24-C25	18.51	1.56	1.33
2	A	501	LUX	C24-C25	18.34	1.56	1.33
2	B	501	LUX	C24-C25	18.20	1.55	1.33
2	A	502	LUX	C24-C25	18.19	1.55	1.33

The worst 5 of 797 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	501	LUX	C27-C26-C25	12.98	136.38	111.96
2	C	501	LUX	C27-C26-C25	12.57	135.60	111.96
2	B	502	LUX	C27-C26-C25	12.25	135.00	111.96
2	B	501	LUX	C27-C26-C25	12.24	134.99	111.96
2	A	502	LUX	C27-C26-C25	12.02	134.56	111.96

5 of 141 chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
2	A	501	LUX	C29
2	A	501	LUX	C26
2	A	501	LUX	C9
2	A	501	LUX	C13
2	A	501	LUX	C33

5 of 657 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	LUX	C11-C10-C9-C19
2	A	501	LUX	C11-C12-C13-C20
2	A	501	LUX	C20-C13-C14-C15
2	A	501	LUX	C25-C26-C27-C28
2	A	501	LUX	C27-C28-C29-C39

There are no ring outliers.

60 monomers are involved in 801 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	C	504	XAT	16	0
5	C	604	CLA	20	0
3	C	503	NEX	4	0
6	C	609	CHL	30	0
2	C	501	LUX	31	0
8	C	802	DGD	4	0
5	B	607	CLA	17	0
2	B	501	LUX	31	0
6	B	610	CHL	28	0
2	B	502	LUX	18	0
5	A	604	CLA	22	0
4	A	504	XAT	18	0
5	B	608	CLA	21	0
5	B	603	CLA	24	0
3	B	503	NEX	4	0
5	C	602	CLA	11	0
6	C	612	CHL	21	0
6	A	613	CHL	6	0
7	C	801	LHG	13	0
5	C	605	CLA	21	0
5	C	607	CLA	19	0
4	B	504	XAT	16	0
6	A	611	CHL	22	0
5	A	608	CLA	16	0

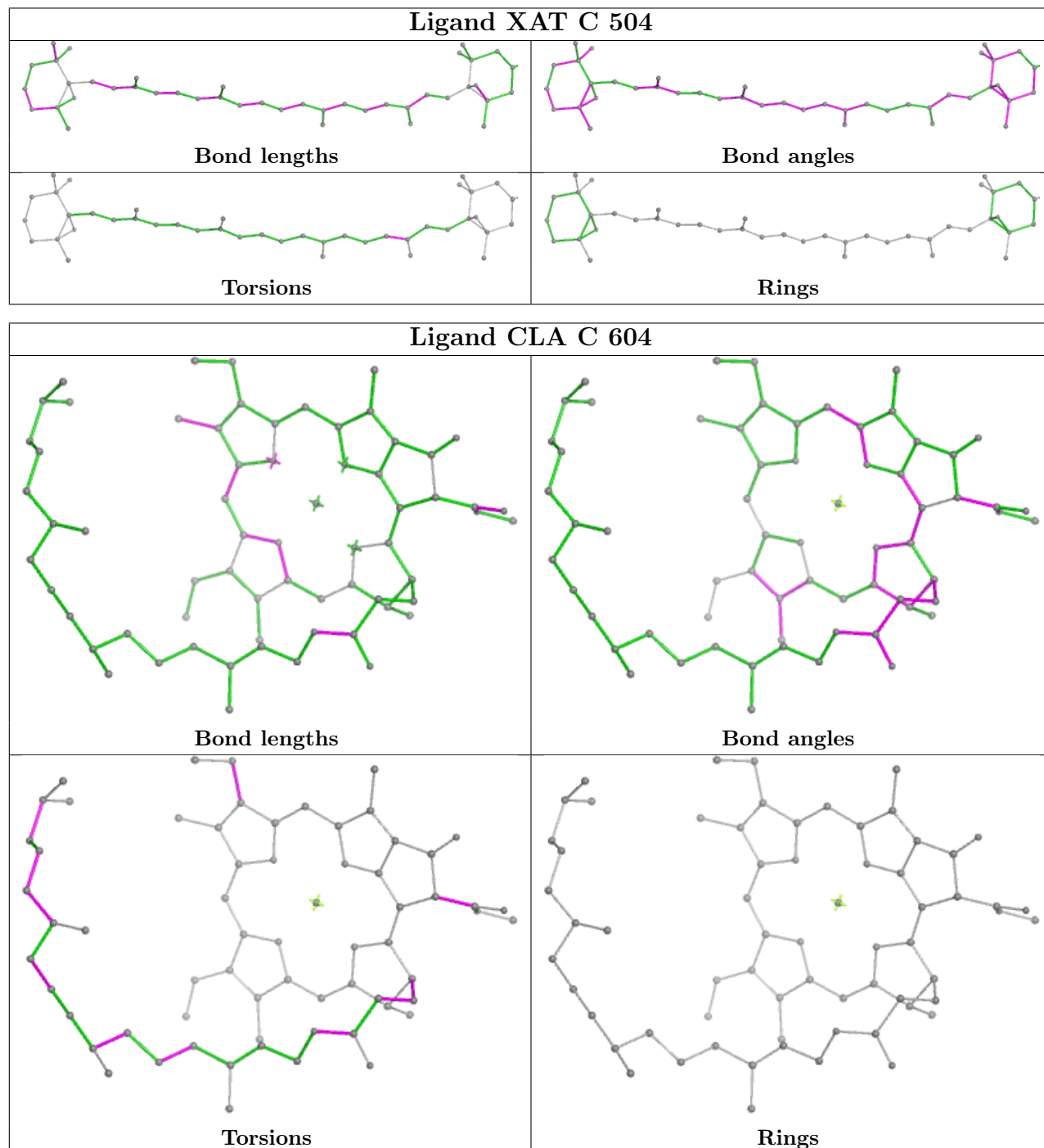
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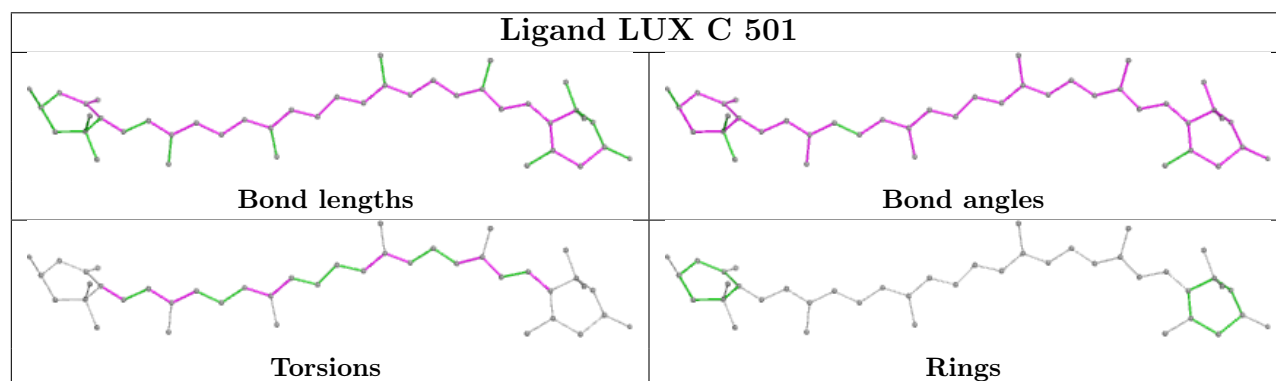
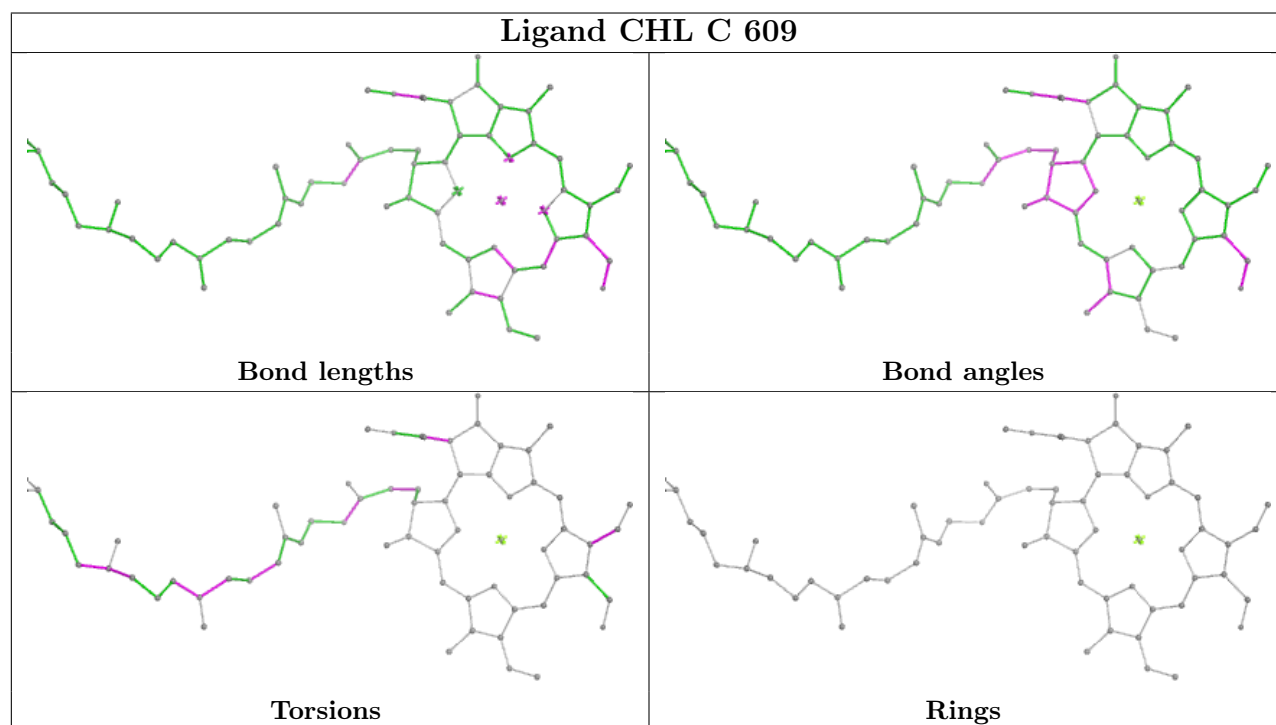
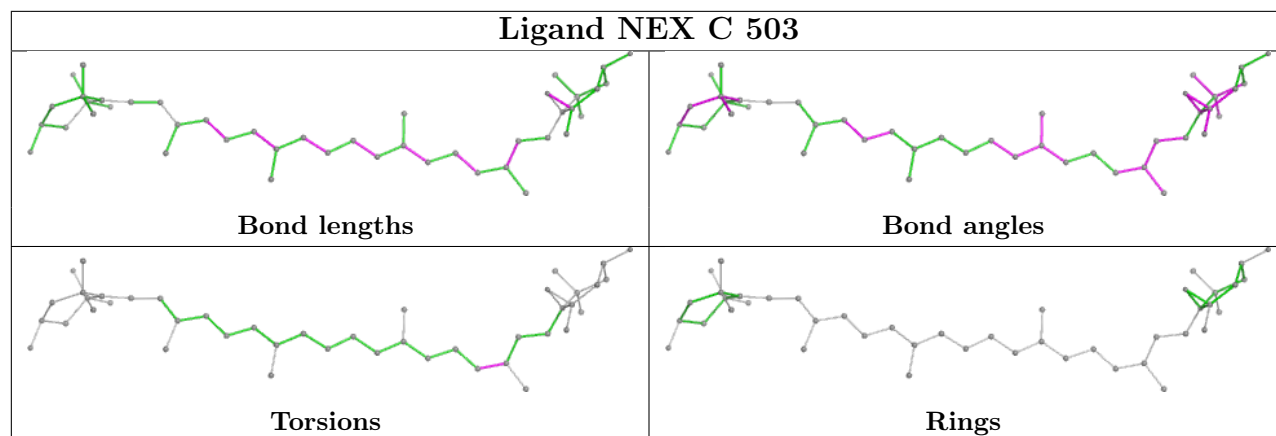
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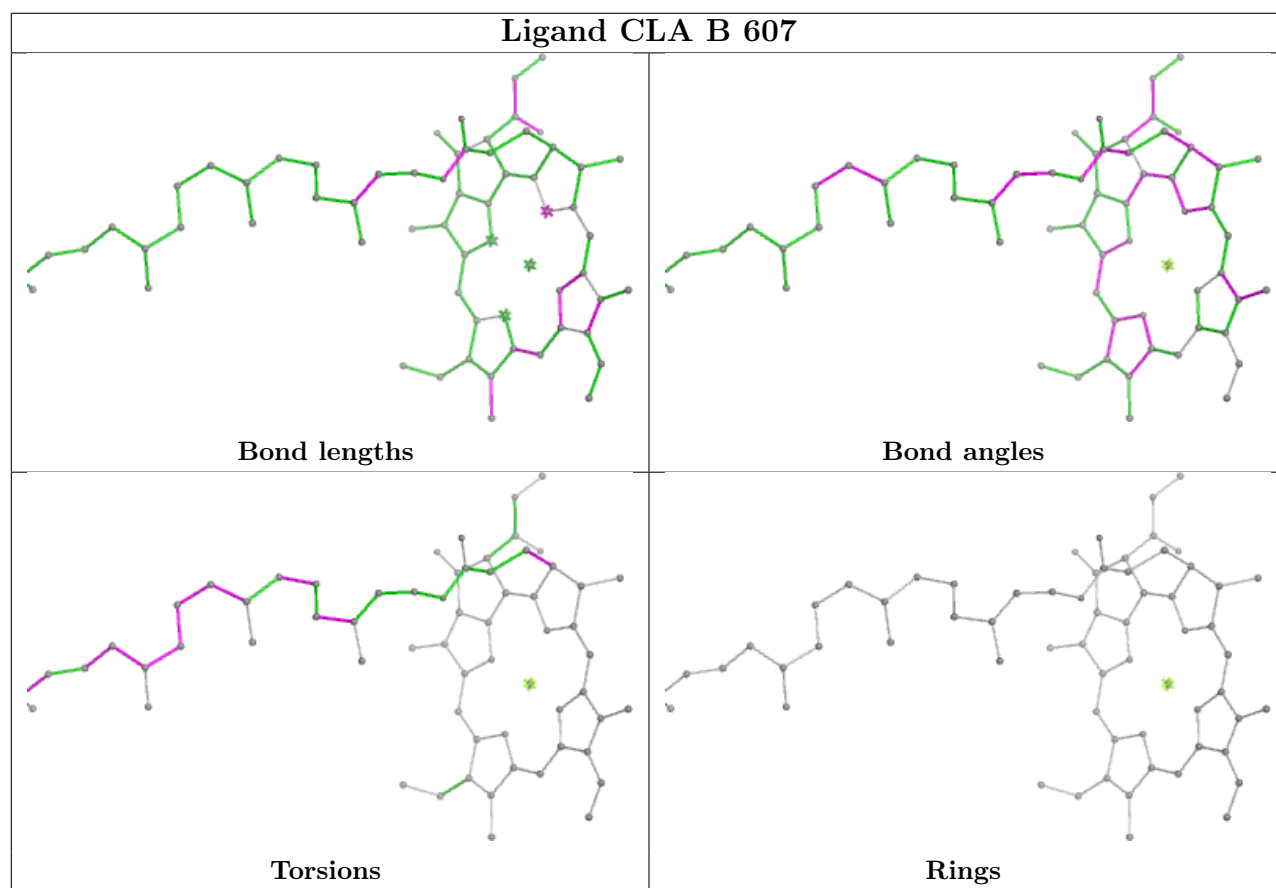
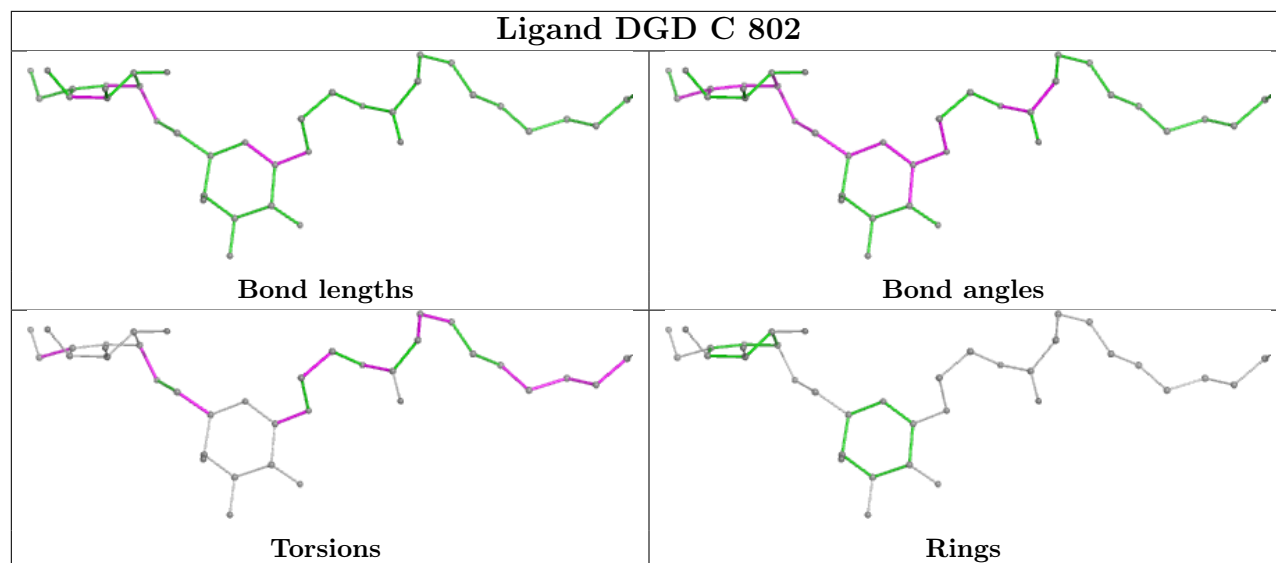
Mol	Chain	Res	Type	Clashes	Symm-Clashes
5	B	604	CLA	24	0
6	A	614	CHL	15	0
2	C	502	LUX	19	0
5	A	602	CLA	13	0
8	B	802	DGD	2	0
5	C	603	CLA	17	0
6	B	614	CHL	16	0
5	A	601	CLA	21	0
5	A	603	CLA	18	0
6	C	614	CHL	14	0
5	A	606	CLA	18	0
3	A	503	NEX	4	0
6	A	609	CHL	31	0
2	A	502	LUX	21	0
5	B	601	CLA	23	0
6	A	610	CHL	31	0
6	A	612	CHL	18	0
5	B	606	CLA	17	0
6	C	611	CHL	23	0
5	B	605	CLA	23	0
6	B	609	CHL	27	0
6	C	613	CHL	5	0
8	A	802	DGD	4	0
2	A	501	LUX	29	0
6	B	612	CHL	22	0
5	C	601	CLA	23	0
5	C	606	CLA	17	0
5	C	608	CLA	16	0
6	C	610	CHL	31	0
6	B	613	CHL	9	0
5	B	602	CLA	11	0
7	A	801	LHG	12	0
6	B	611	CHL	22	0
5	A	605	CLA	21	0
7	B	801	LHG	13	0
5	A	607	CLA	17	0

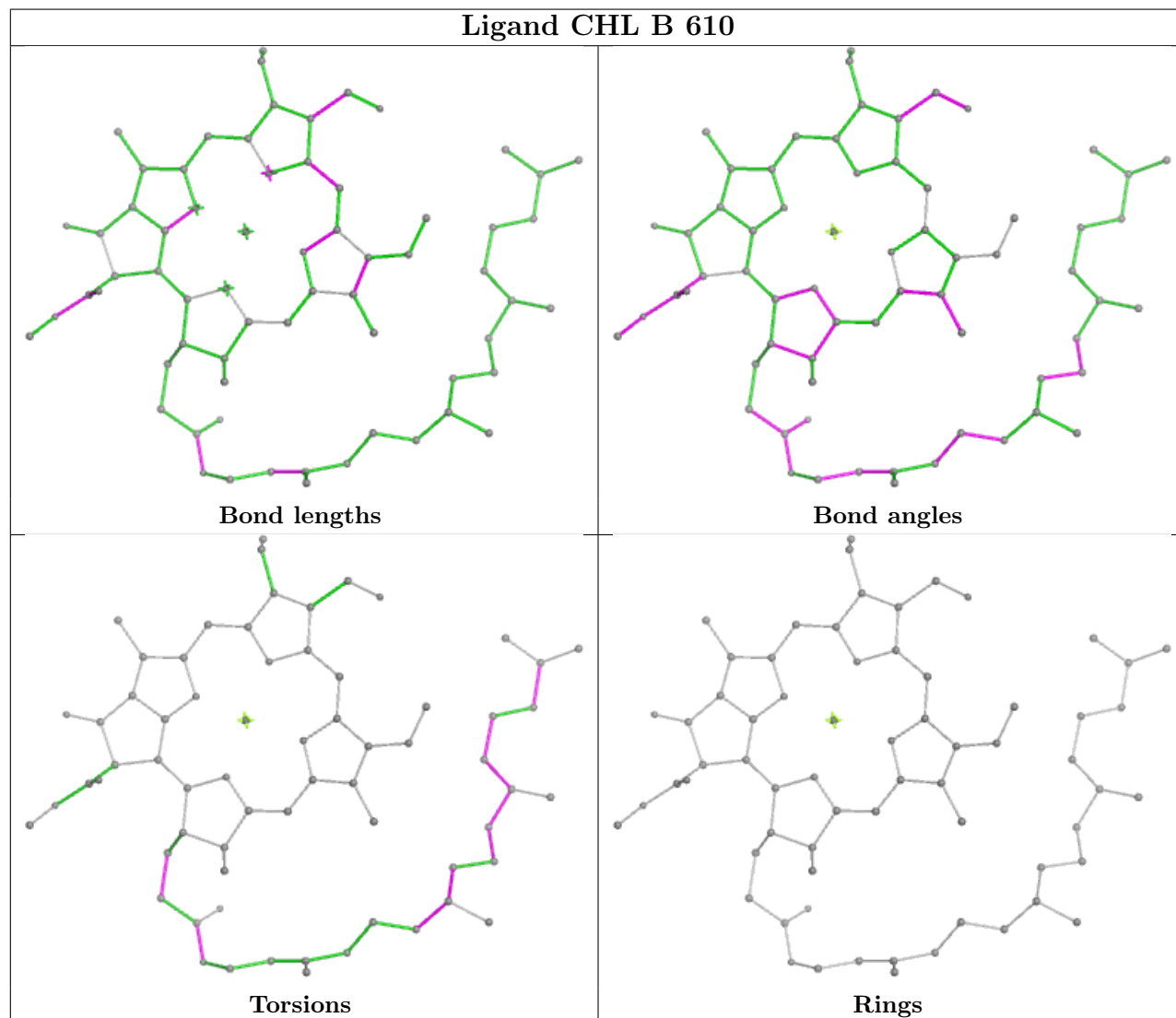
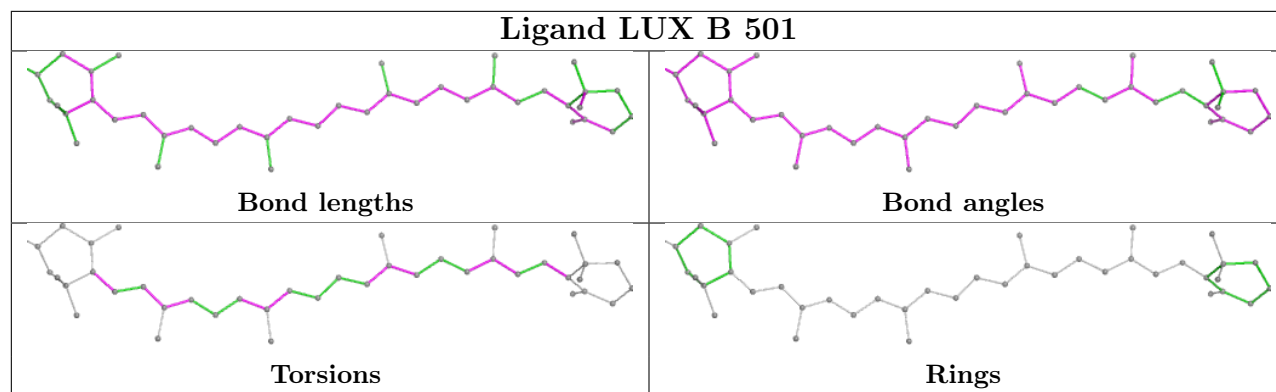
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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

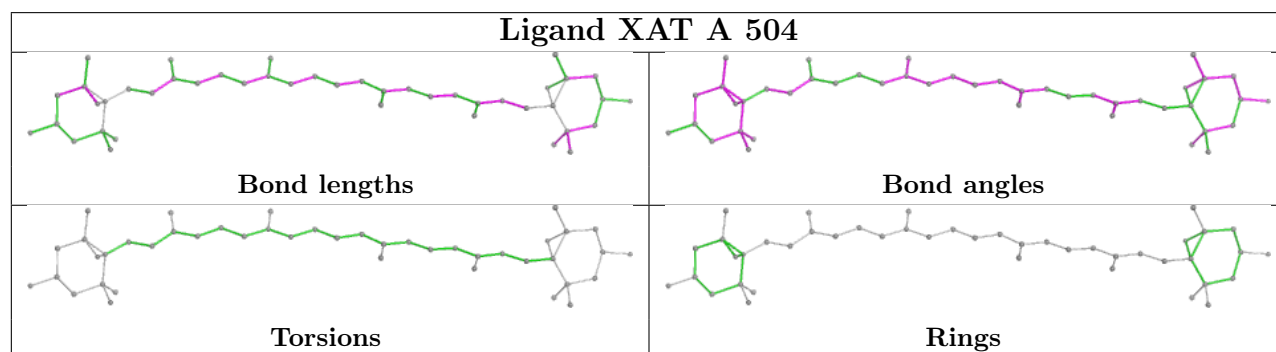
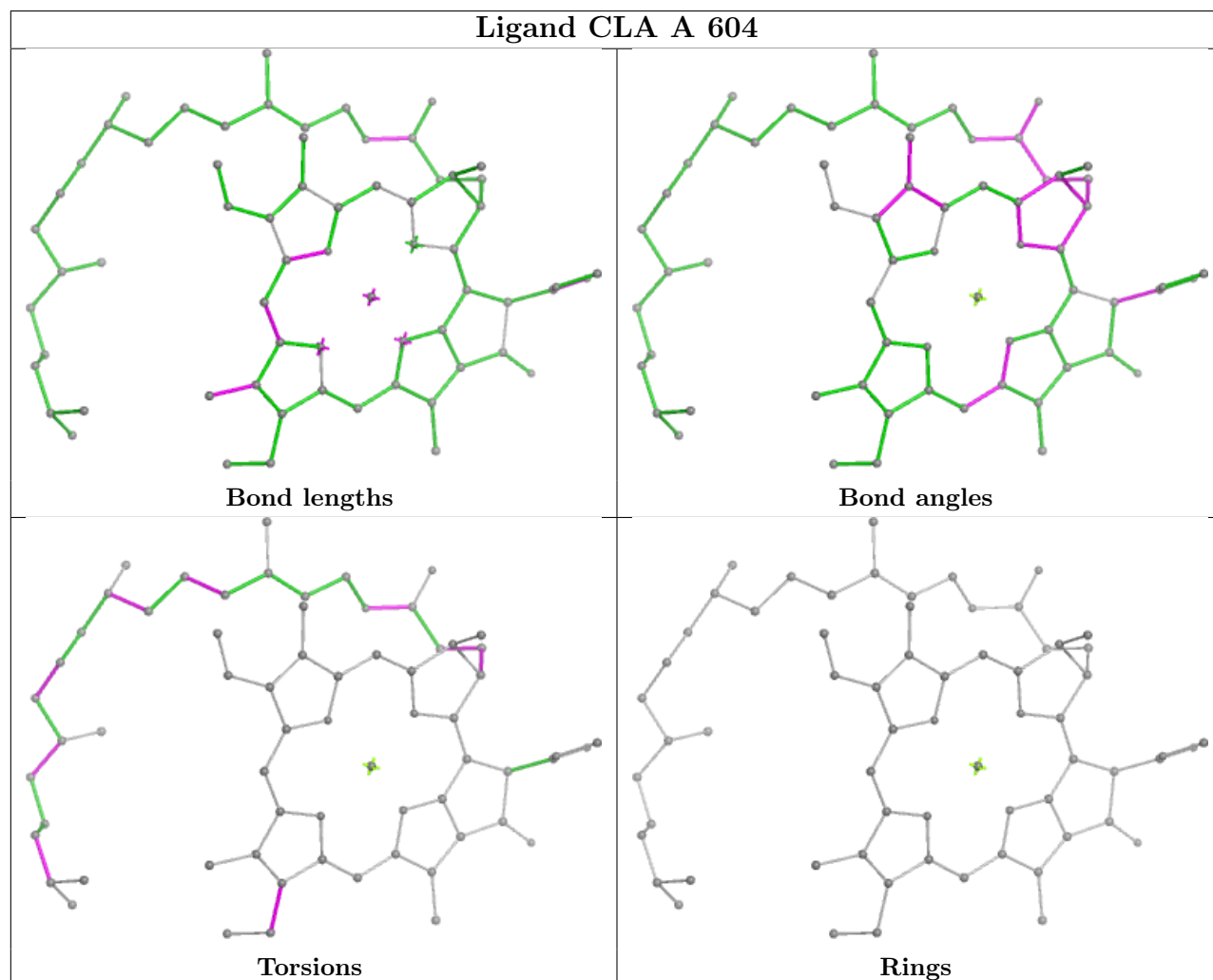
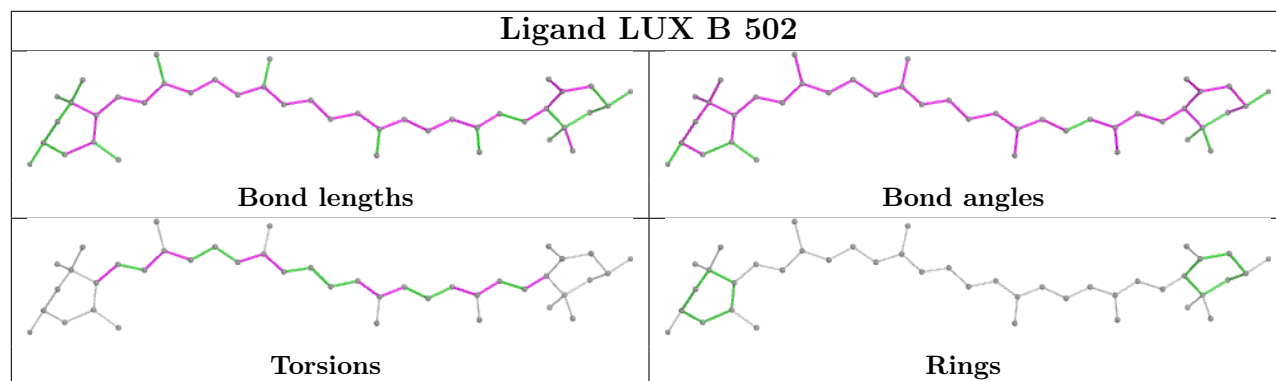
highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified 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.

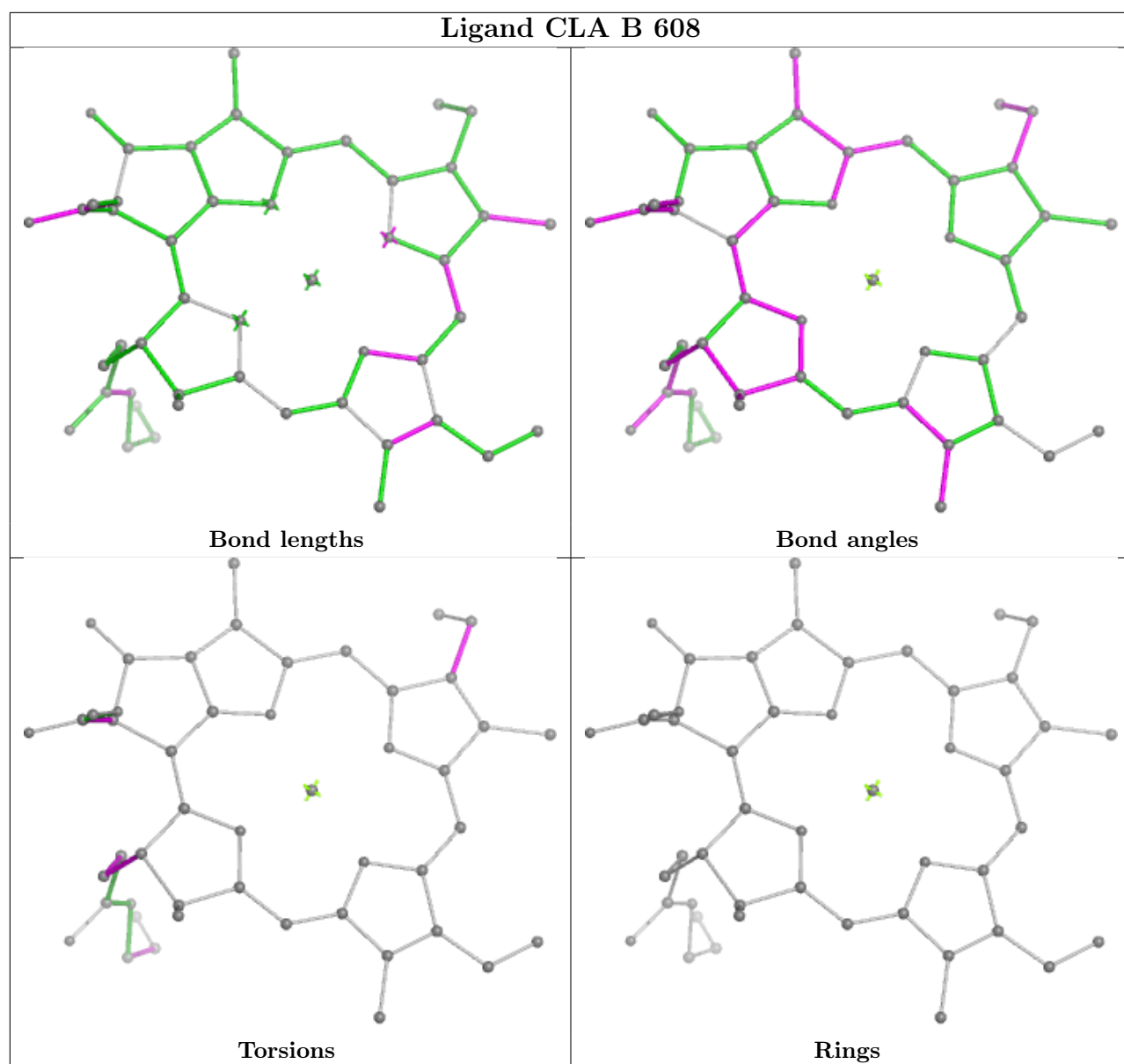


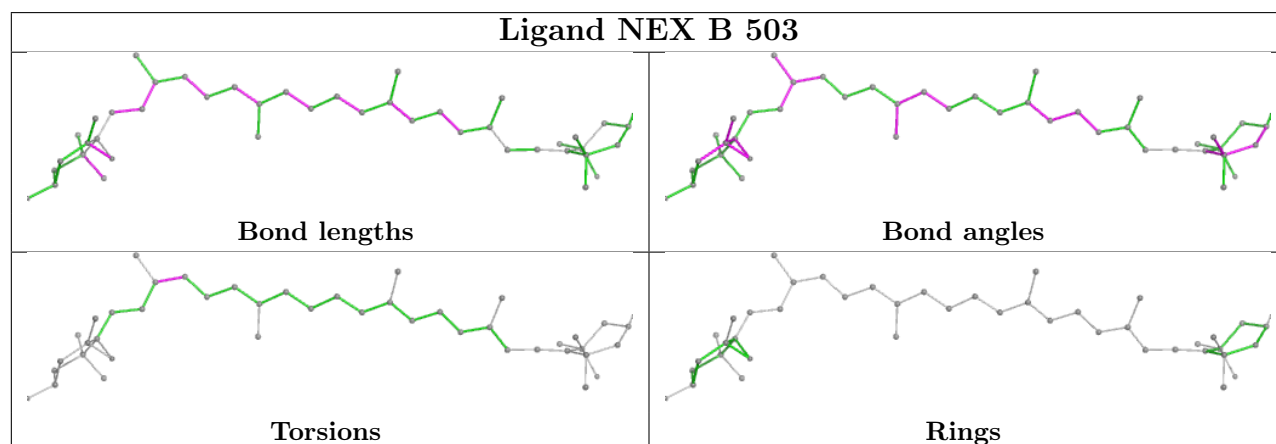
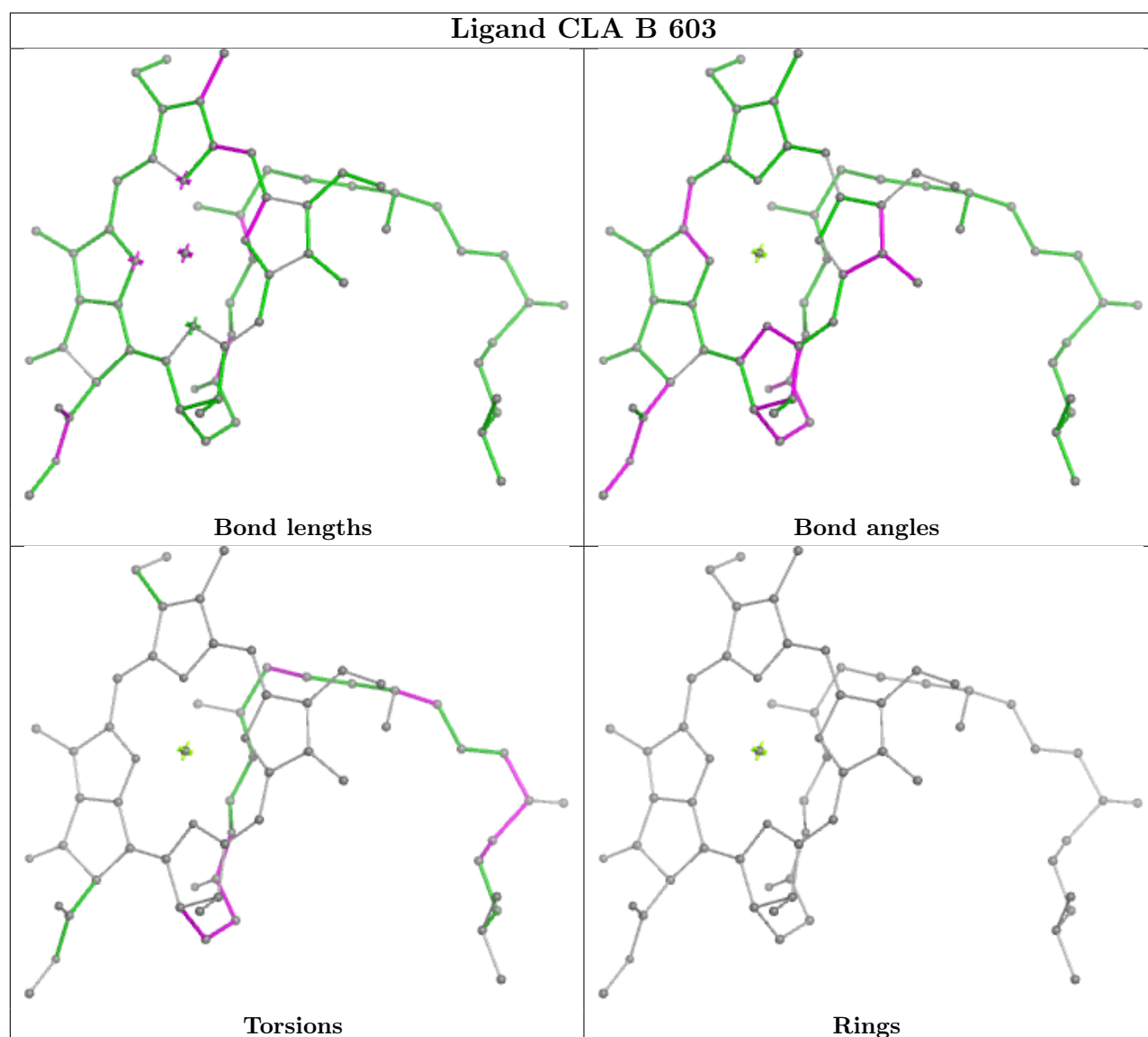




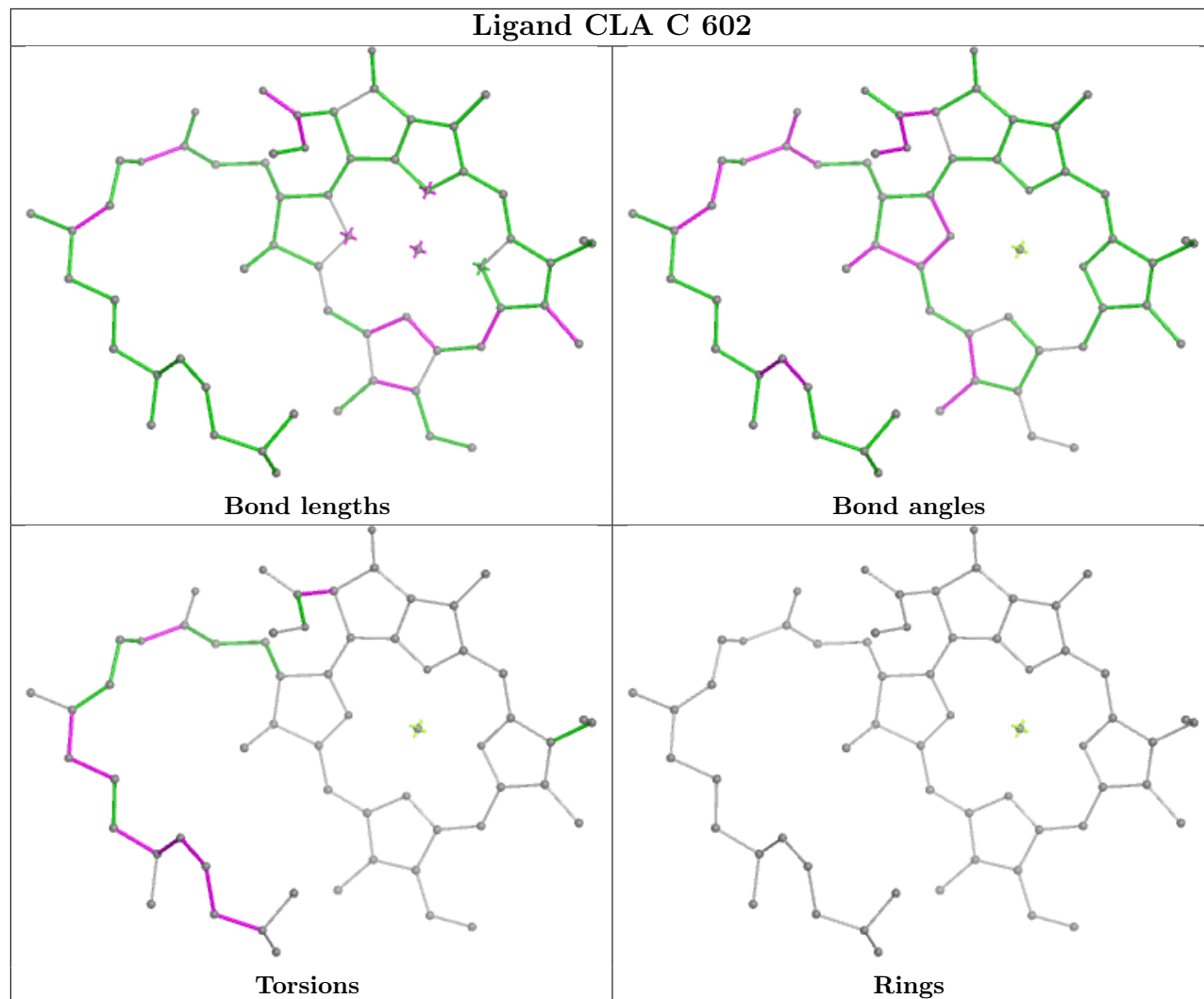


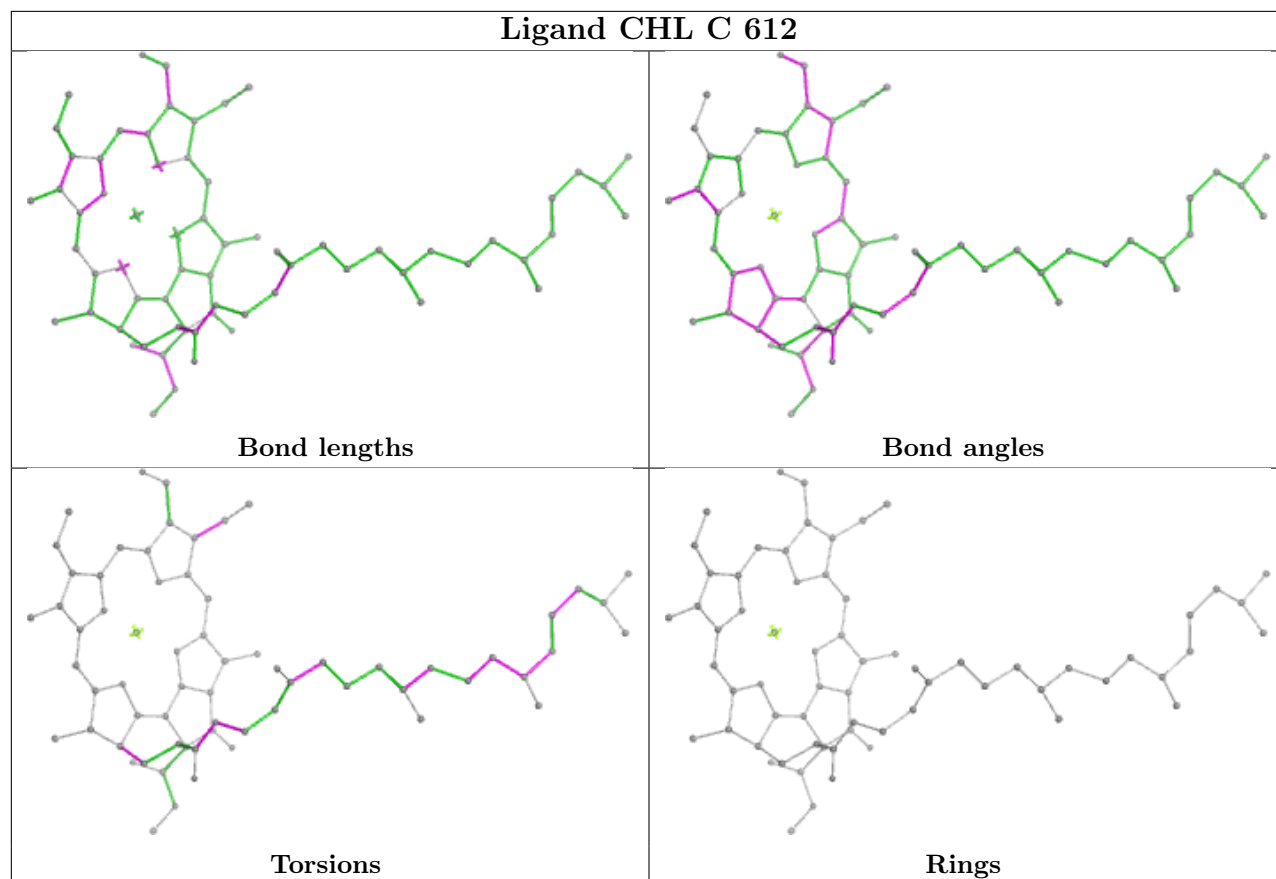


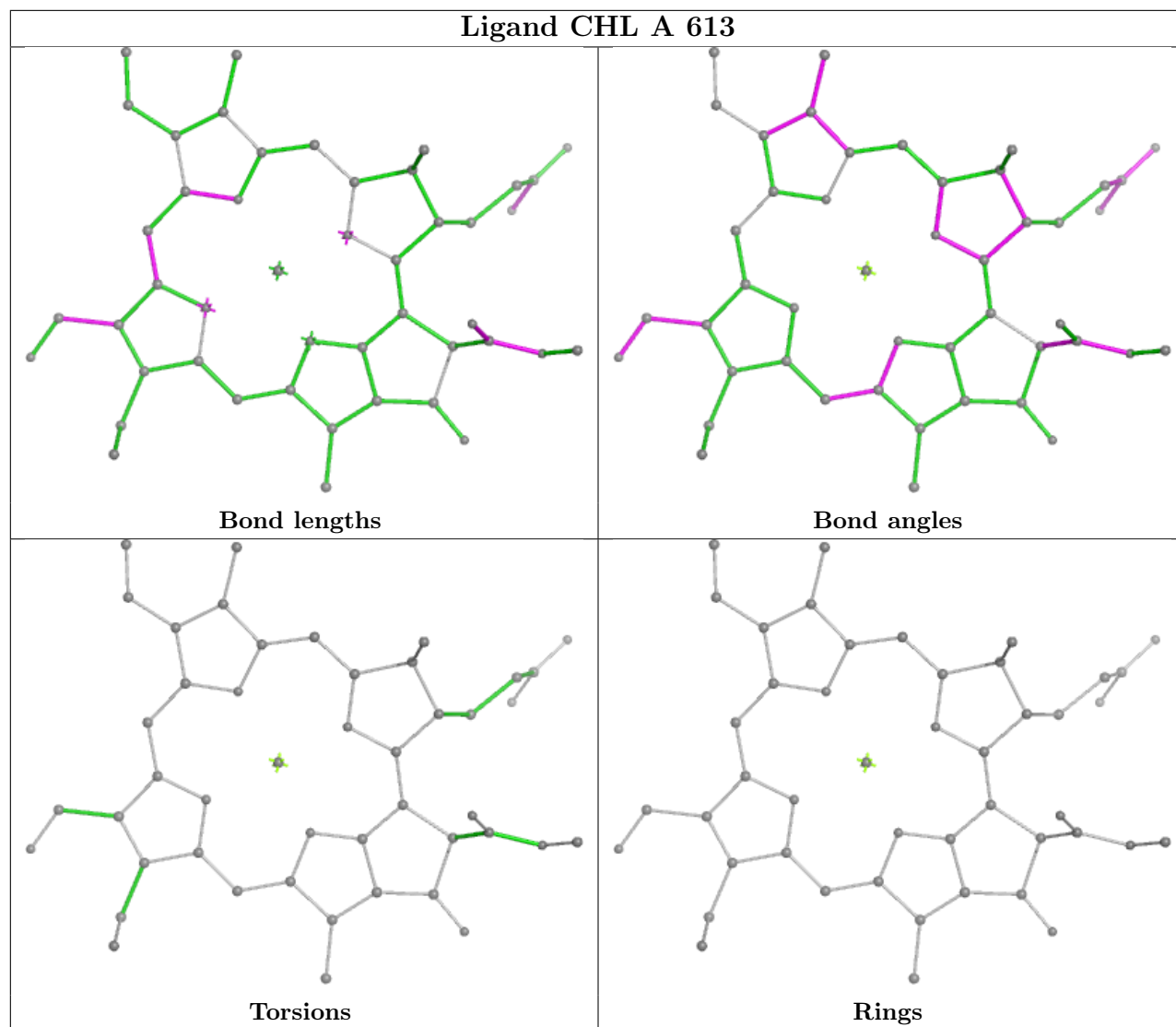


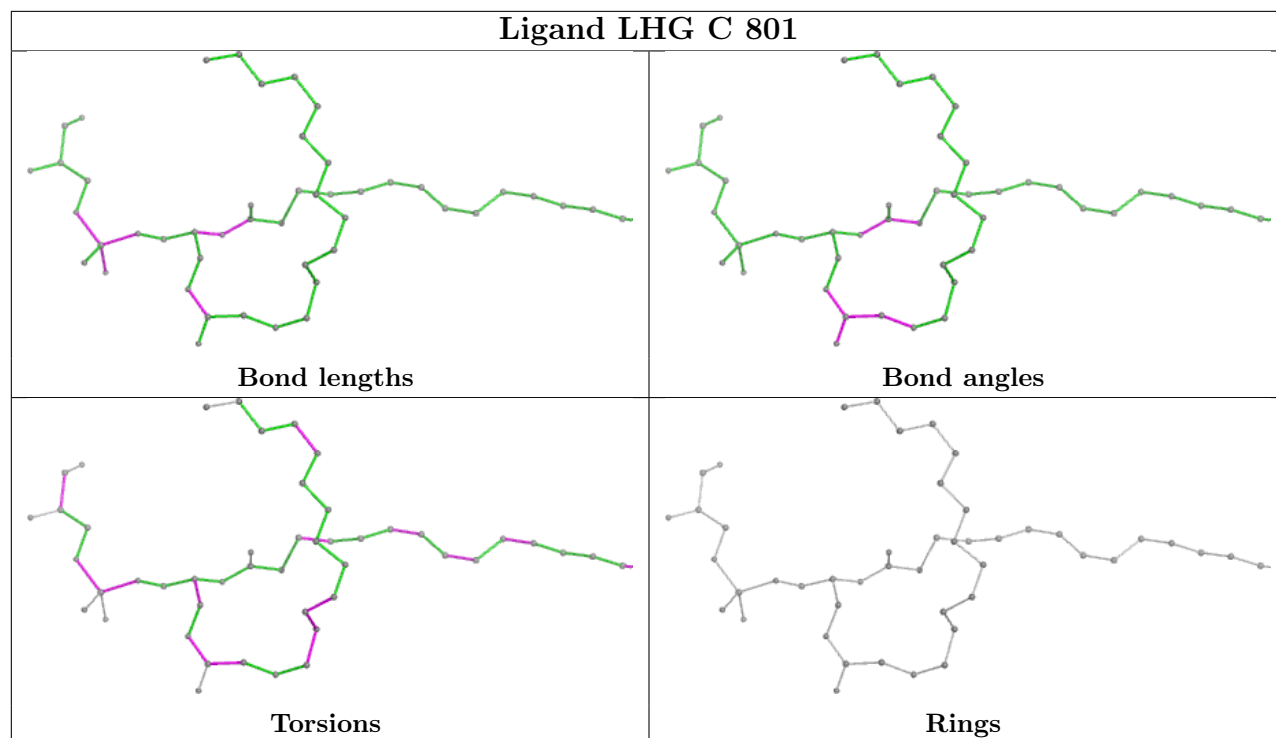


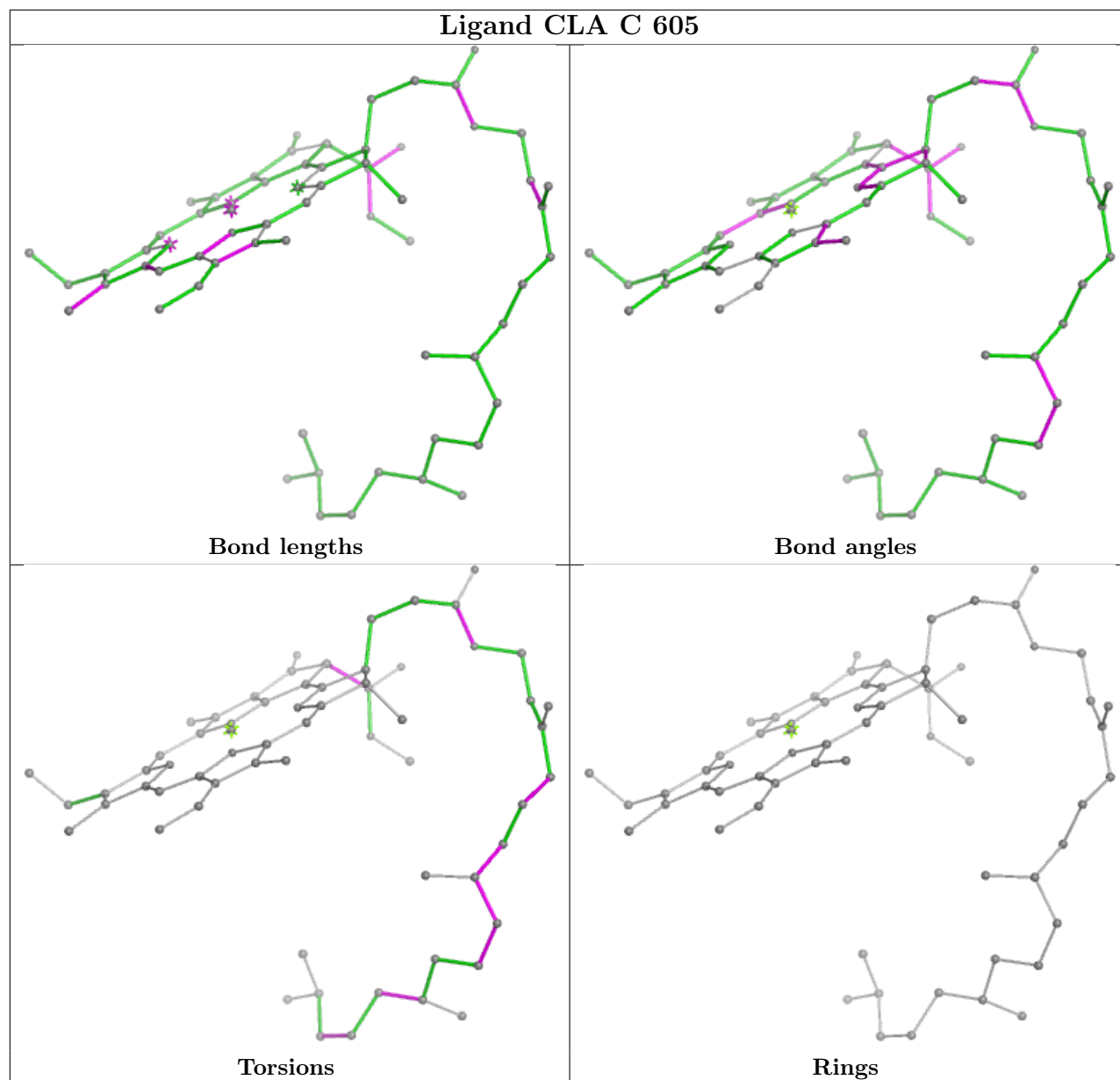
Ligand CLA C 602

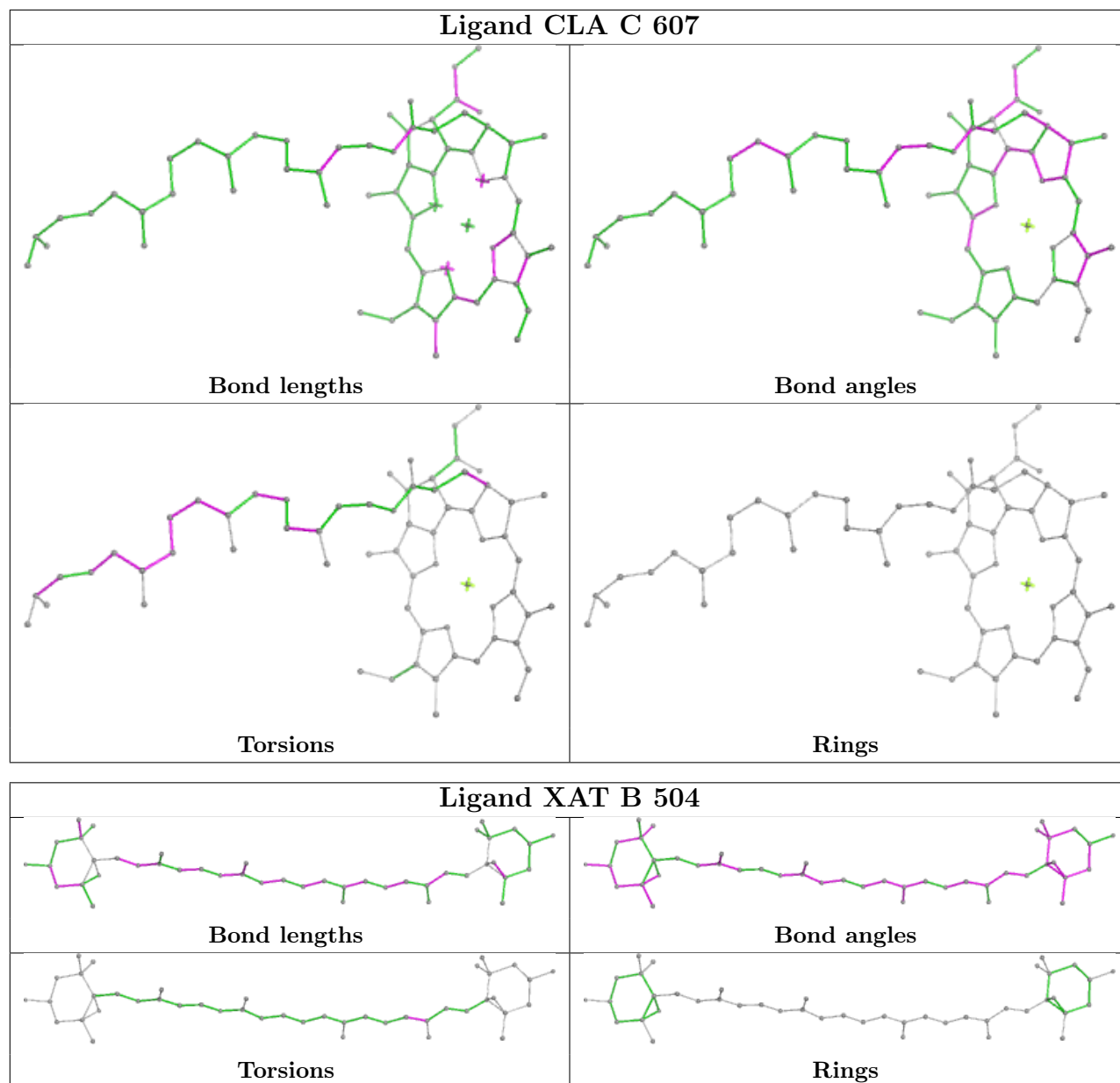


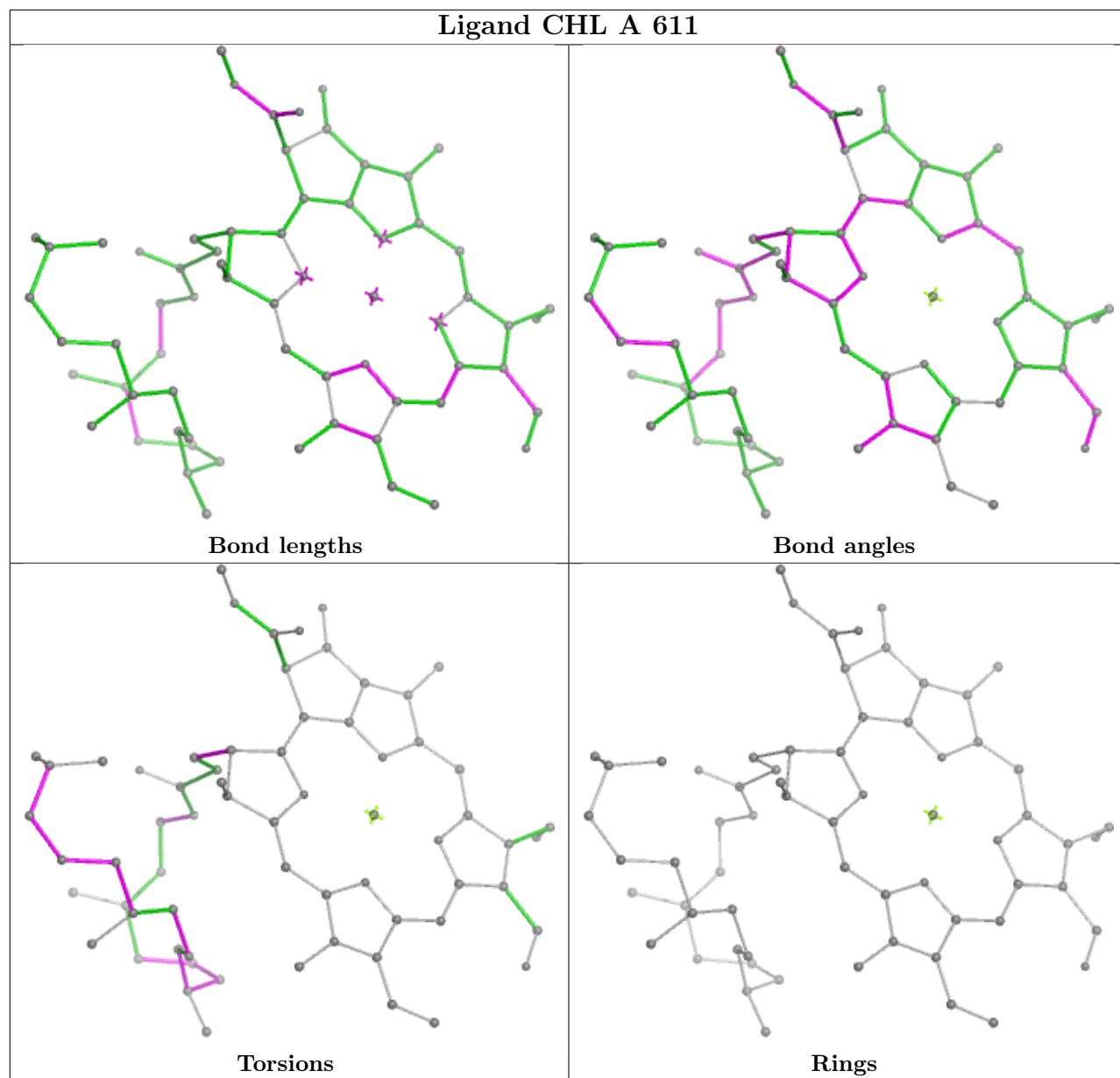


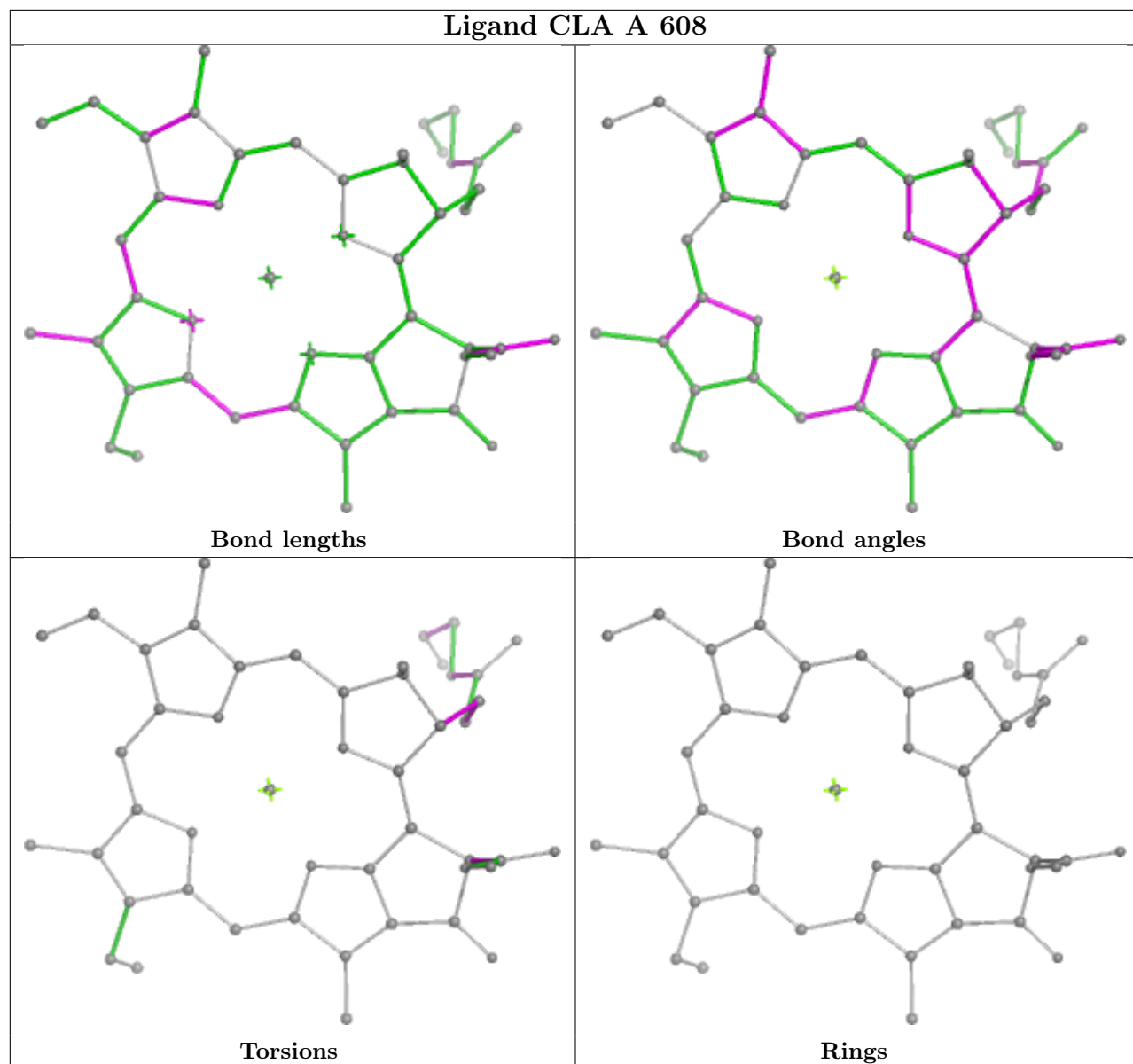


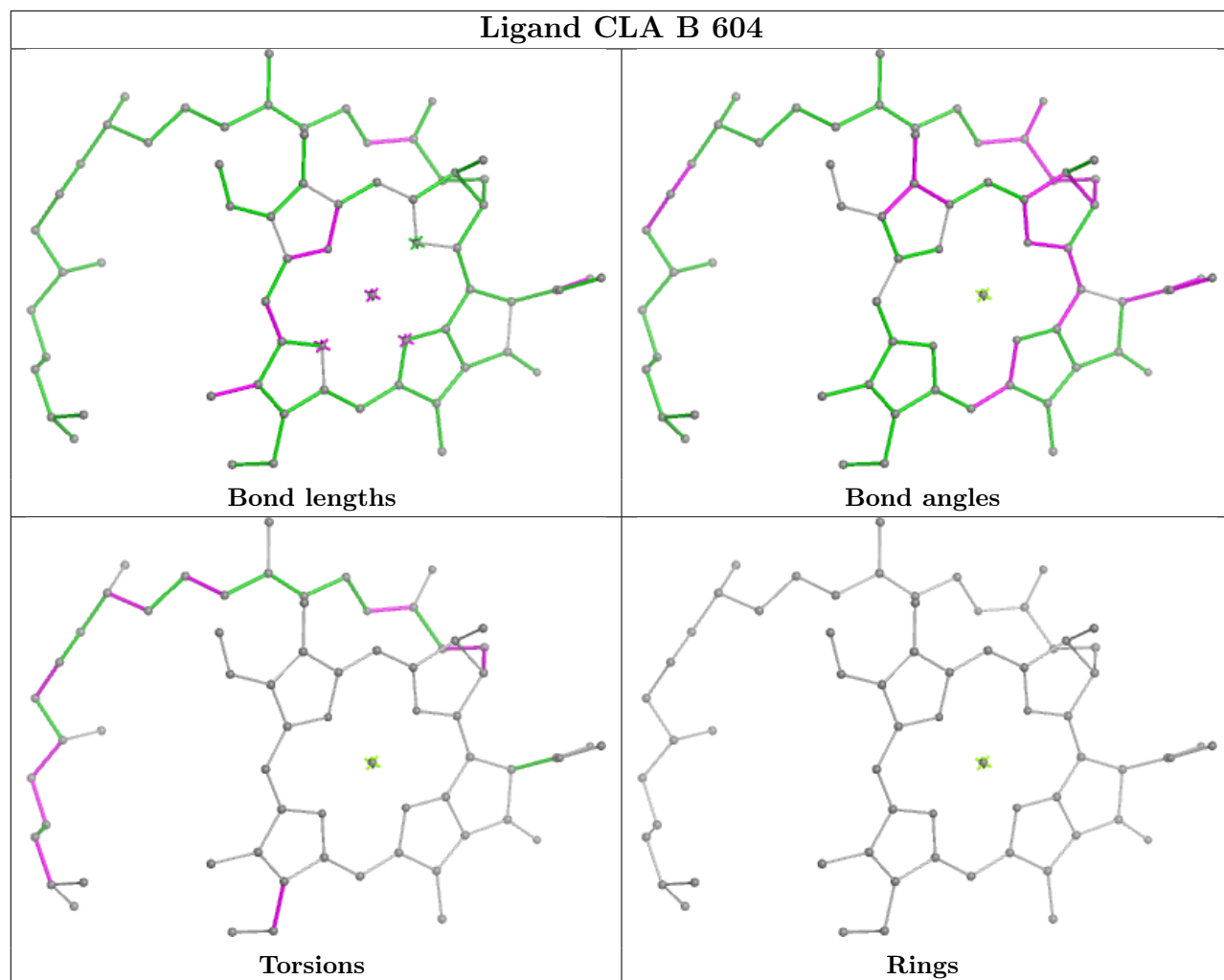


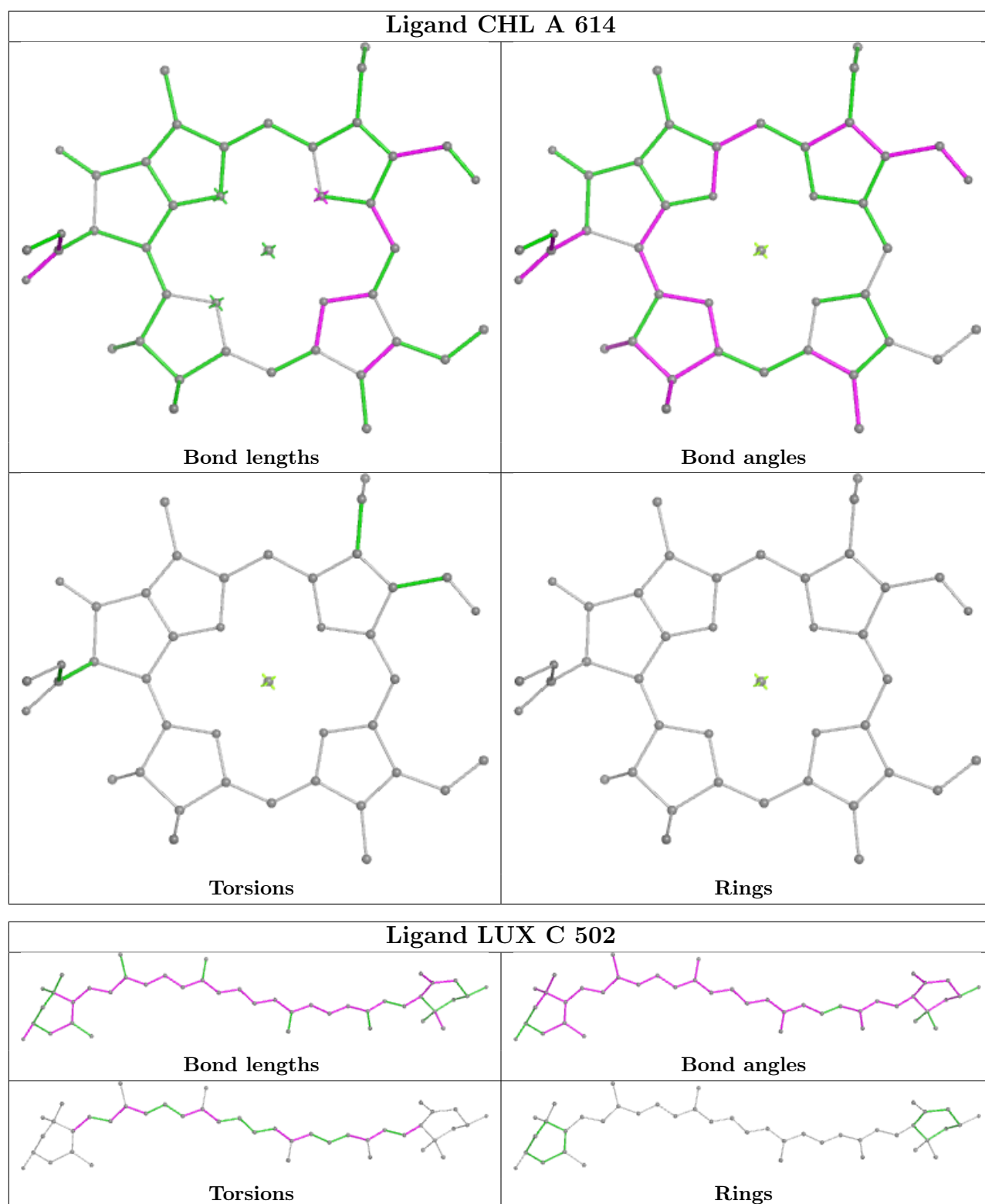


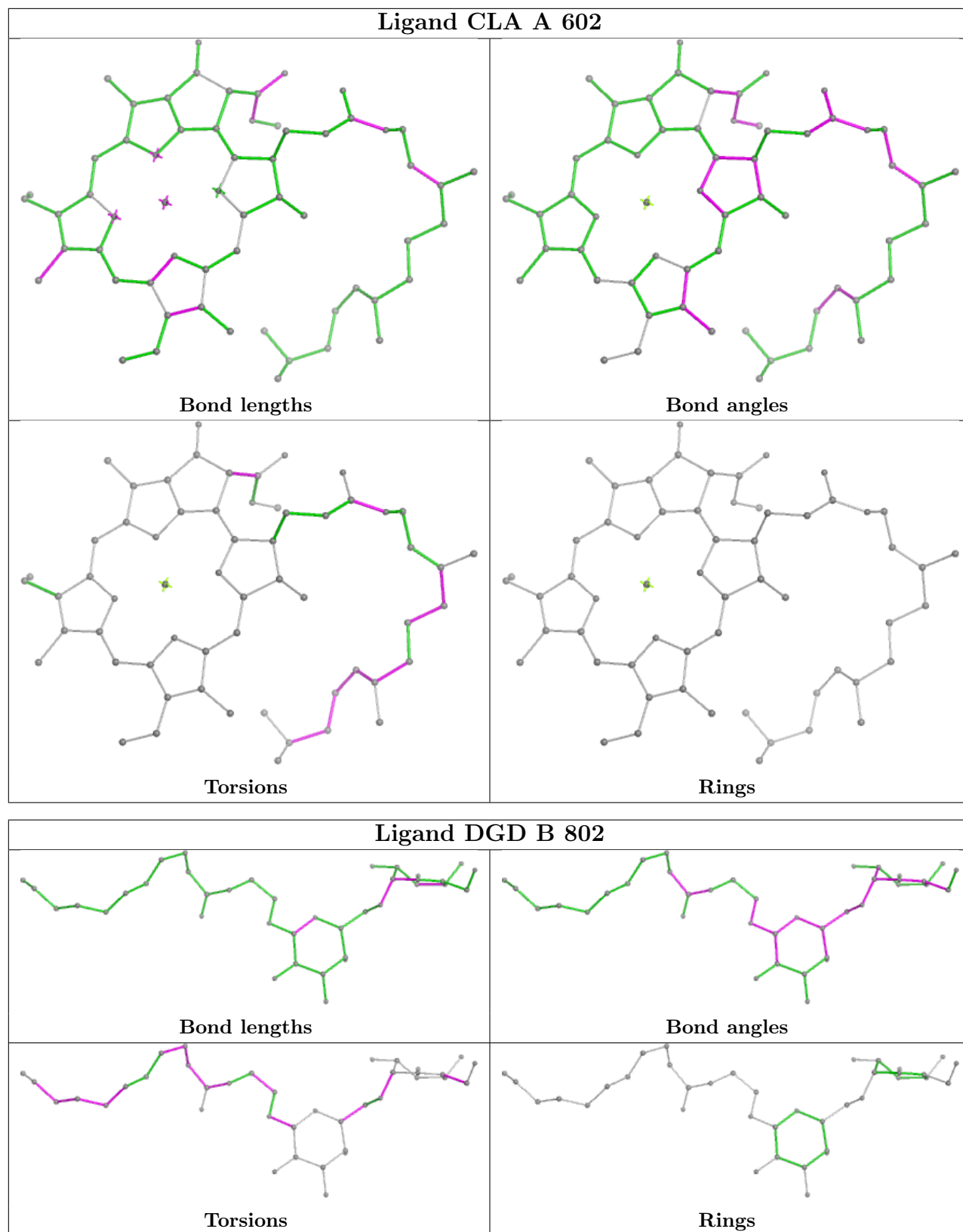


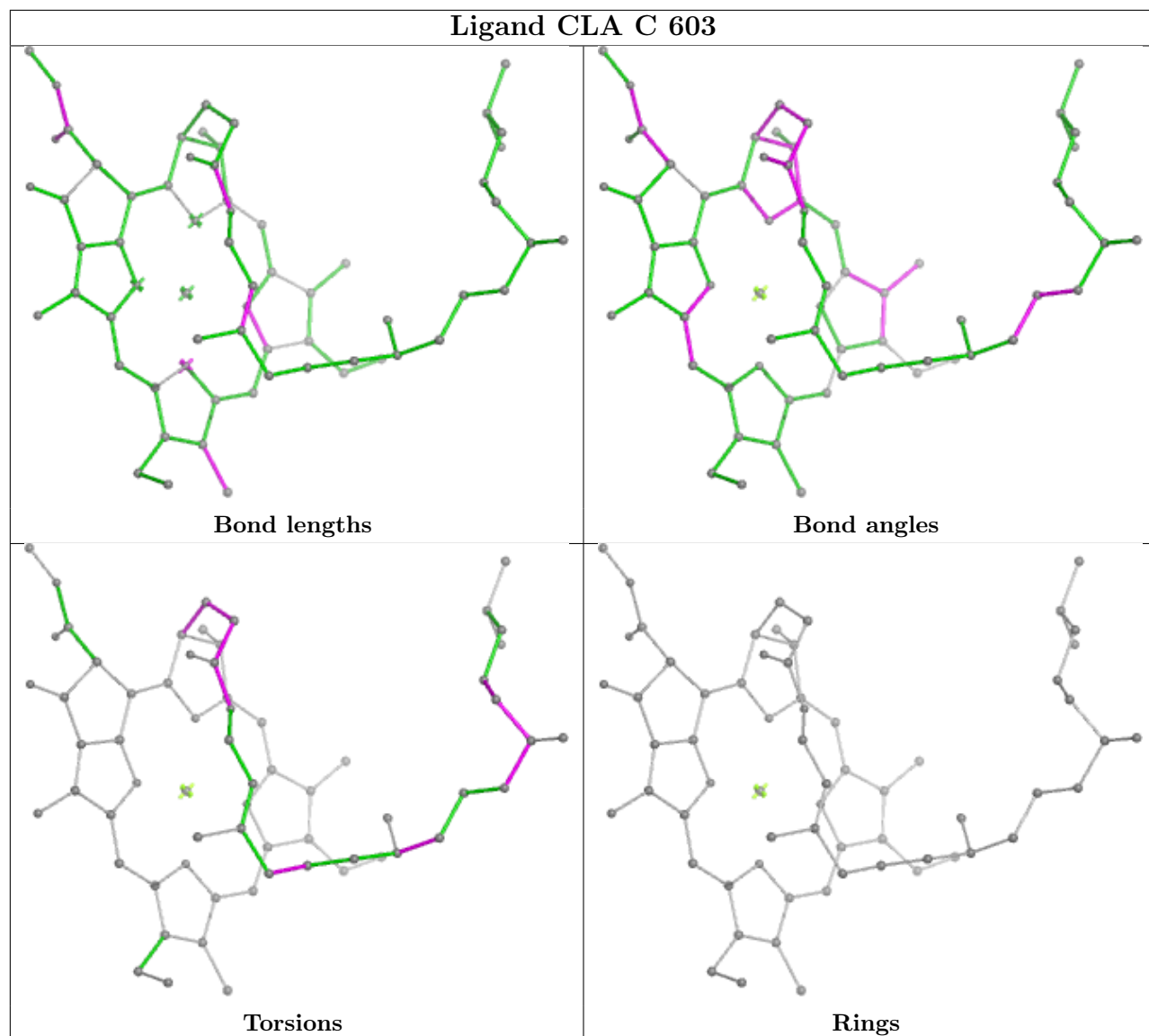


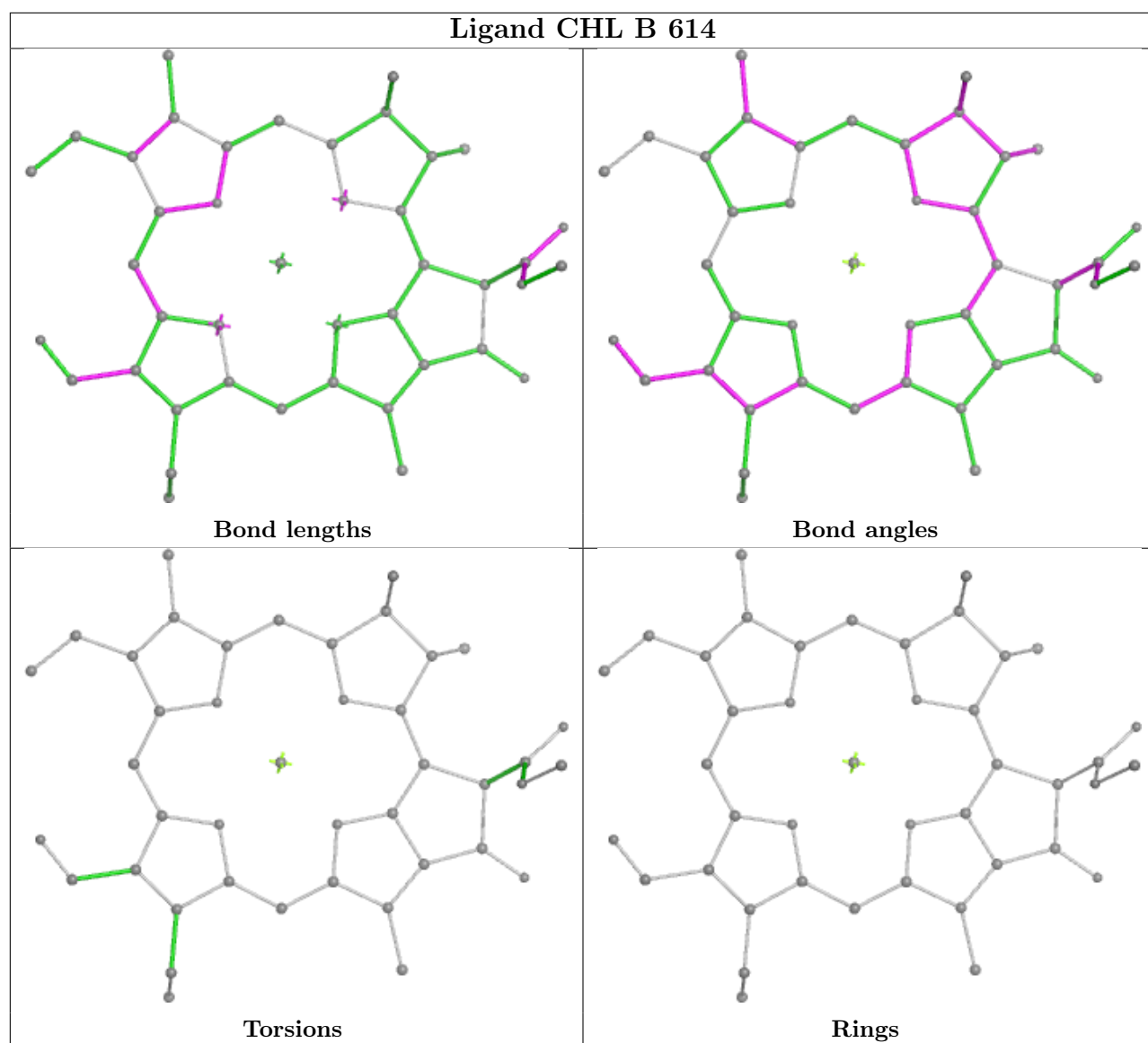


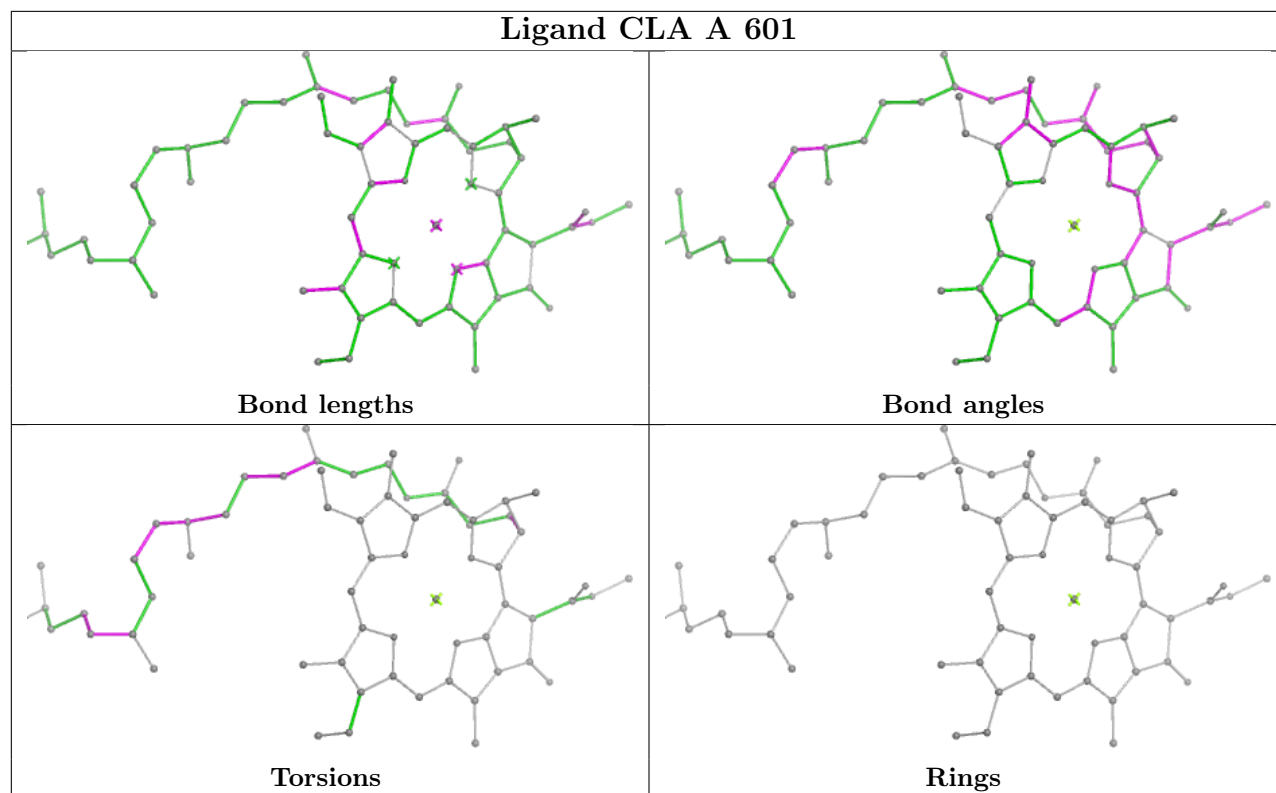


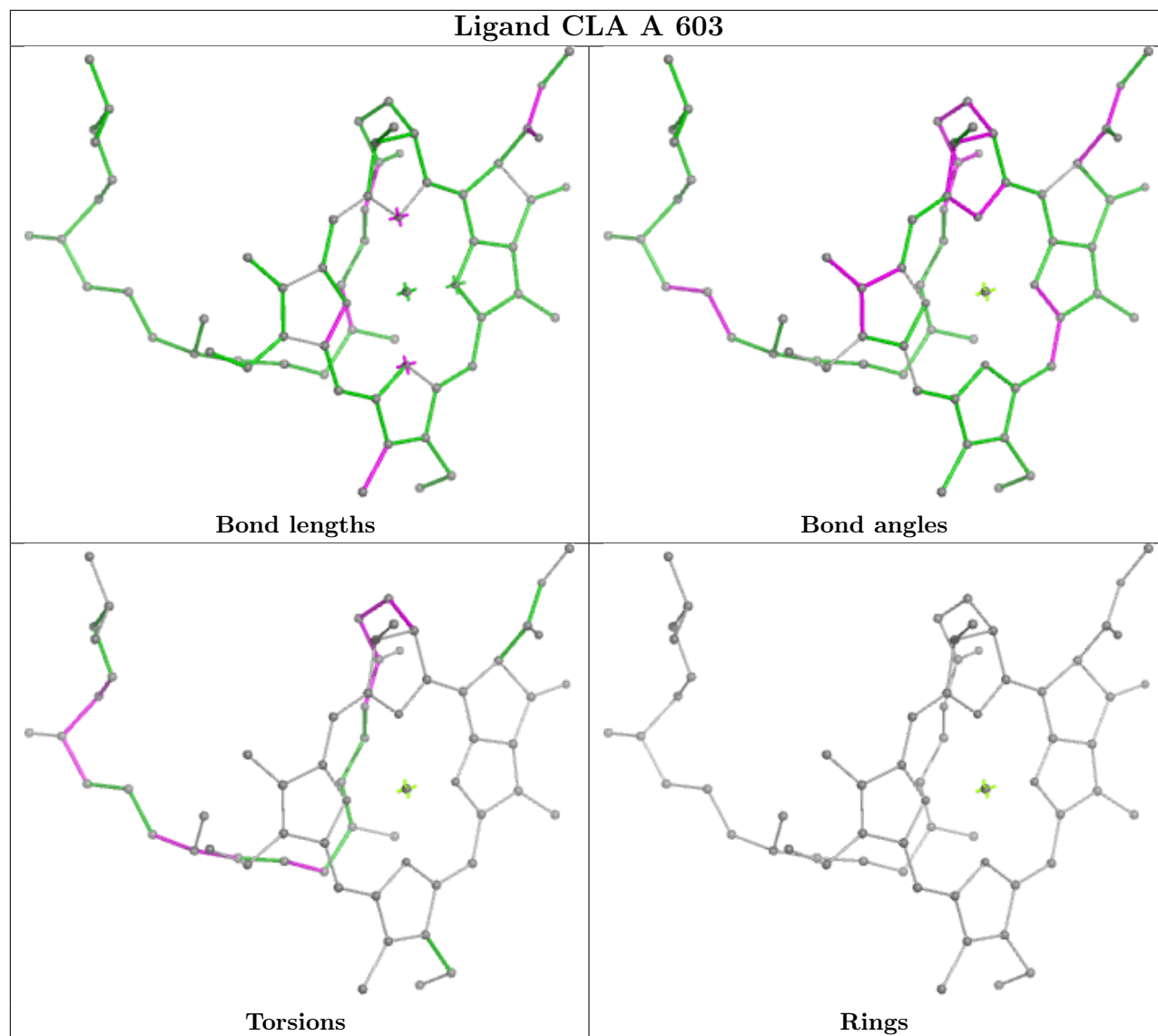


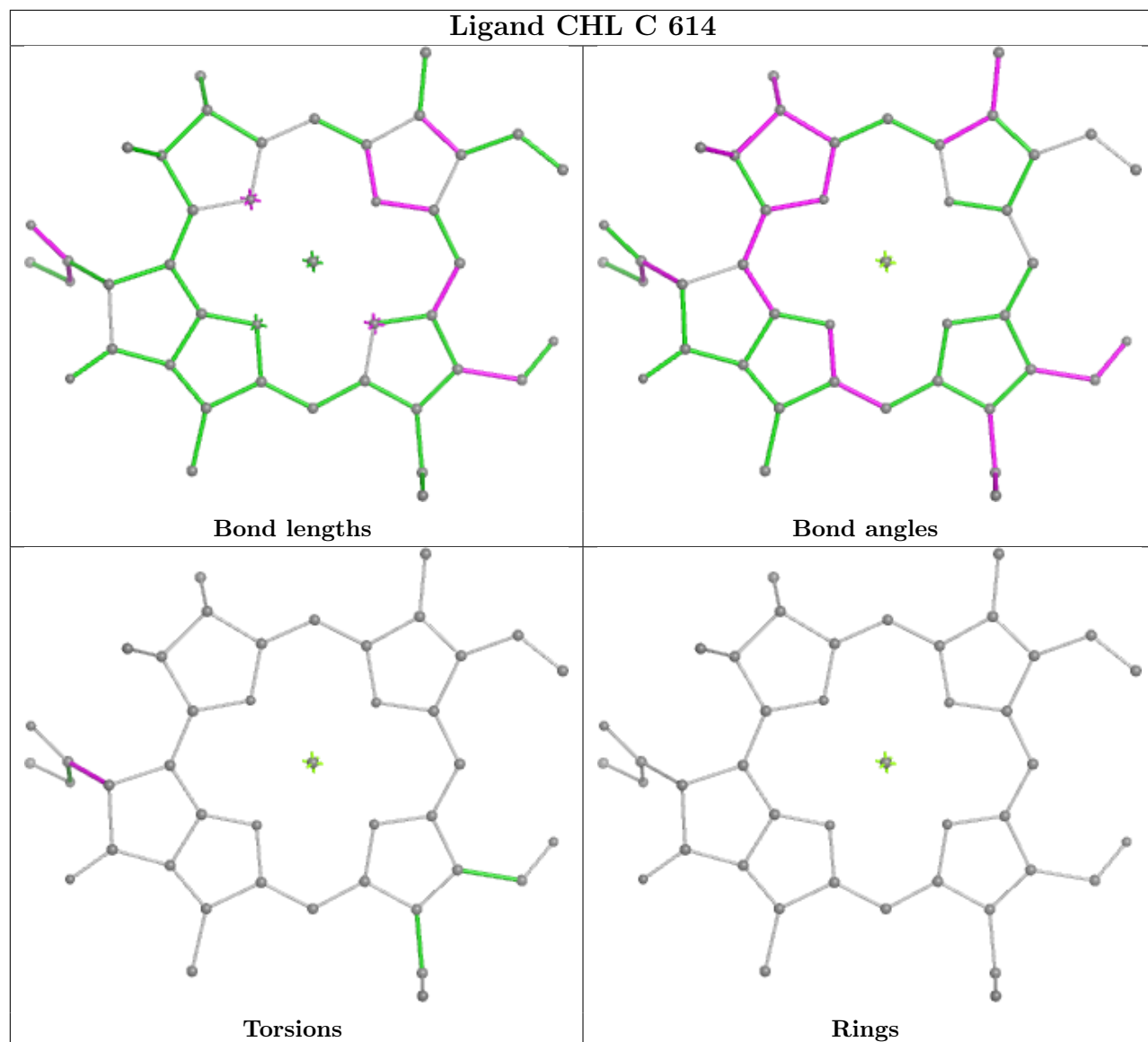


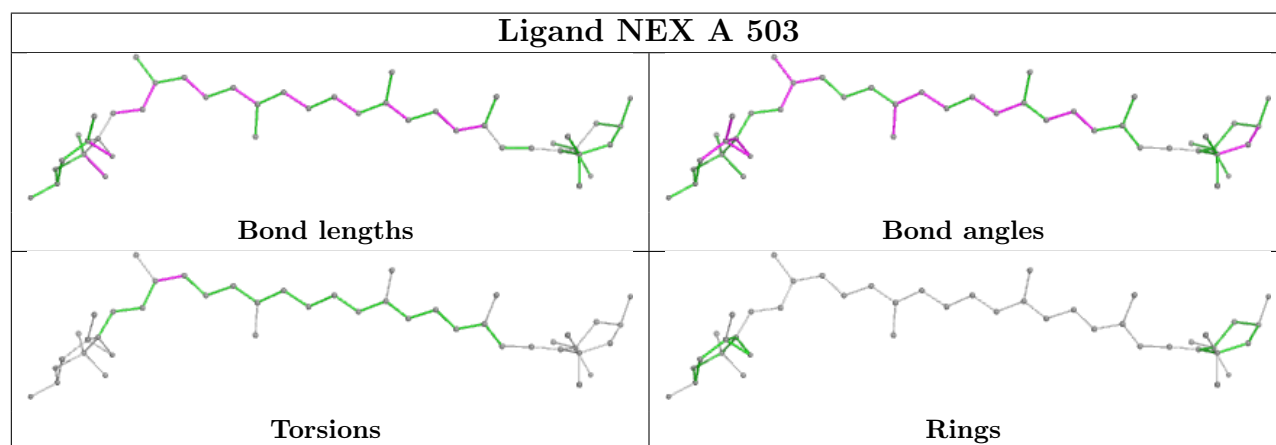
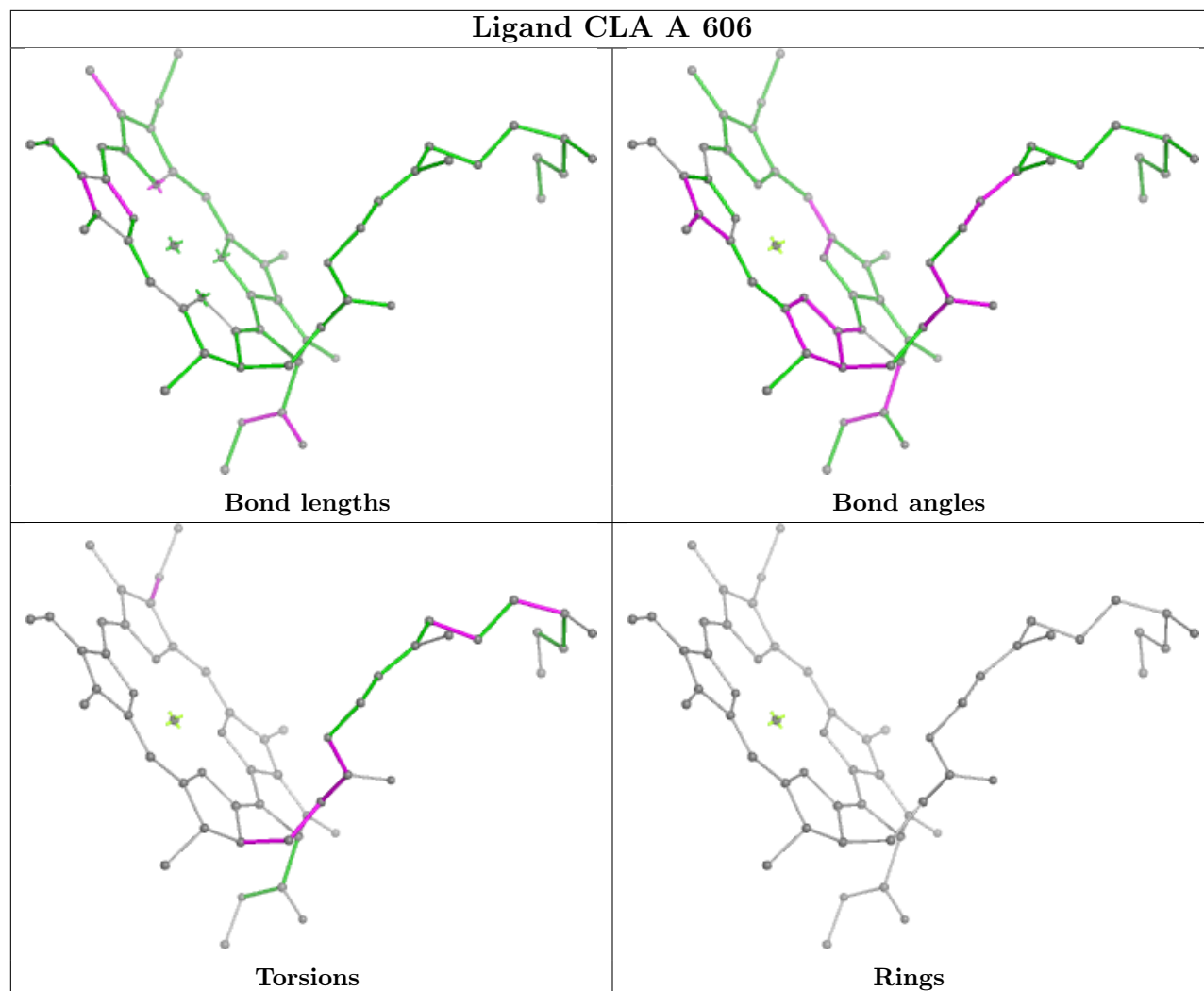


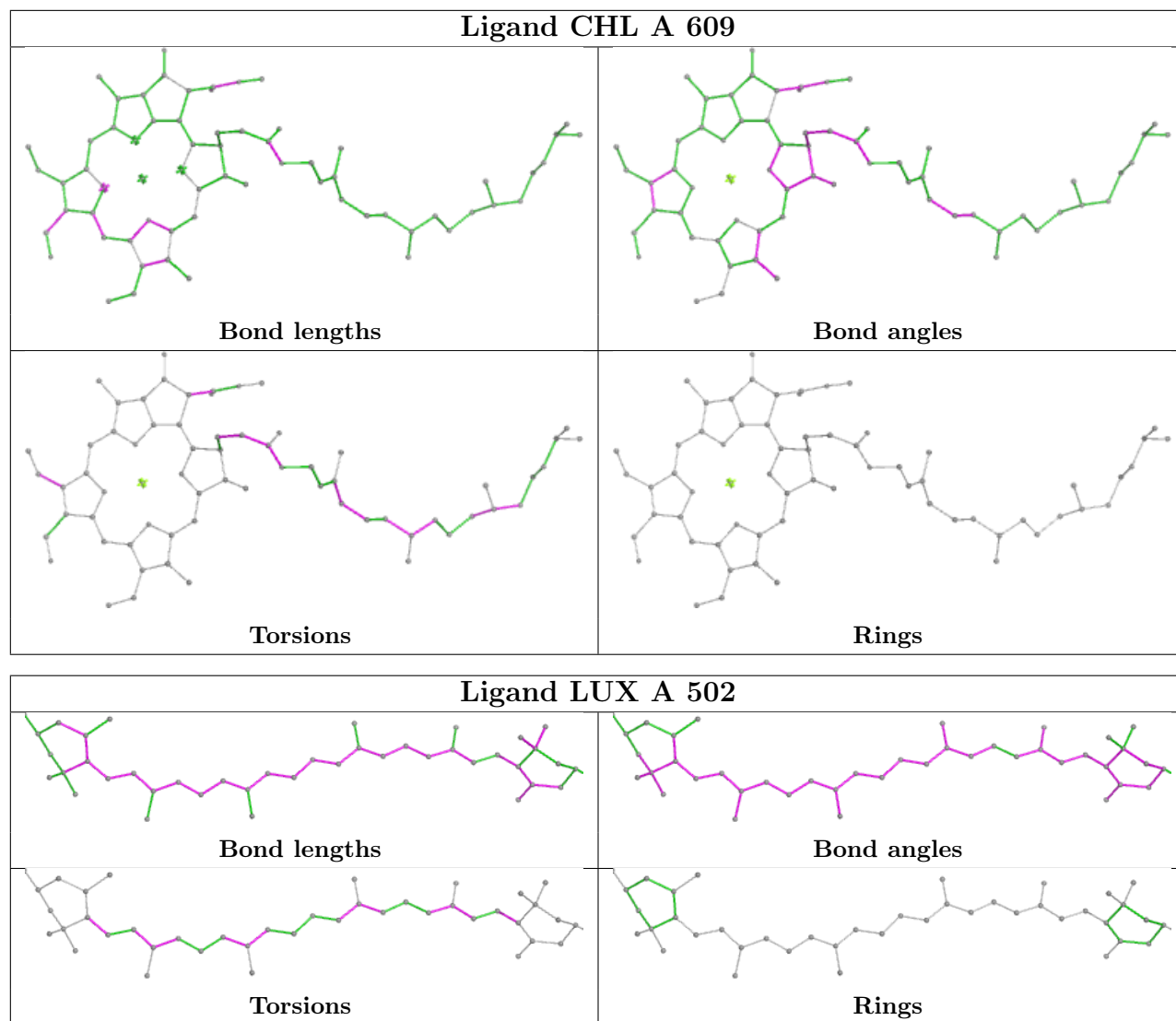


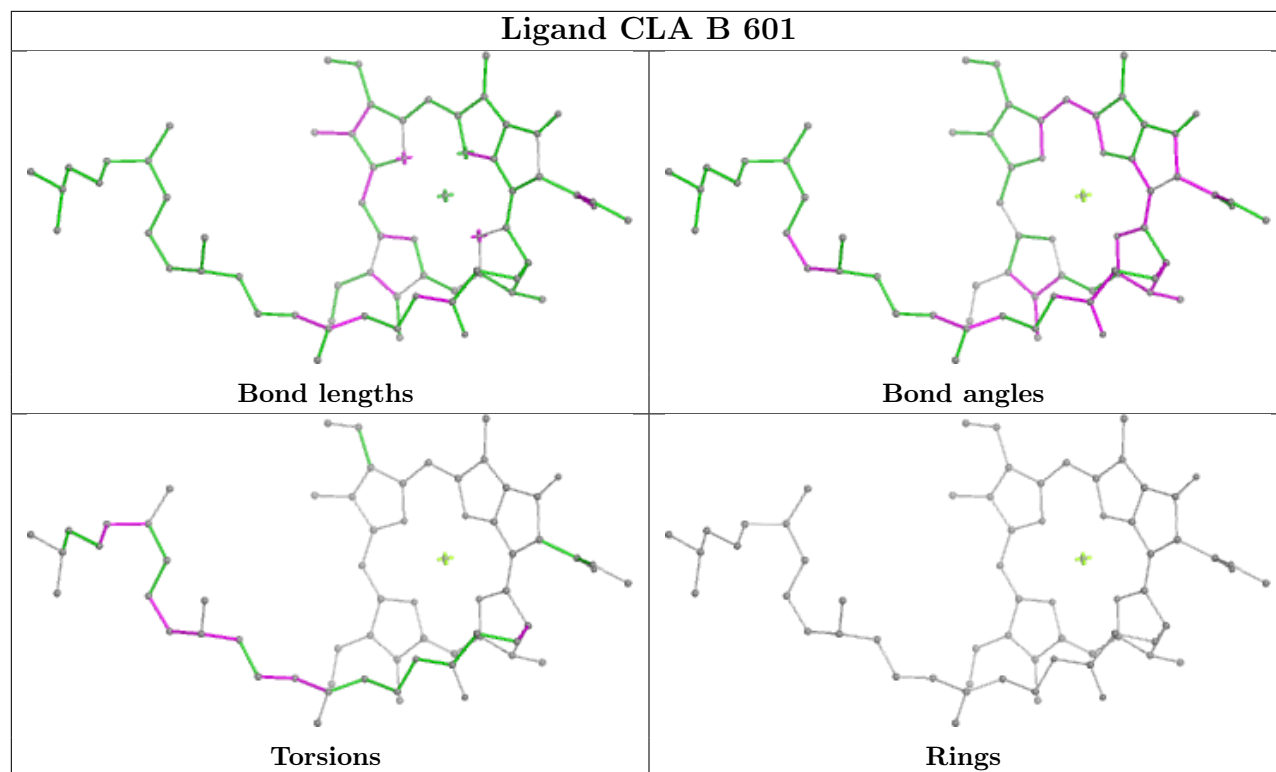


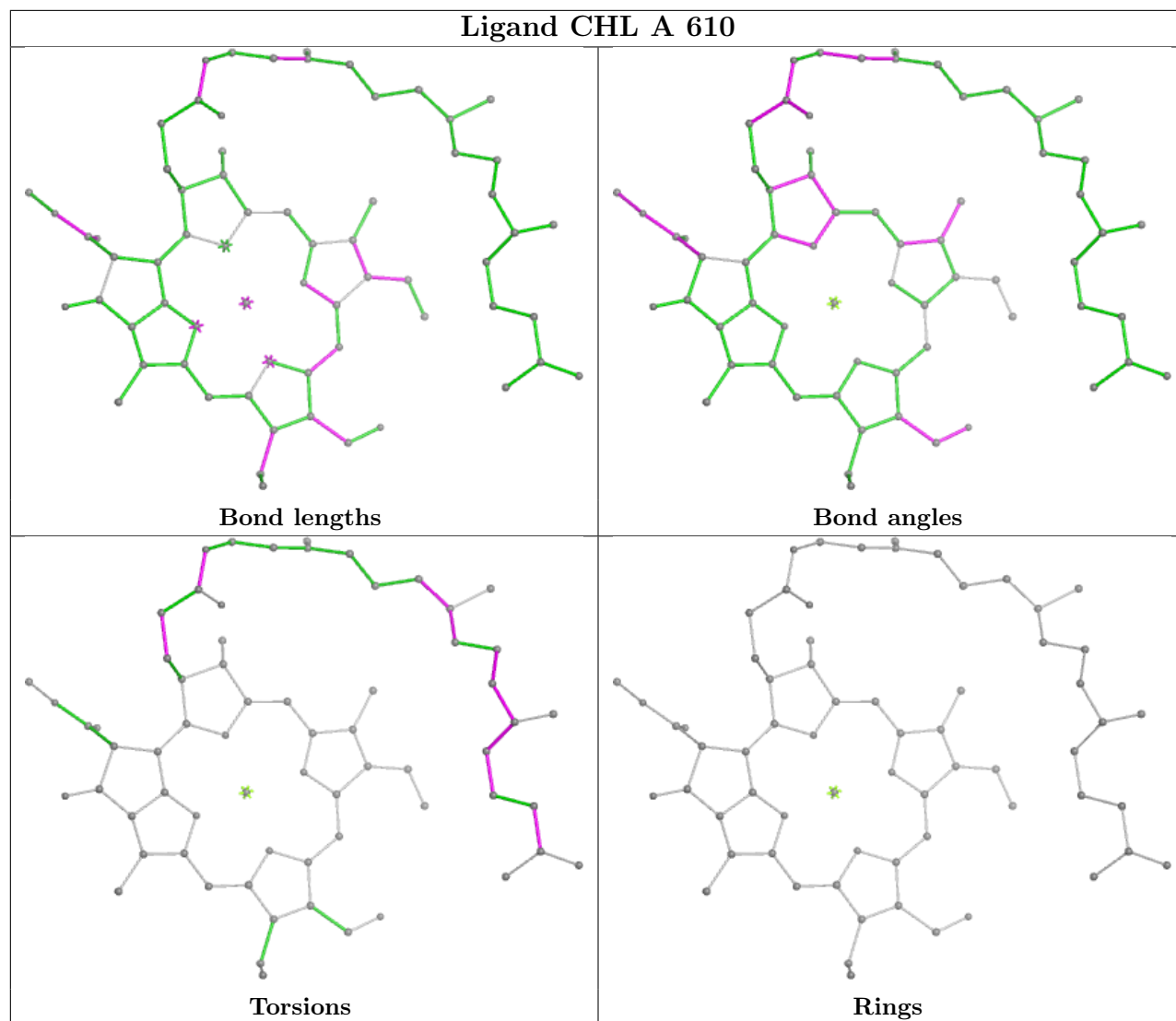


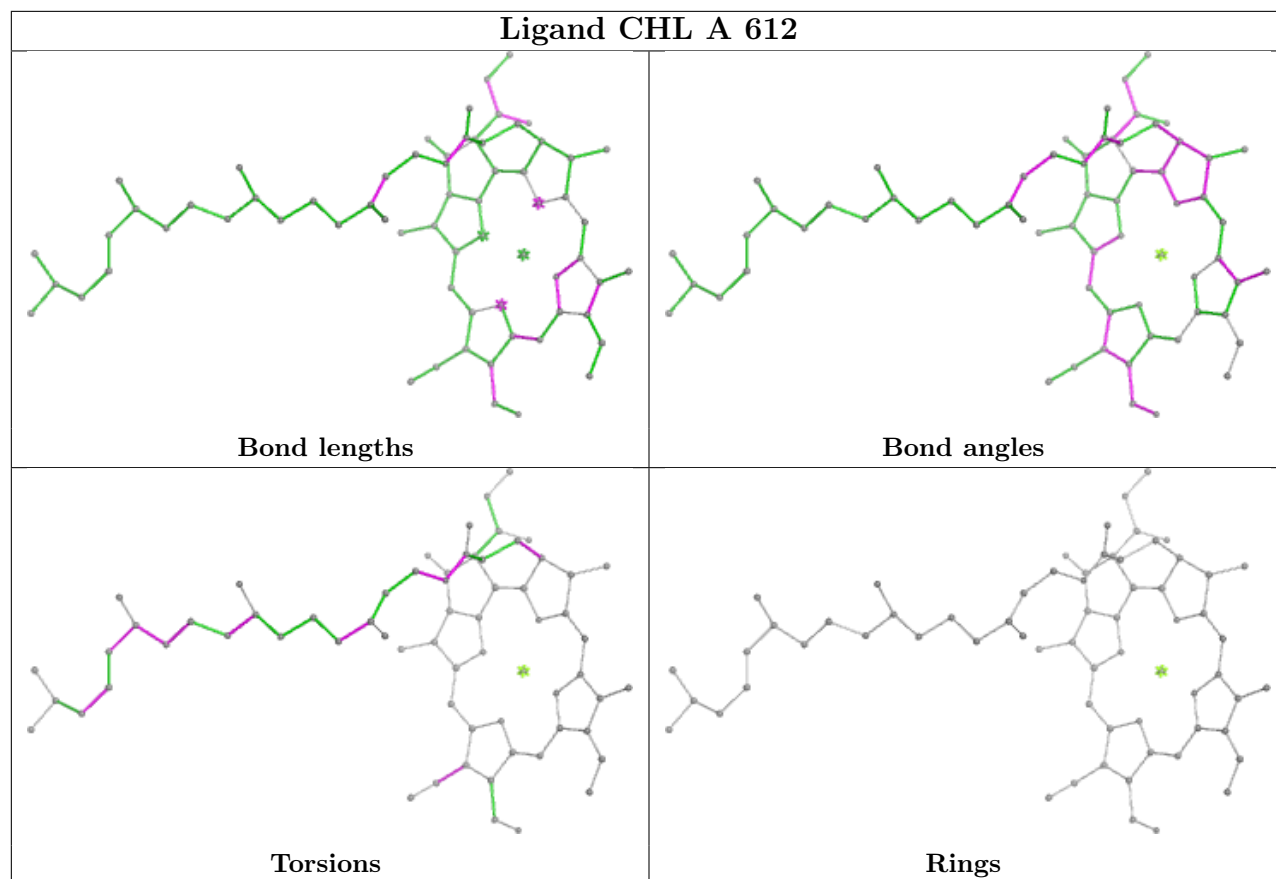


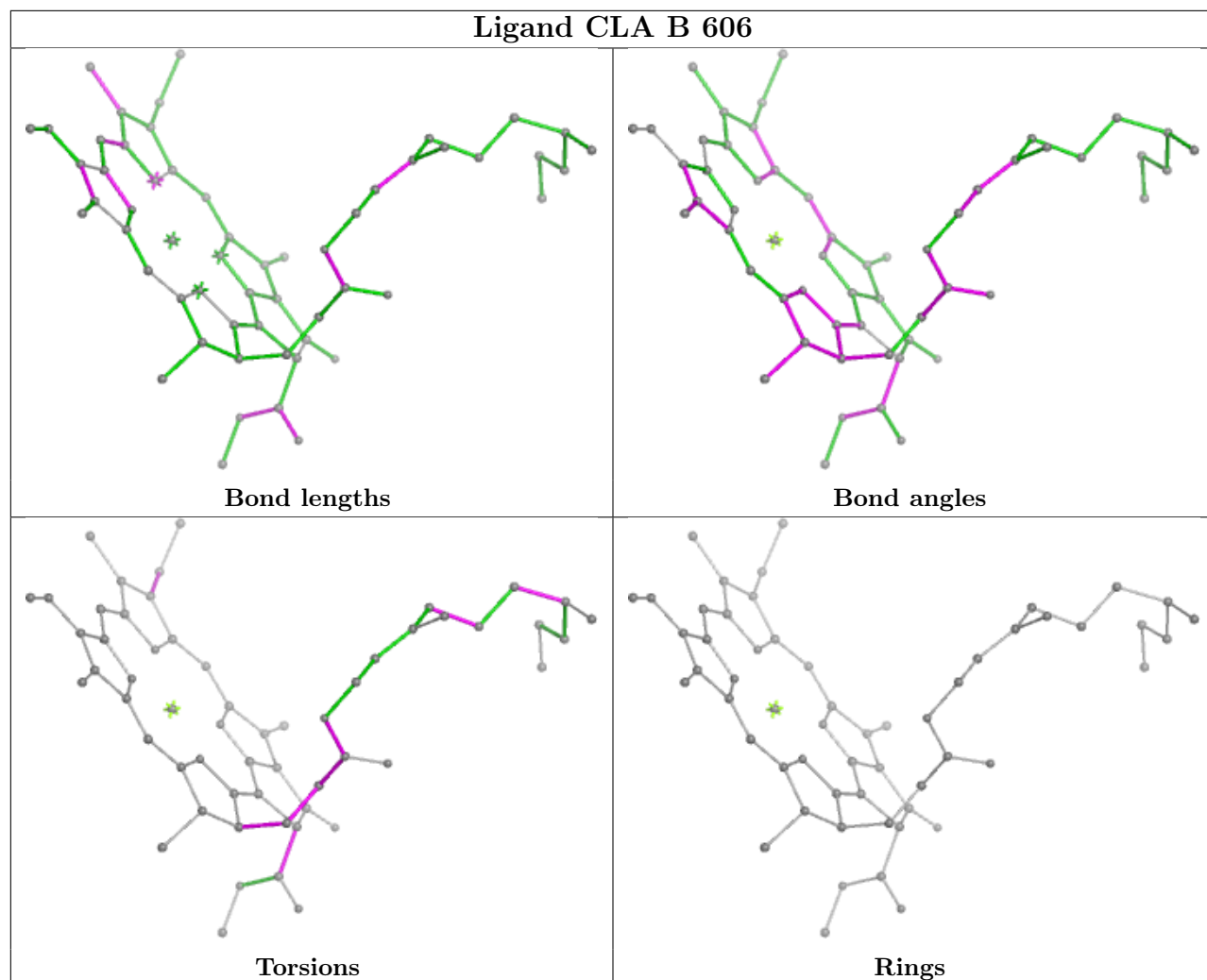


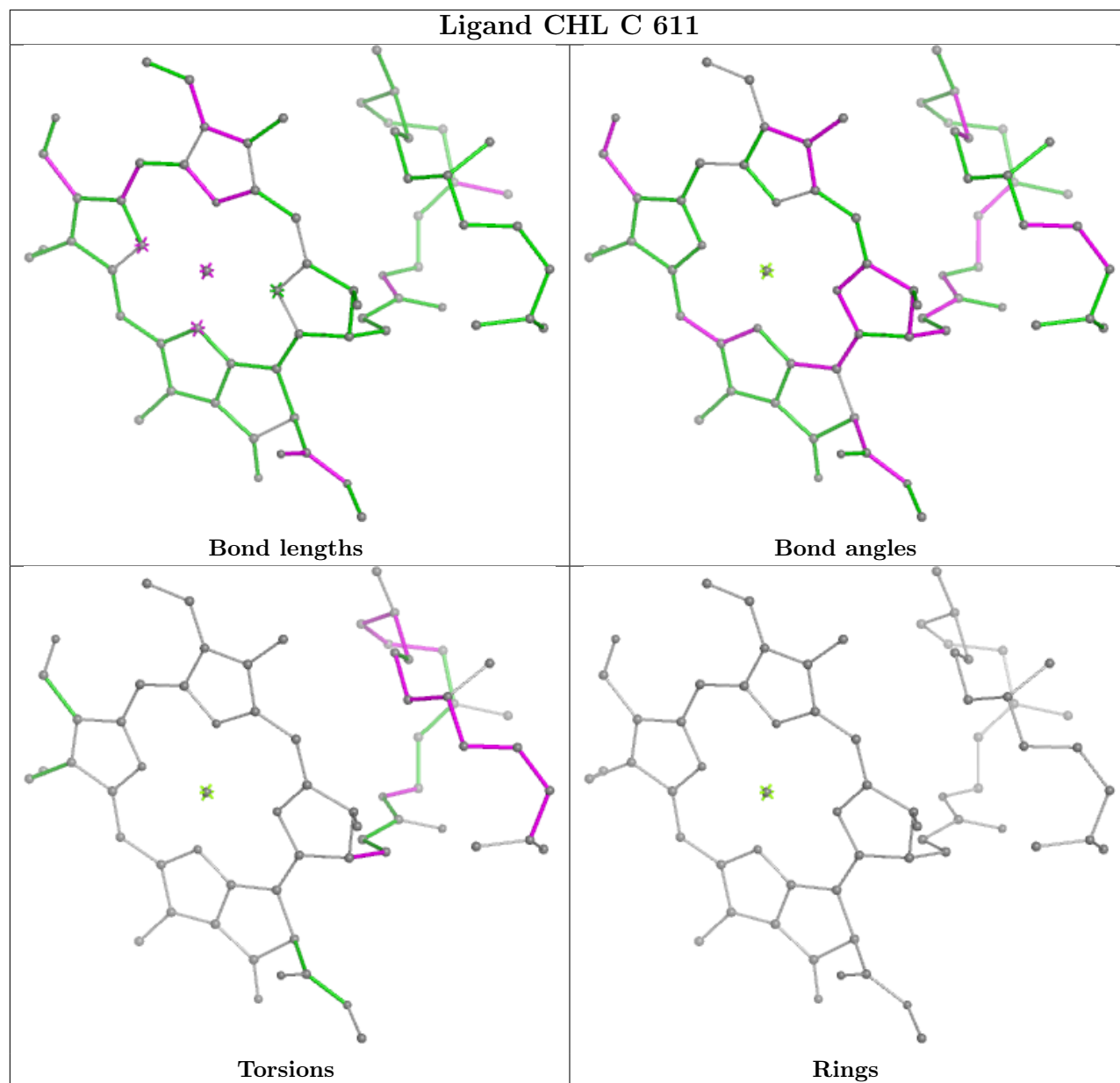


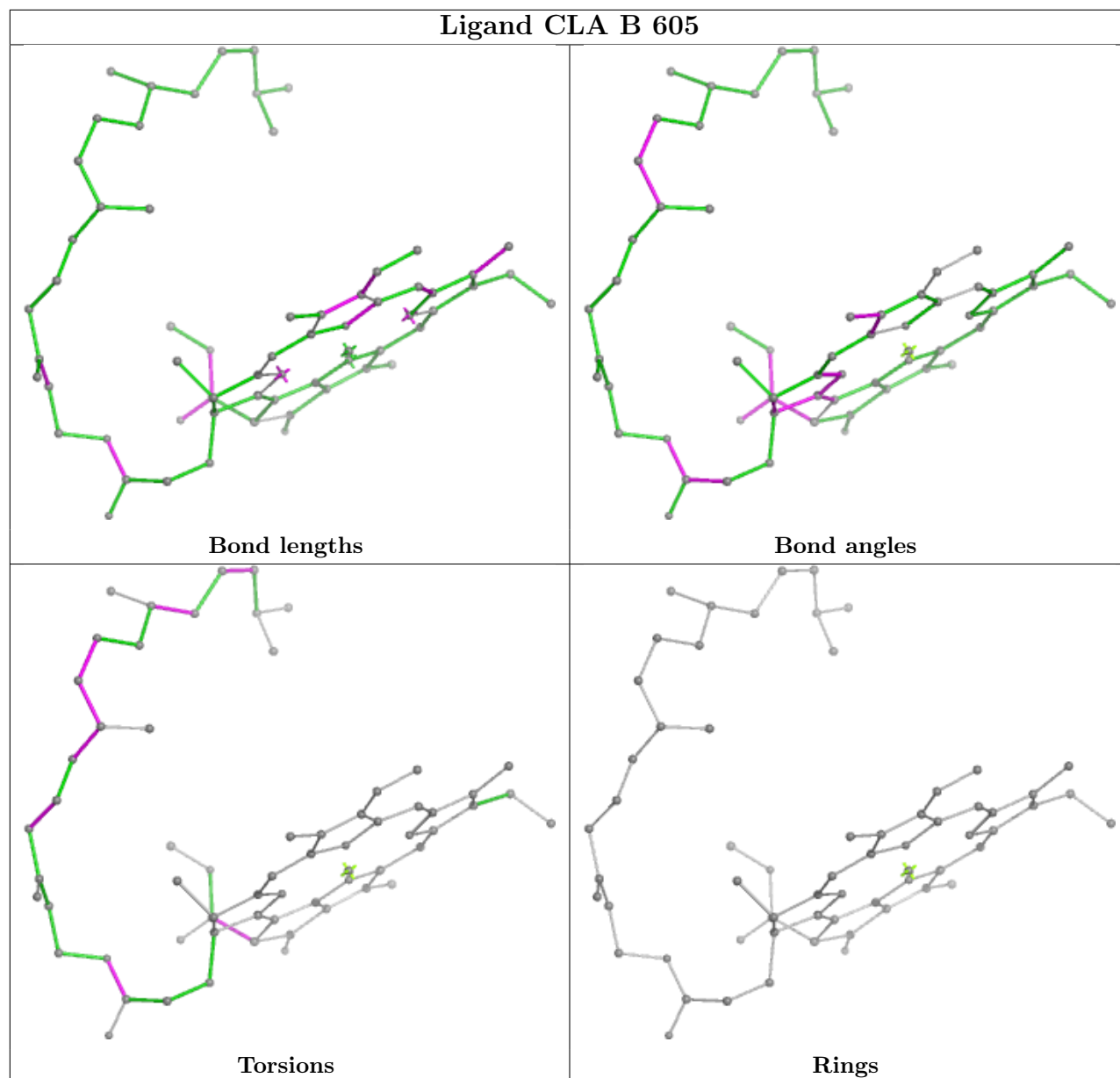


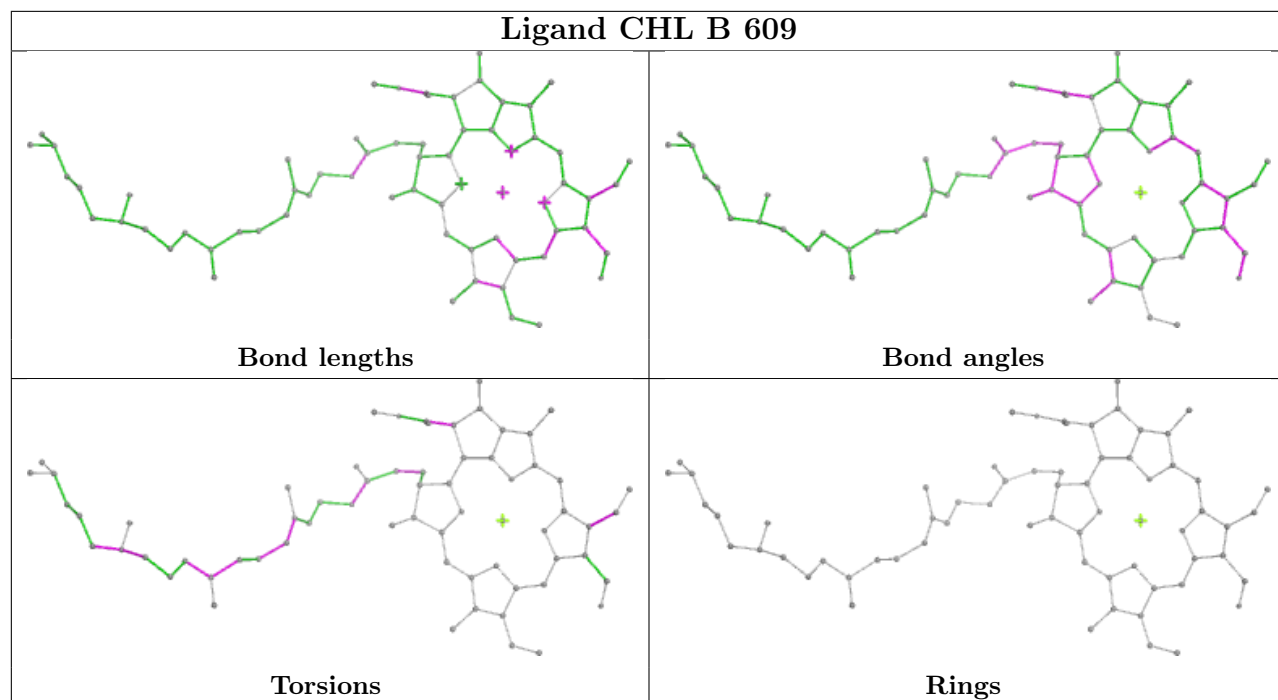


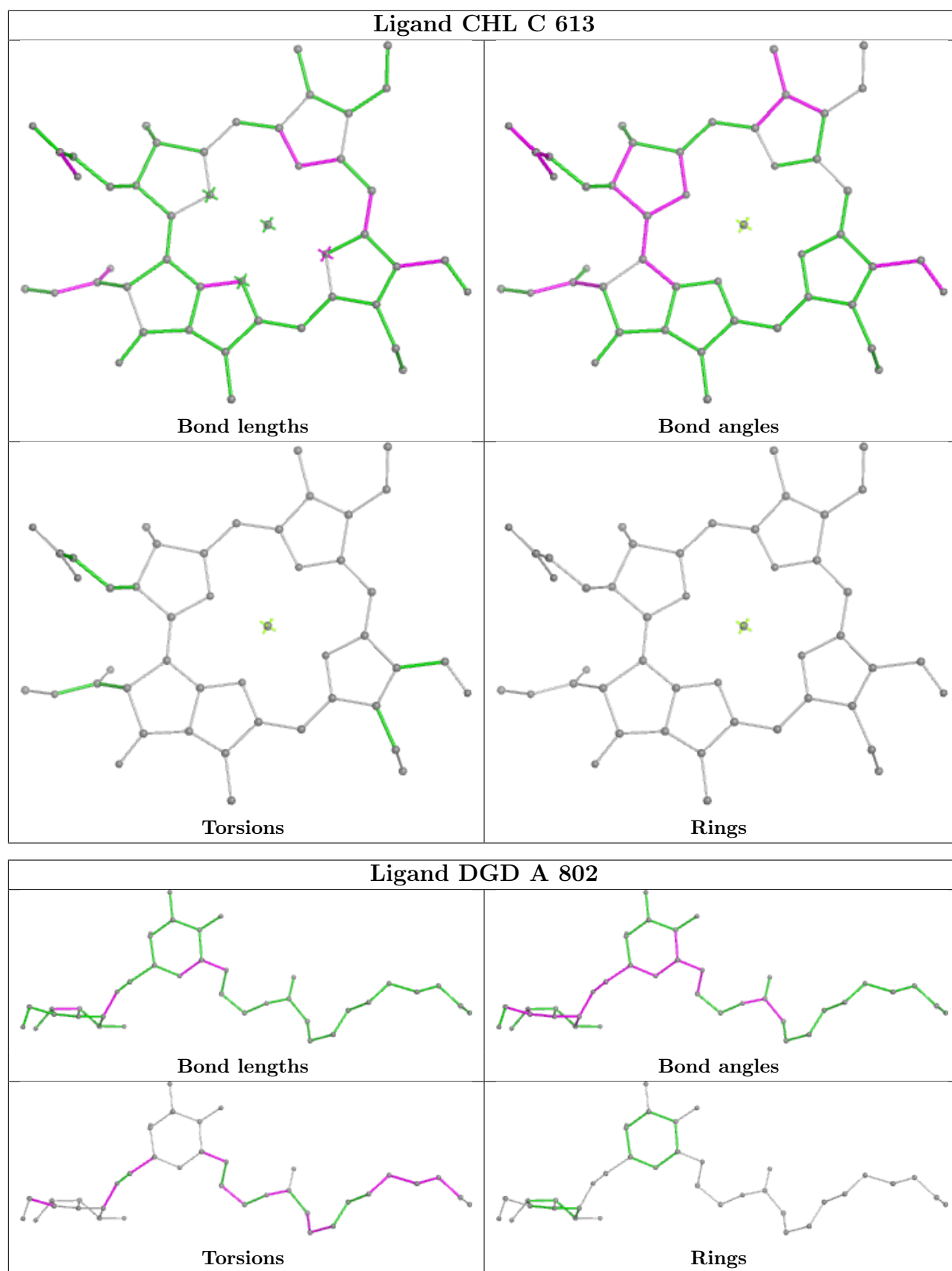


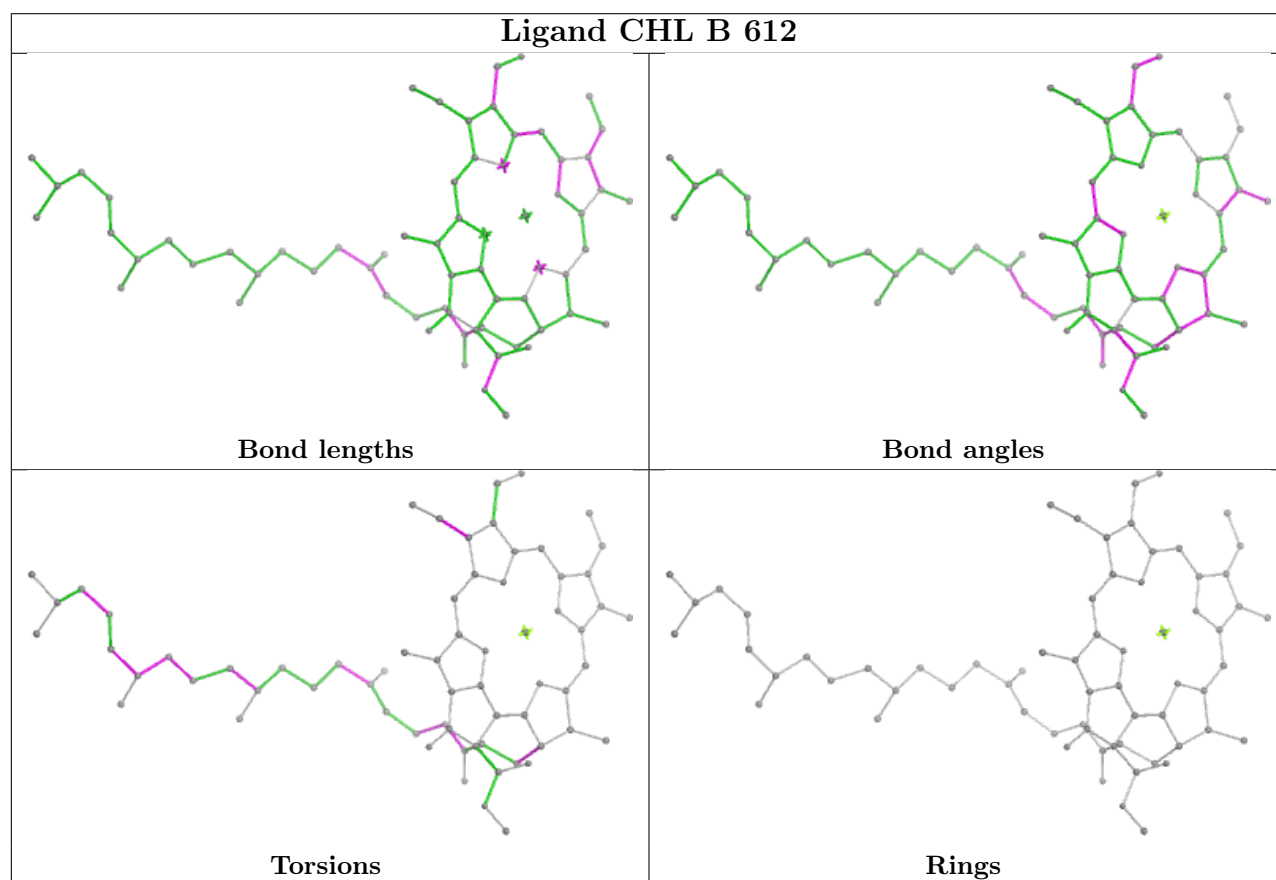
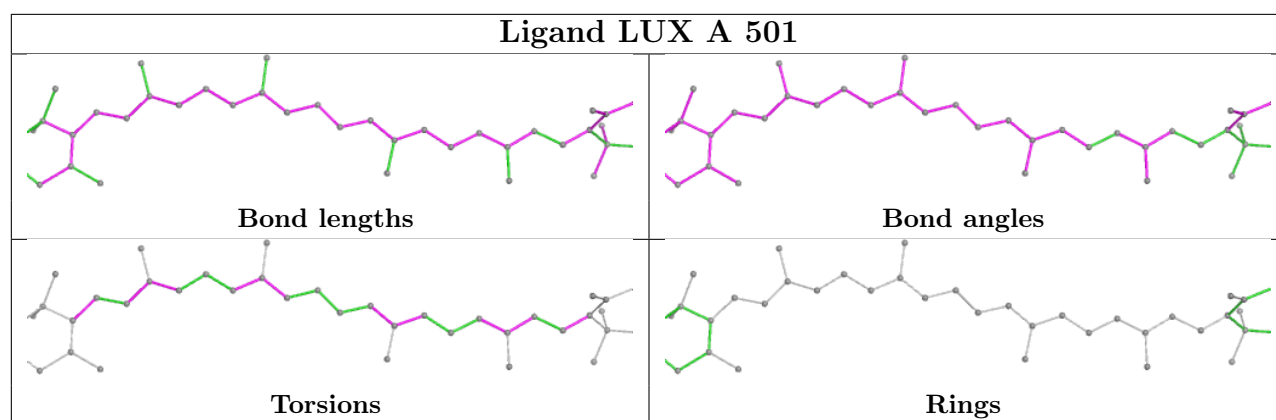


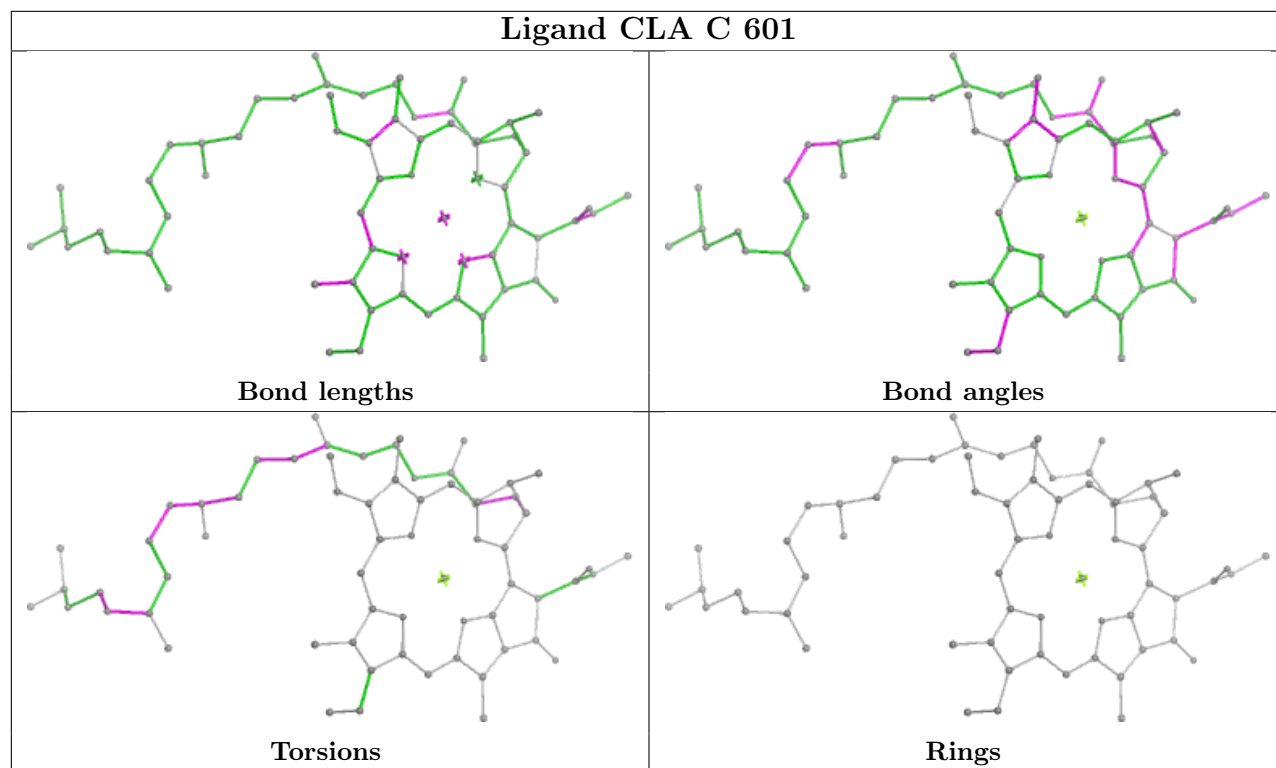


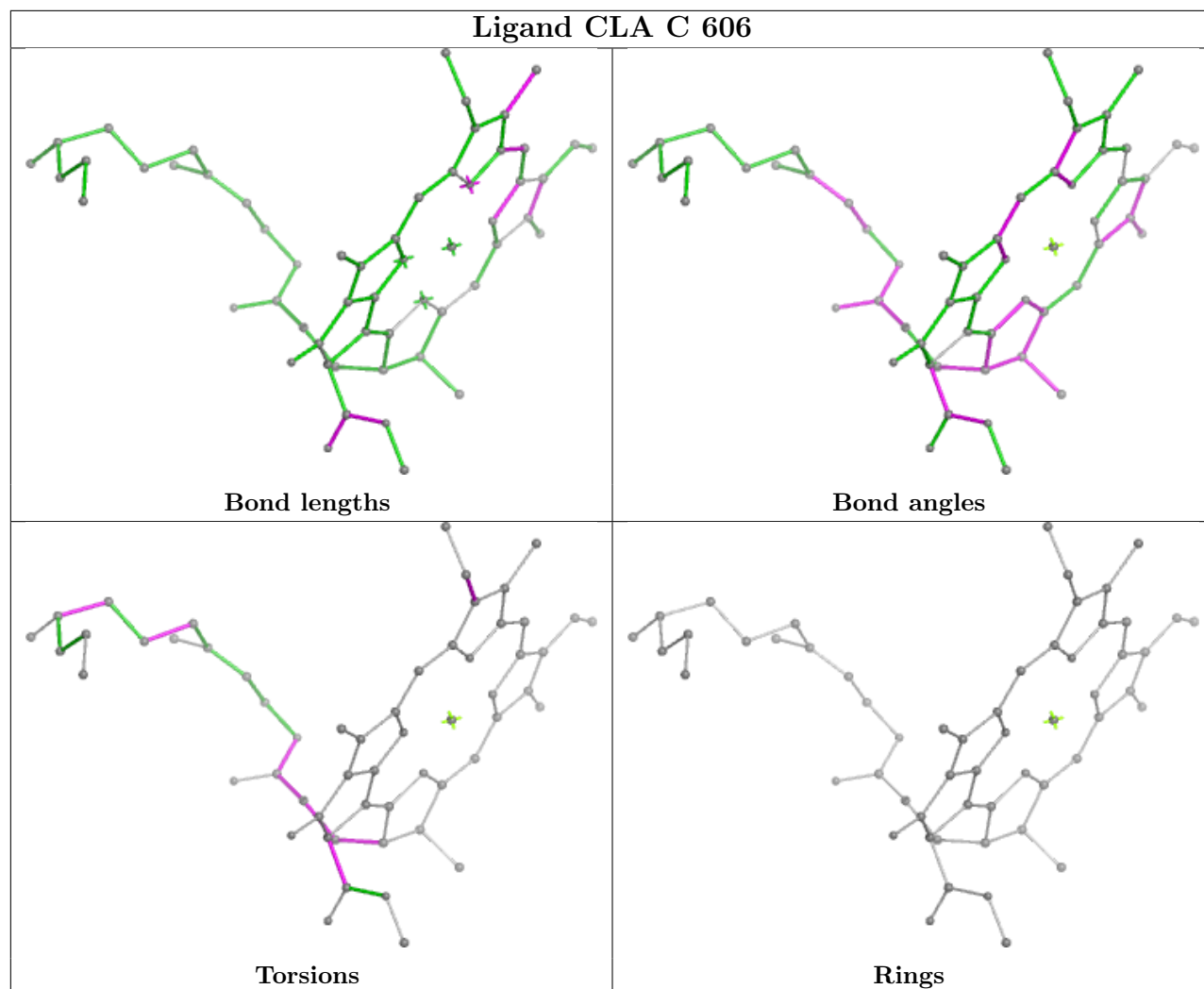


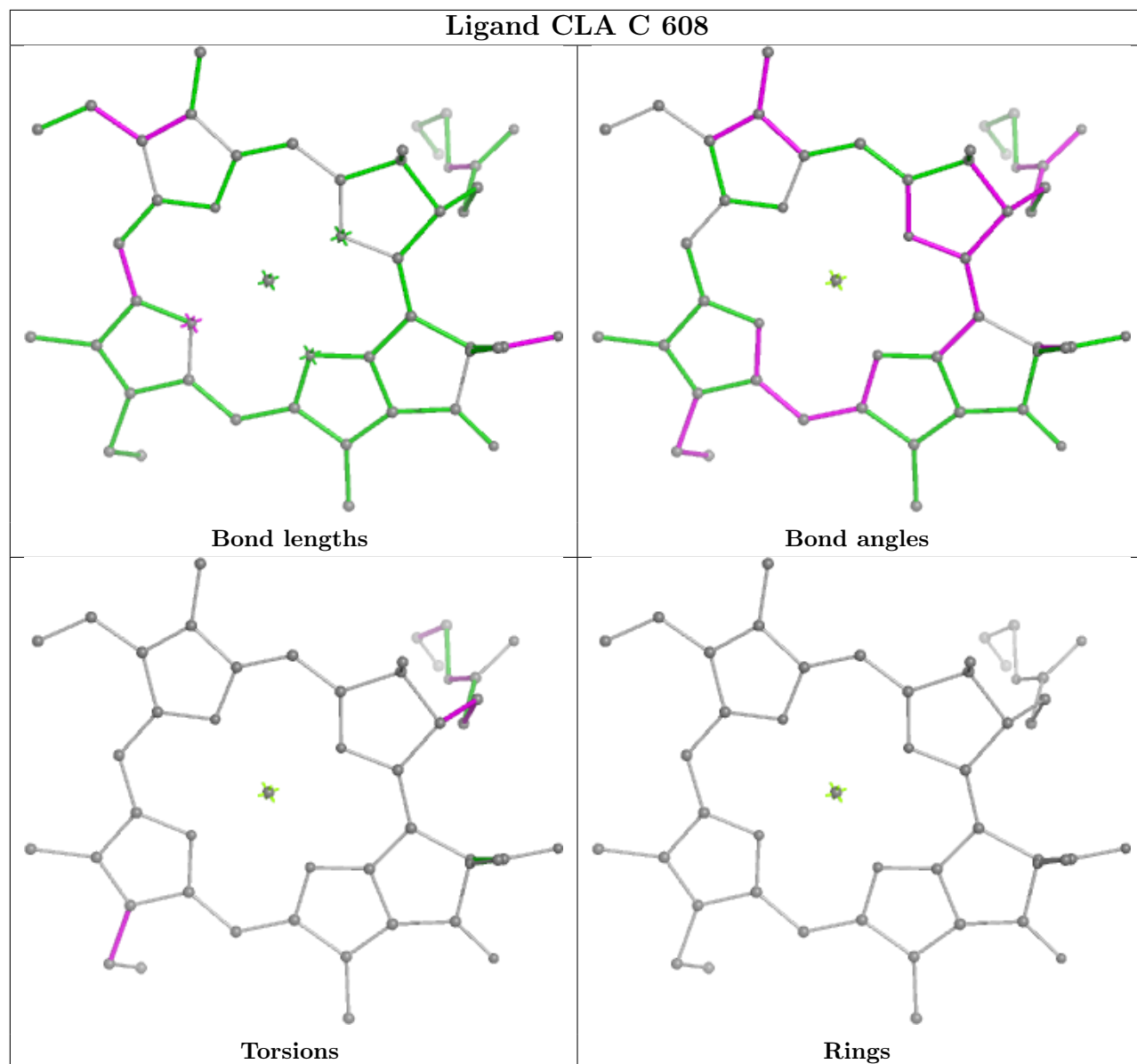


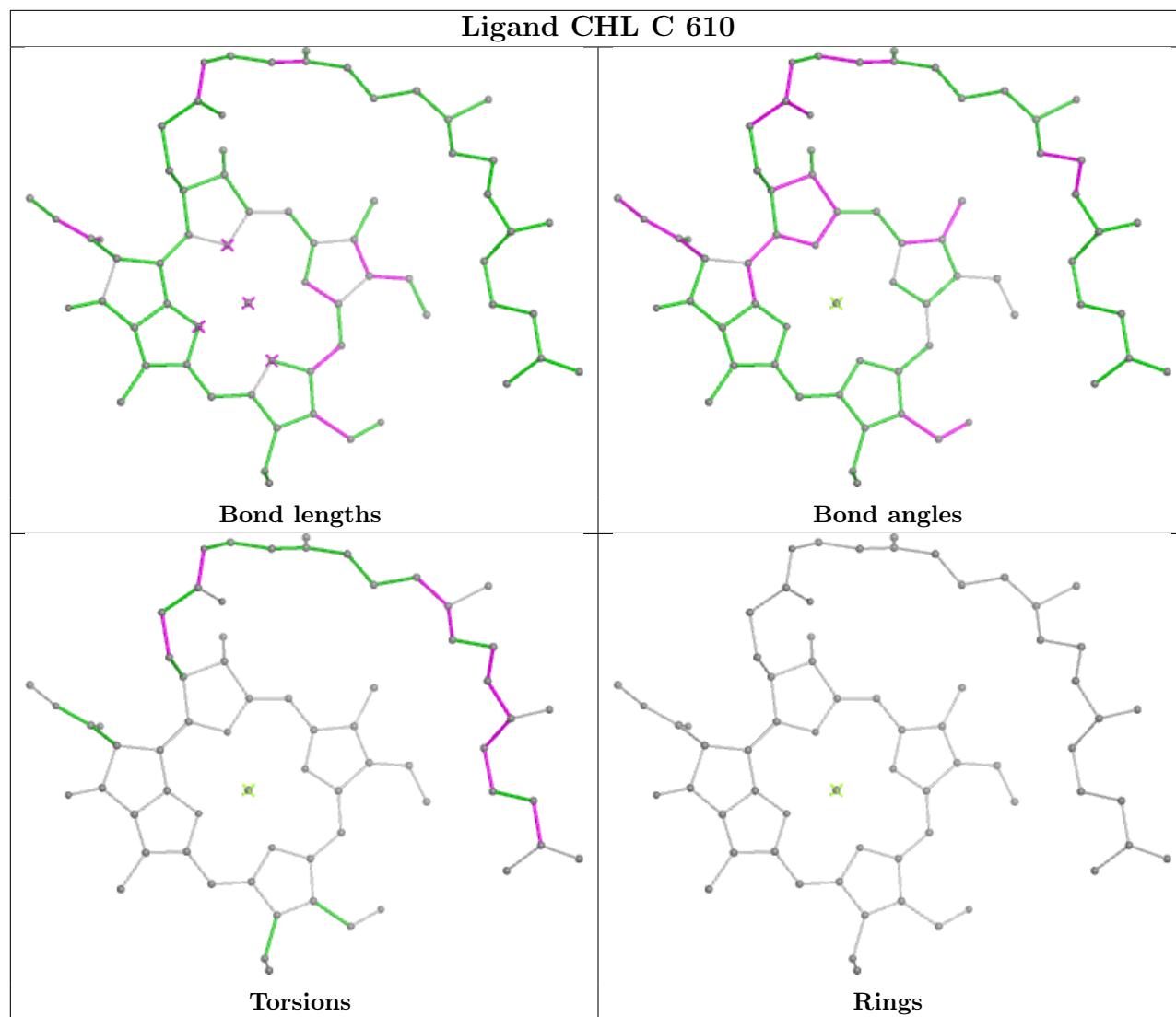


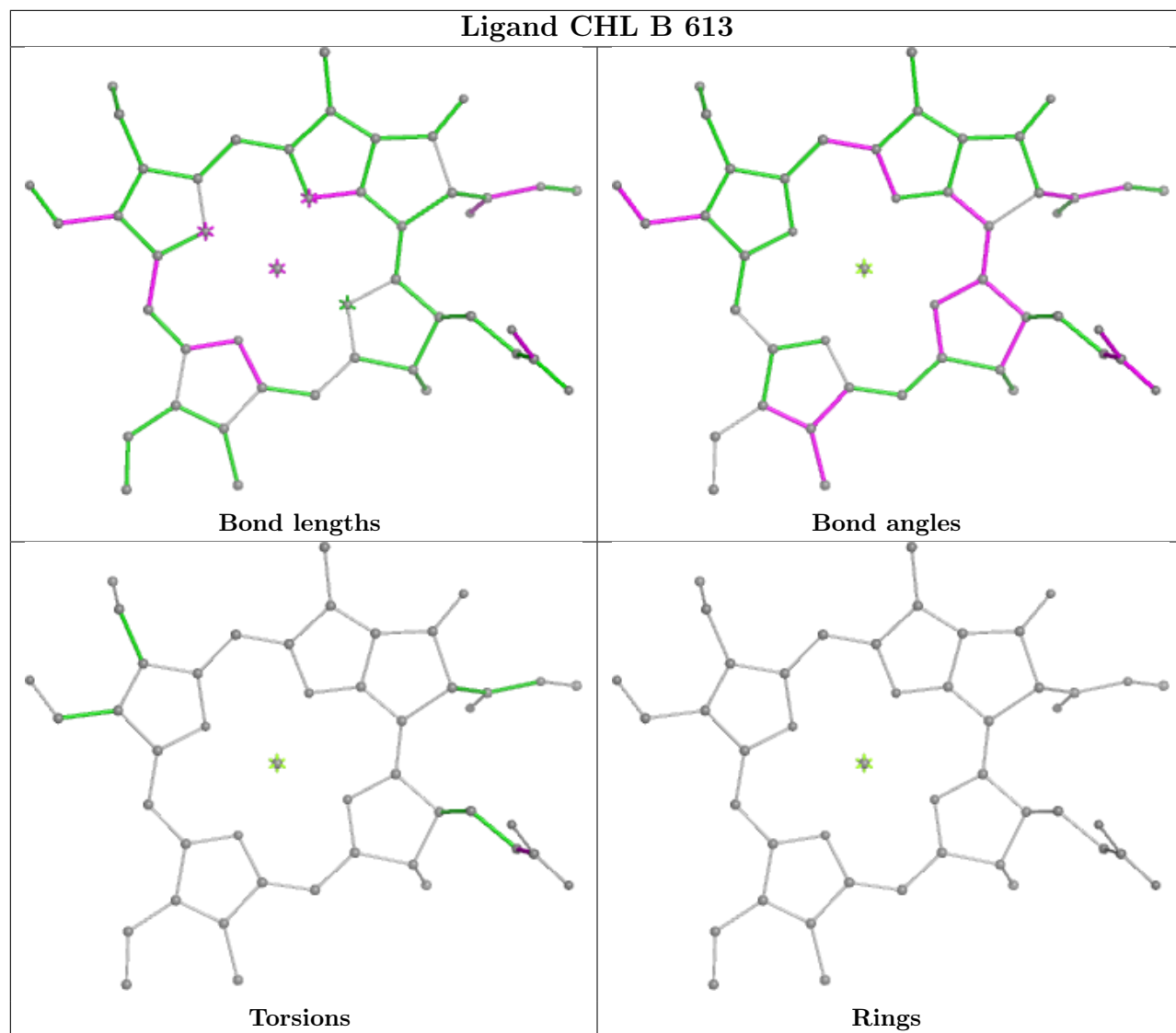


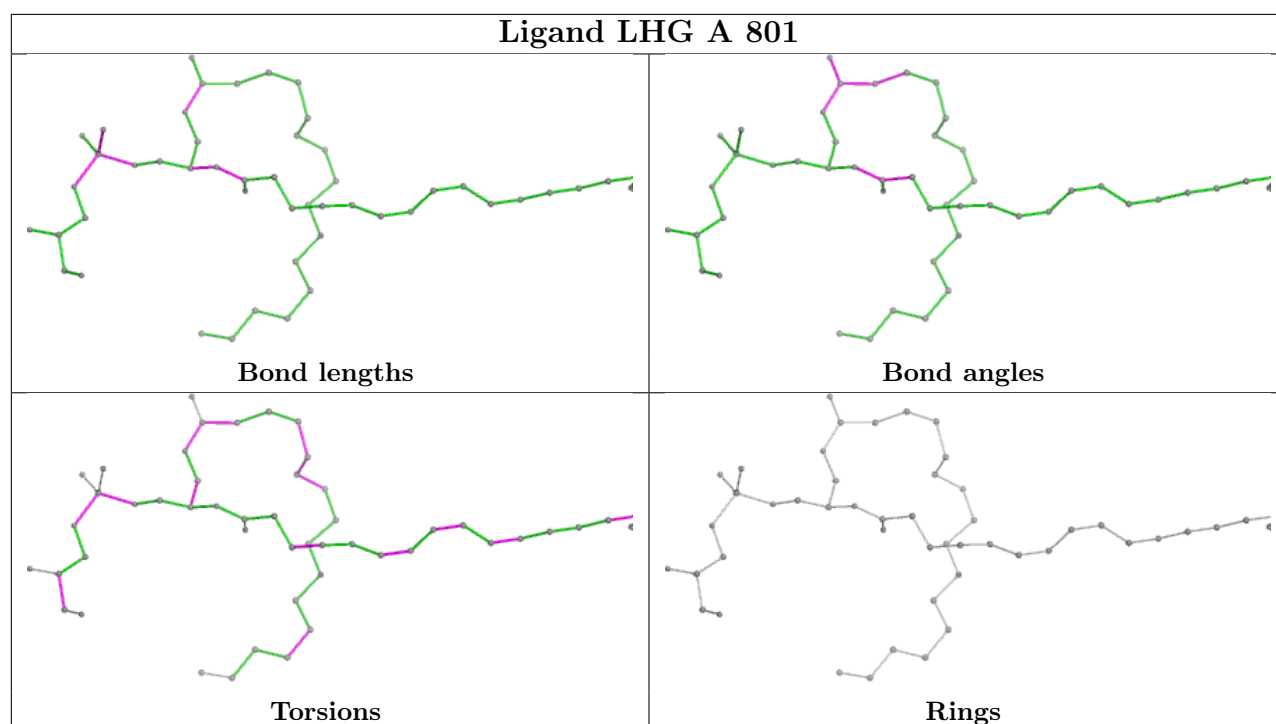
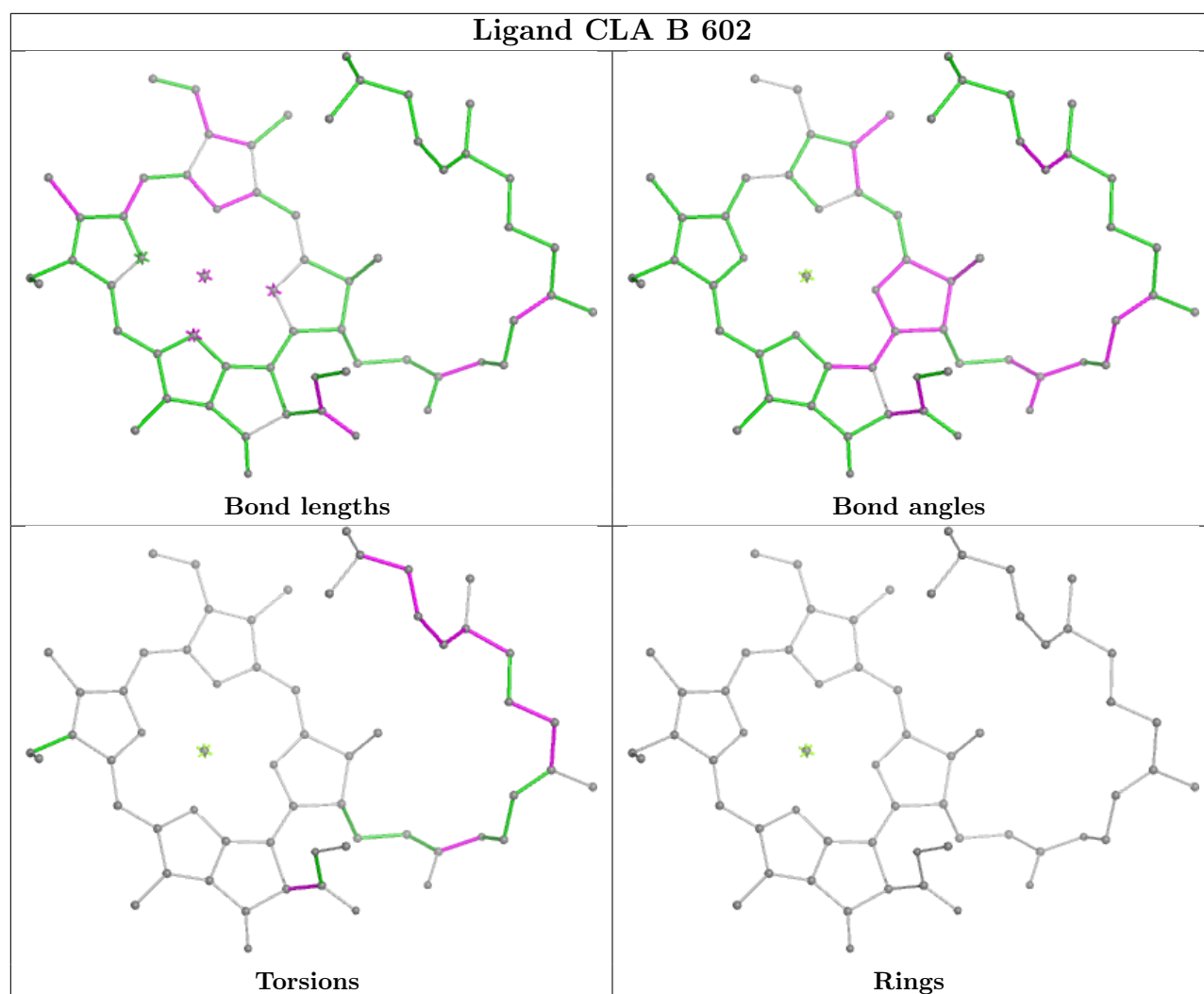


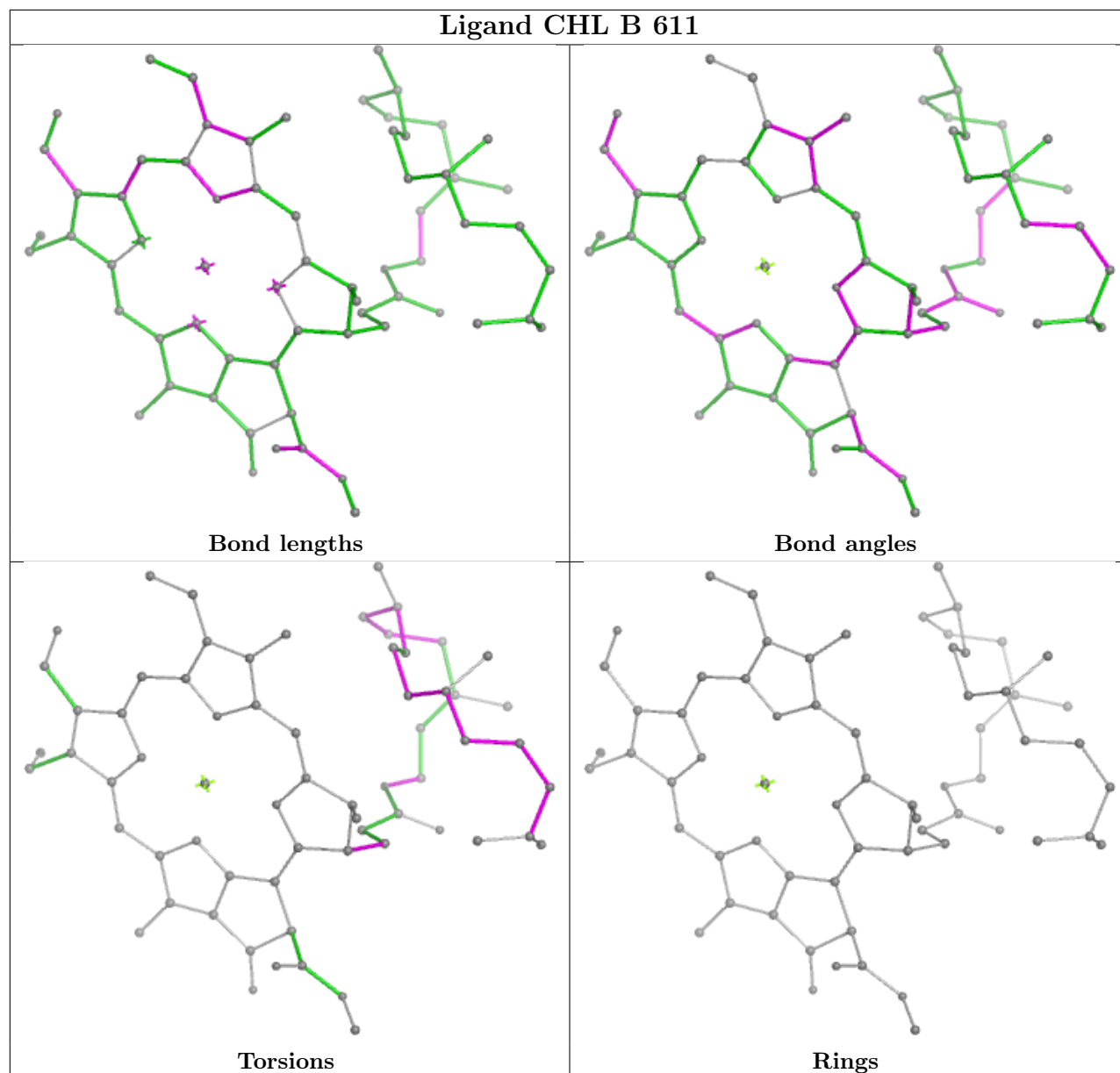


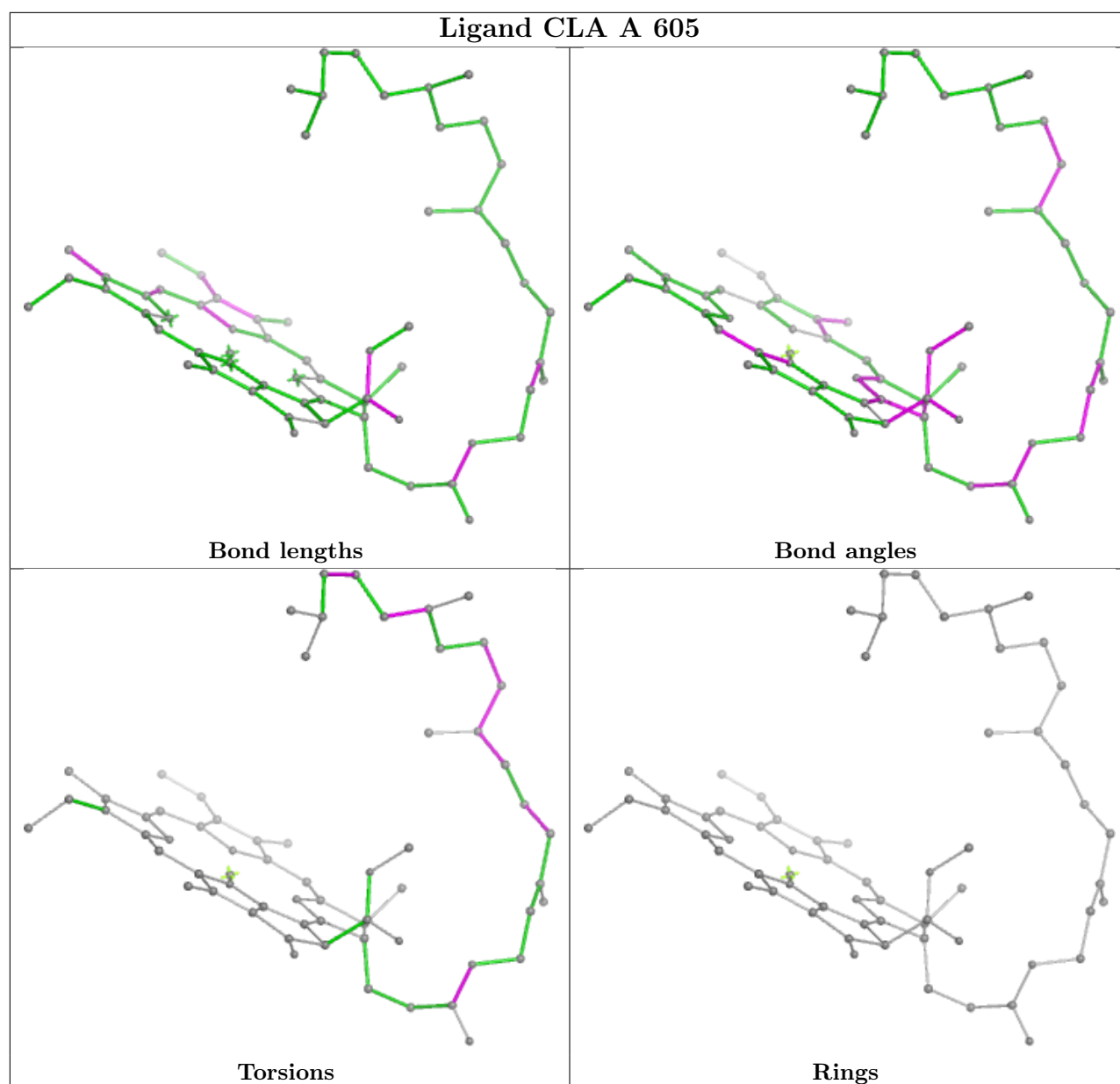


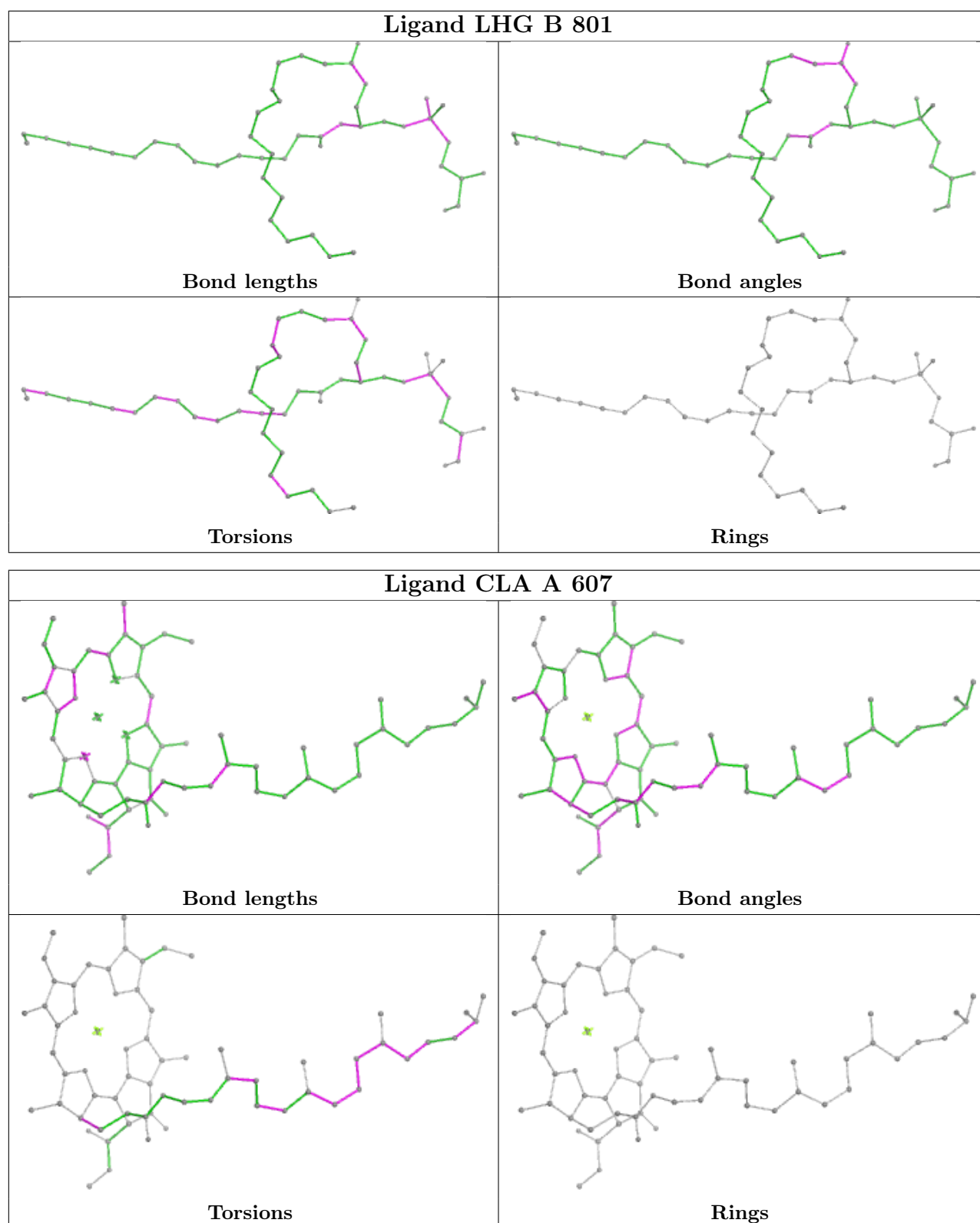












5.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

5.8 Polymer linkage issues

There are no chain breaks in this entry.

6 Fit of model and data [i](#)

6.1 Protein, DNA and RNA chains [i](#)

In the following table, the column labelled ‘#RSRZ > 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95th percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q < 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å ²)	Q<0.9
1	A	223/232 (96%)	1.16	49 (21%) 0 0	72, 91, 109, 128	0
1	B	223/232 (96%)	1.01	37 (16%) 1 1	70, 89, 107, 126	0
1	C	223/232 (96%)	1.05	37 (16%) 1 1	73, 91, 109, 129	0
All	All	669/696 (96%)	1.07	123 (18%) 1 1	70, 90, 109, 129	0

The worst 5 of 123 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	C	11	SER	9.7
1	B	12	SER	9.3
1	B	10	ALA	7.8
1	C	10	ALA	7.3
1	B	11	SER	6.7

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

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

6.3 Carbohydrates [i](#)

There are no monosaccharides in this entry.

6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled ‘Q < 0.9’ lists the number of atoms with occupancy less than 0.9.

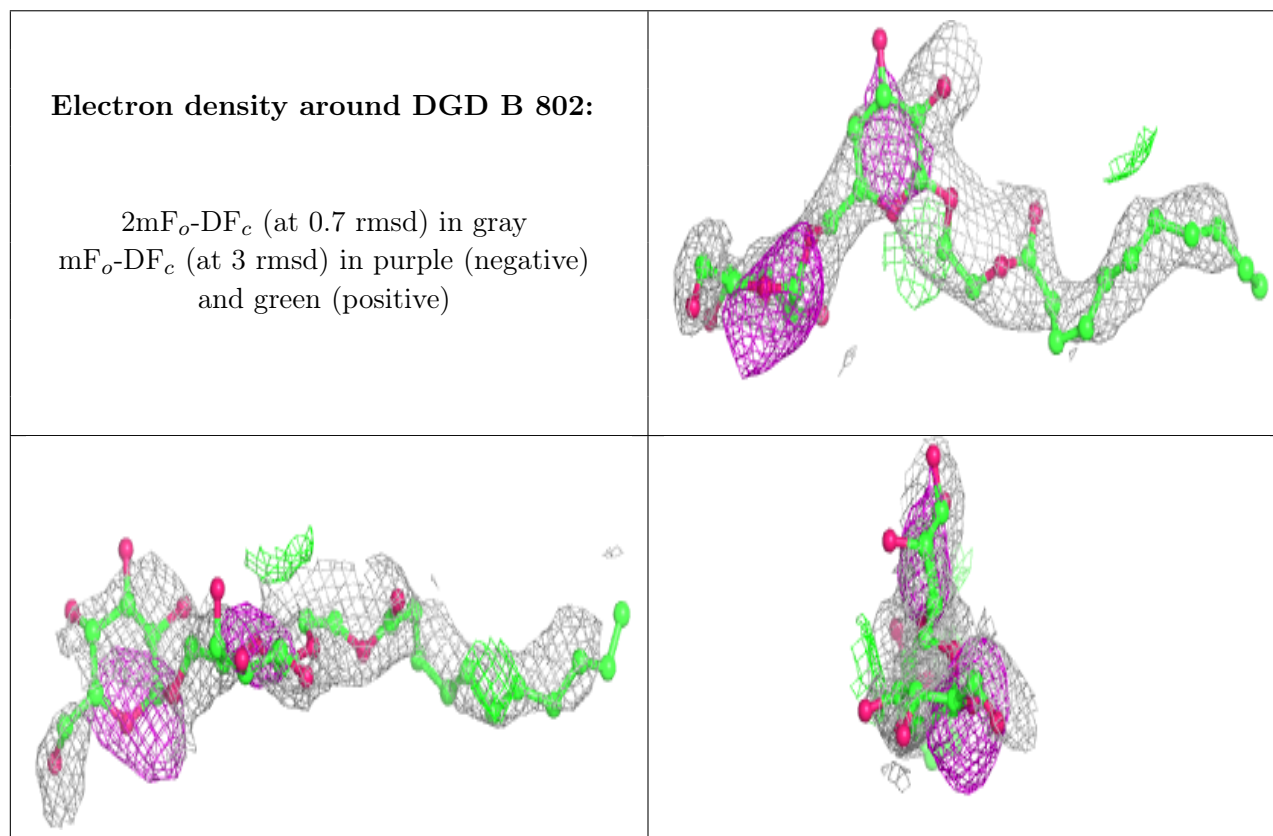
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
8	DGD	B	802	38/66	0.41	0.50	104,118,139,140	0
8	DGD	A	802	38/66	0.49	0.46	106,119,140,140	0
8	DGD	C	802	38/66	0.58	0.43	106,119,141,141	0
3	NEX	C	503	44/44	0.65	0.40	67,98,128,133	0
3	NEX	A	503	44/44	0.80	0.30	66,97,128,132	0
3	NEX	B	503	44/44	0.81	0.35	64,96,126,131	0
6	CHL	C	612	66/66	0.83	0.31	77,85,102,104	0
6	CHL	C	614	42/66	0.83	0.41	99,107,113,137	0
5	CLA	A	607	65/65	0.83	0.23	77,90,103,106	0
5	CLA	C	607	65/65	0.83	0.22	77,91,104,106	0
6	CHL	A	612	66/66	0.83	0.36	76,85,102,104	0
5	CLA	B	607	65/65	0.84	0.24	75,88,101,104	0
2	LUX	B	502	42/42	0.84	0.35	53,76,113,126	0
5	CLA	A	604	65/65	0.85	0.27	59,71,96,101	0
5	CLA	C	604	65/65	0.85	0.30	60,72,97,102	0
7	LHG	B	801	49/49	0.85	0.28	72,88,109,115	0
5	CLA	B	601	65/65	0.86	0.23	68,77,111,113	0
6	CHL	B	614	42/66	0.86	0.42	97,105,111,134	0
4	XAT	C	504	44/44	0.87	0.20	65,91,117,149	0
6	CHL	C	609	66/66	0.87	0.24	77,85,98,101	0
6	CHL	A	609	66/66	0.87	0.22	75,85,97,100	0
5	CLA	B	604	65/65	0.87	0.30	56,69,95,100	0
4	XAT	B	504	44/44	0.88	0.21	63,89,115,147	0
2	LUX	C	502	42/42	0.88	0.35	58,79,115,128	0
5	CLA	A	601	65/65	0.88	0.23	69,79,112,115	0
2	LUX	C	501	42/42	0.88	0.26	62,88,120,136	0
6	CHL	A	614	42/66	0.88	0.43	99,106,113,136	0
6	CHL	B	609	66/66	0.88	0.22	75,83,95,99	0
5	CLA	C	601	65/65	0.88	0.24	70,79,113,115	0
5	CLA	B	606	57/65	0.89	0.28	78,89,105,109	0
7	LHG	A	801	49/49	0.89	0.25	74,90,110,116	0
6	CHL	A	611	66/66	0.89	0.24	70,85,117,123	0
2	LUX	A	502	42/42	0.89	0.32	57,78,115,128	0
6	CHL	C	611	66/66	0.89	0.28	69,85,118,124	0
5	CLA	B	608	48/65	0.89	0.22	88,97,115,117	0
5	CLA	C	608	48/65	0.90	0.21	90,99,117,119	0
5	CLA	B	603	65/65	0.90	0.20	74,83,100,103	0
2	LUX	B	501	42/42	0.90	0.26	59,86,118,134	0
7	LHG	C	801	49/49	0.90	0.24	74,90,111,117	0
6	CHL	C	610	66/66	0.90	0.24	67,81,108,110	0
5	CLA	A	608	48/65	0.90	0.16	90,99,116,118	0
5	CLA	A	606	57/65	0.90	0.26	79,90,107,110	0
4	XAT	A	504	44/44	0.91	0.18	65,91,117,148	0

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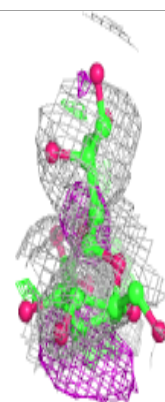
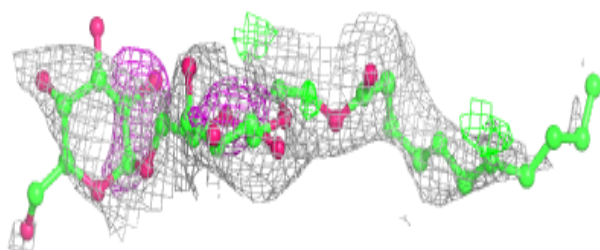
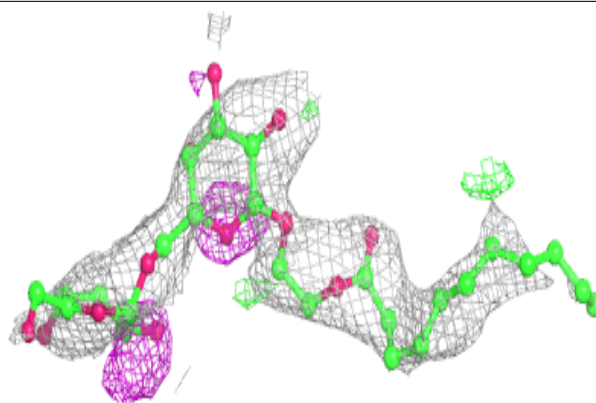
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(\AA^2)	Q<0.9
5	CLA	C	606	57/65	0.91	0.25	80,91,107,111	0
6	CHL	C	613	46/66	0.91	0.17	73,82,107,121	0
5	CLA	A	605	65/65	0.91	0.49	77,86,107,110	0
6	CHL	B	611	66/66	0.91	0.26	68,82,115,121	0
6	CHL	B	612	66/66	0.91	0.27	74,83,100,102	0
6	CHL	B	613	46/66	0.91	0.20	71,80,106,119	0
5	CLA	A	602	60/65	0.91	0.24	79,87,115,116	0
5	CLA	A	603	65/65	0.91	0.21	75,85,102,105	0
5	CLA	C	602	60/65	0.91	0.21	80,87,116,117	0
5	CLA	C	603	65/65	0.92	0.22	75,85,102,105	0
6	CHL	A	613	46/66	0.92	0.20	73,81,107,120	0
5	CLA	B	602	60/65	0.92	0.19	77,85,114,114	0
5	CLA	C	605	65/65	0.92	0.41	77,86,108,110	0
6	CHL	B	610	66/66	0.92	0.23	65,79,106,108	0
6	CHL	A	610	66/66	0.92	0.29	65,80,108,110	0
2	LUX	A	501	42/42	0.92	0.27	61,87,120,136	0
5	CLA	B	605	65/65	0.93	0.45	75,84,105,108	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.

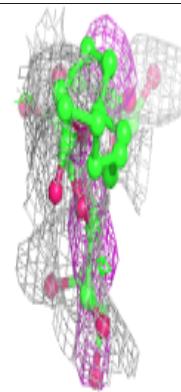
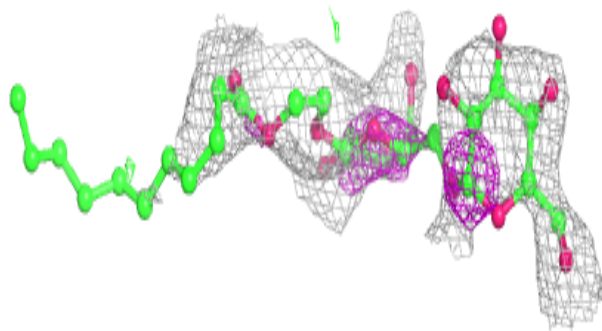
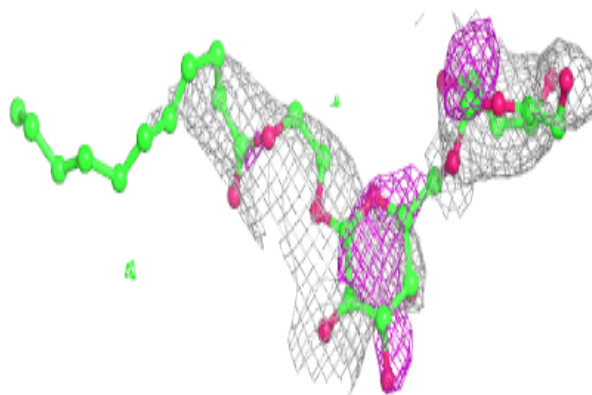


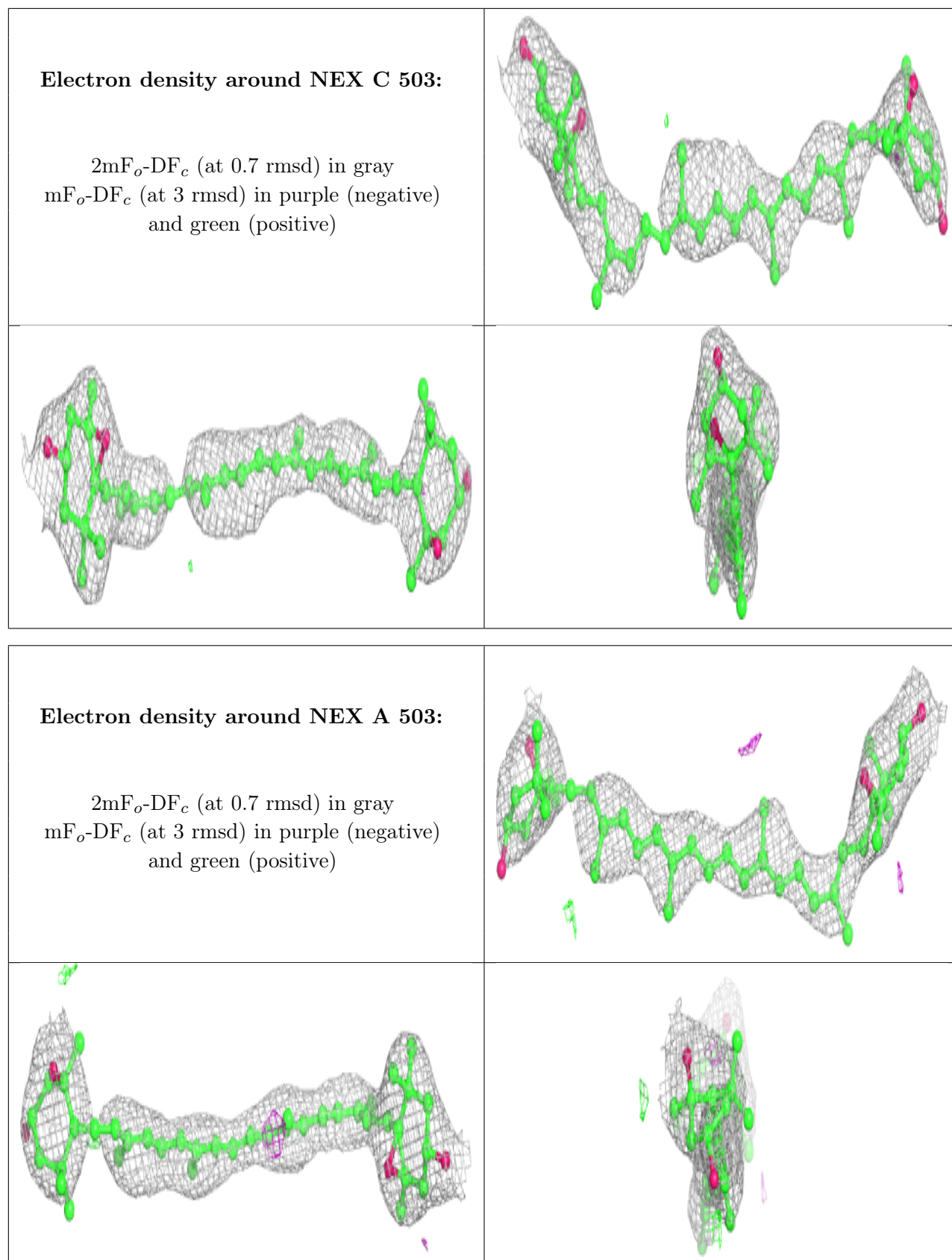
Electron density around DGD A 802:

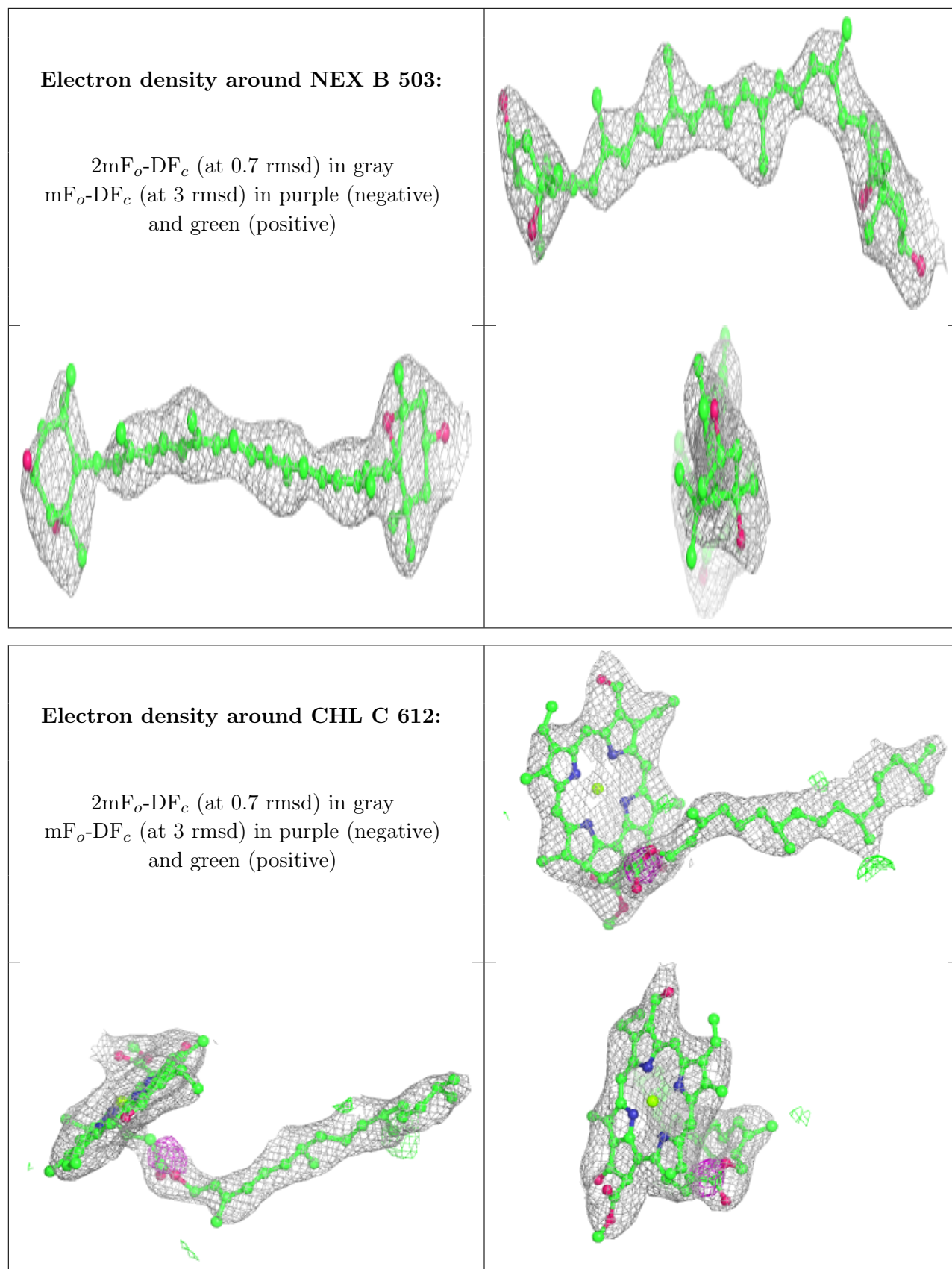
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

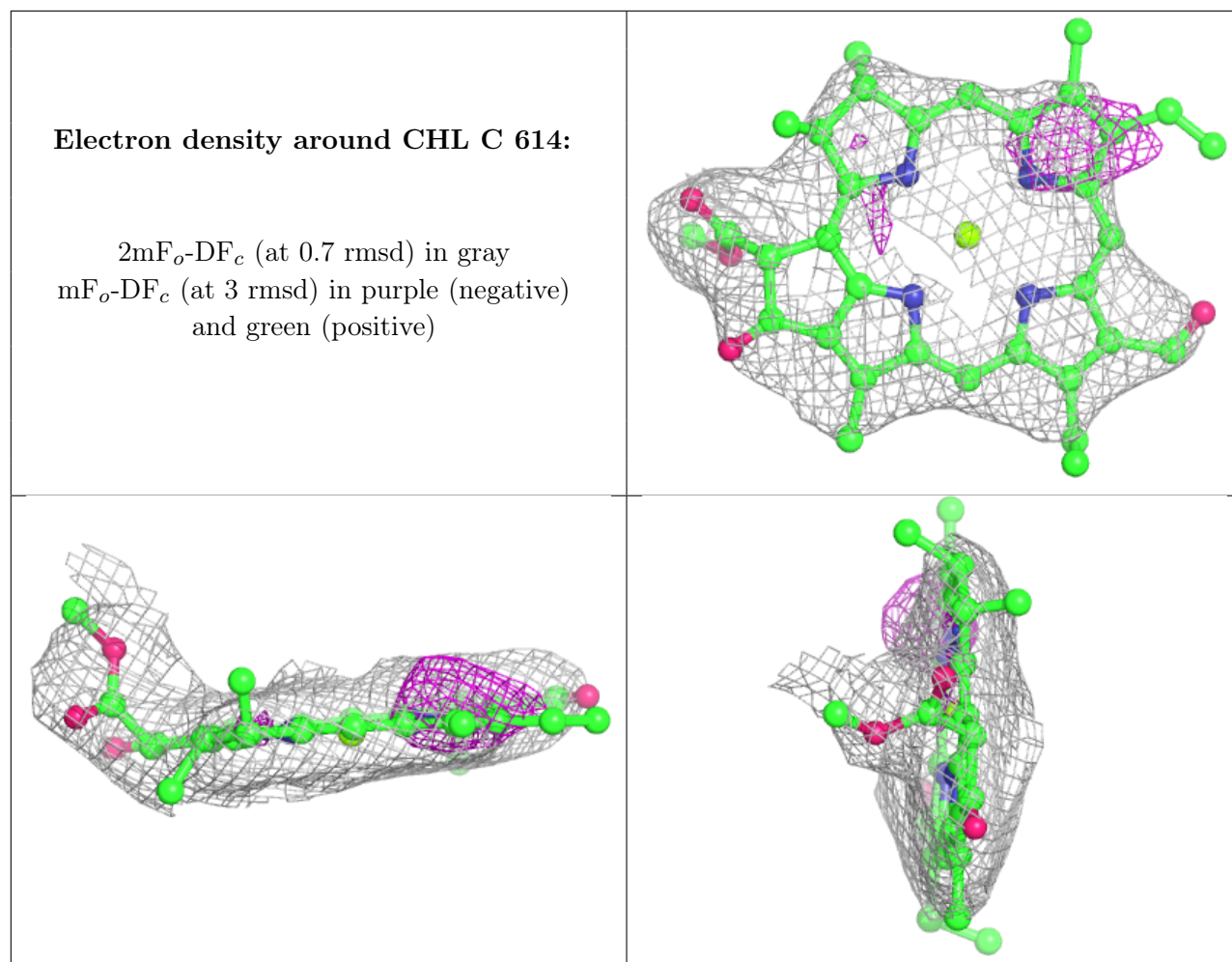
**Electron density around DGD C 802:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



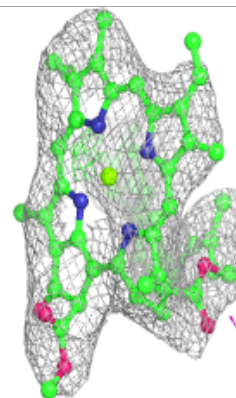
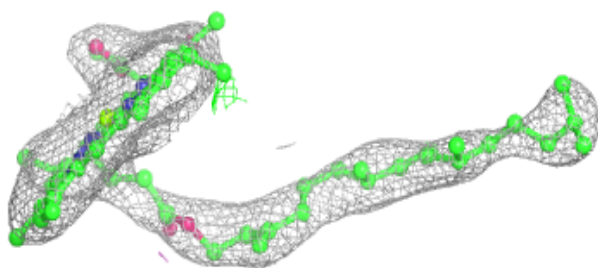
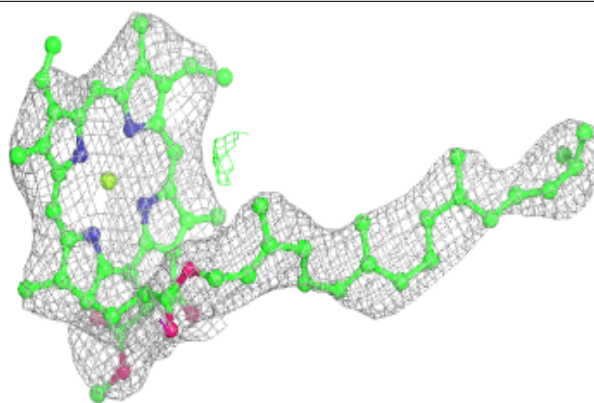




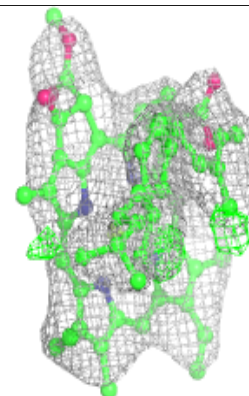
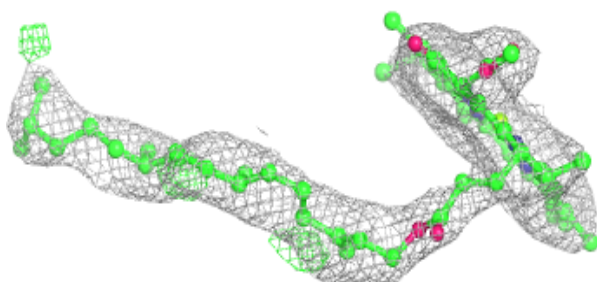
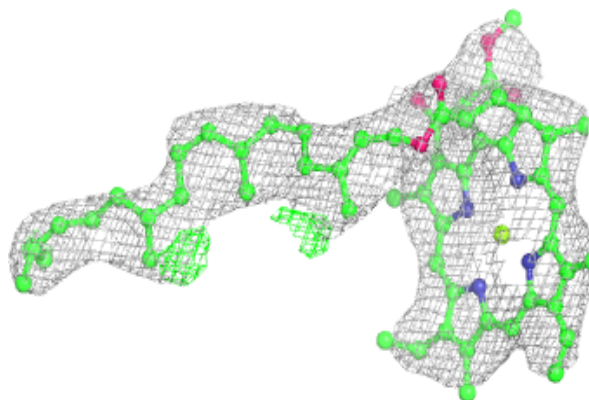


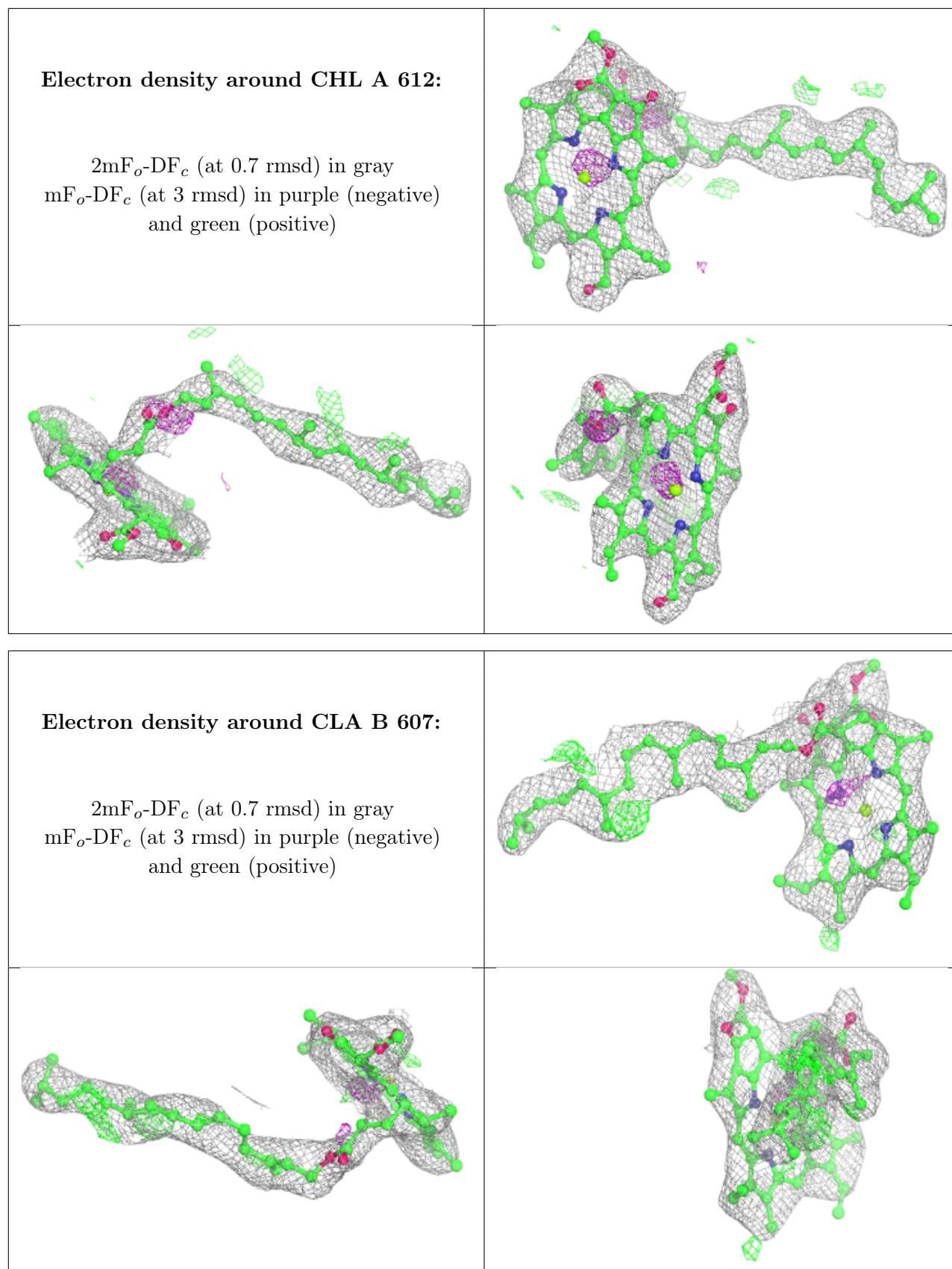
Electron density around CLA A 607:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around CLA C 607:**

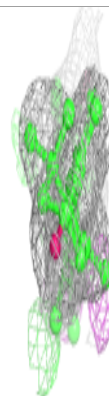
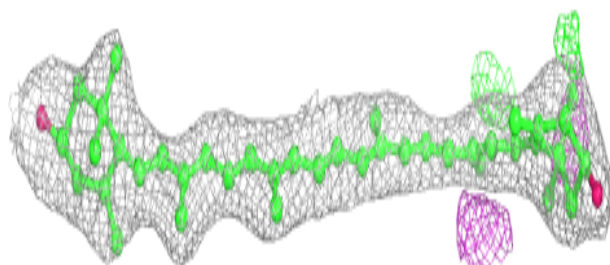
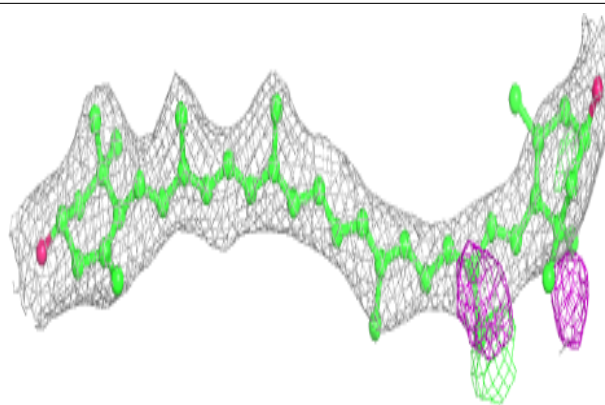
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



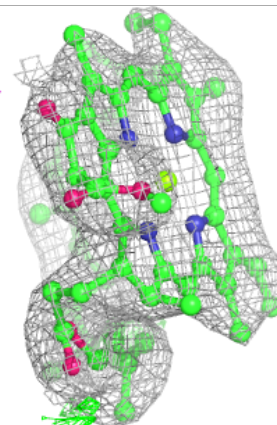
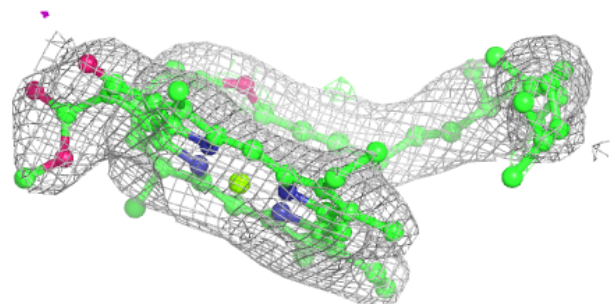
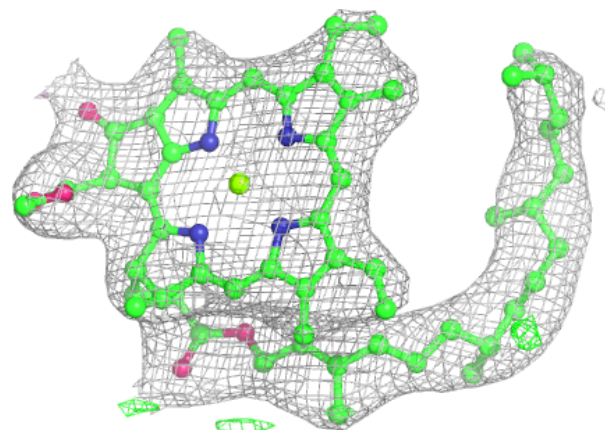


Electron density around LUX B 502:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

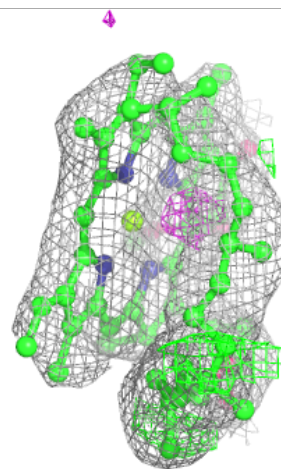
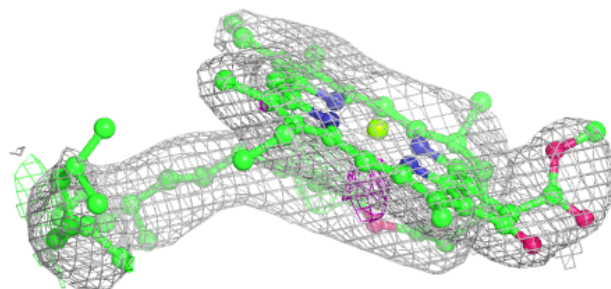
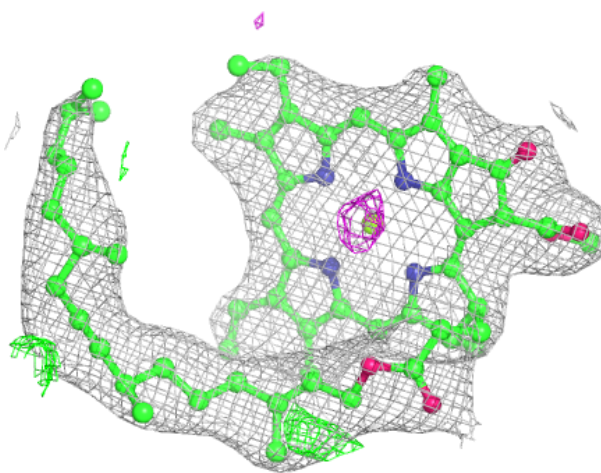
**Electron density around CLA A 604:**

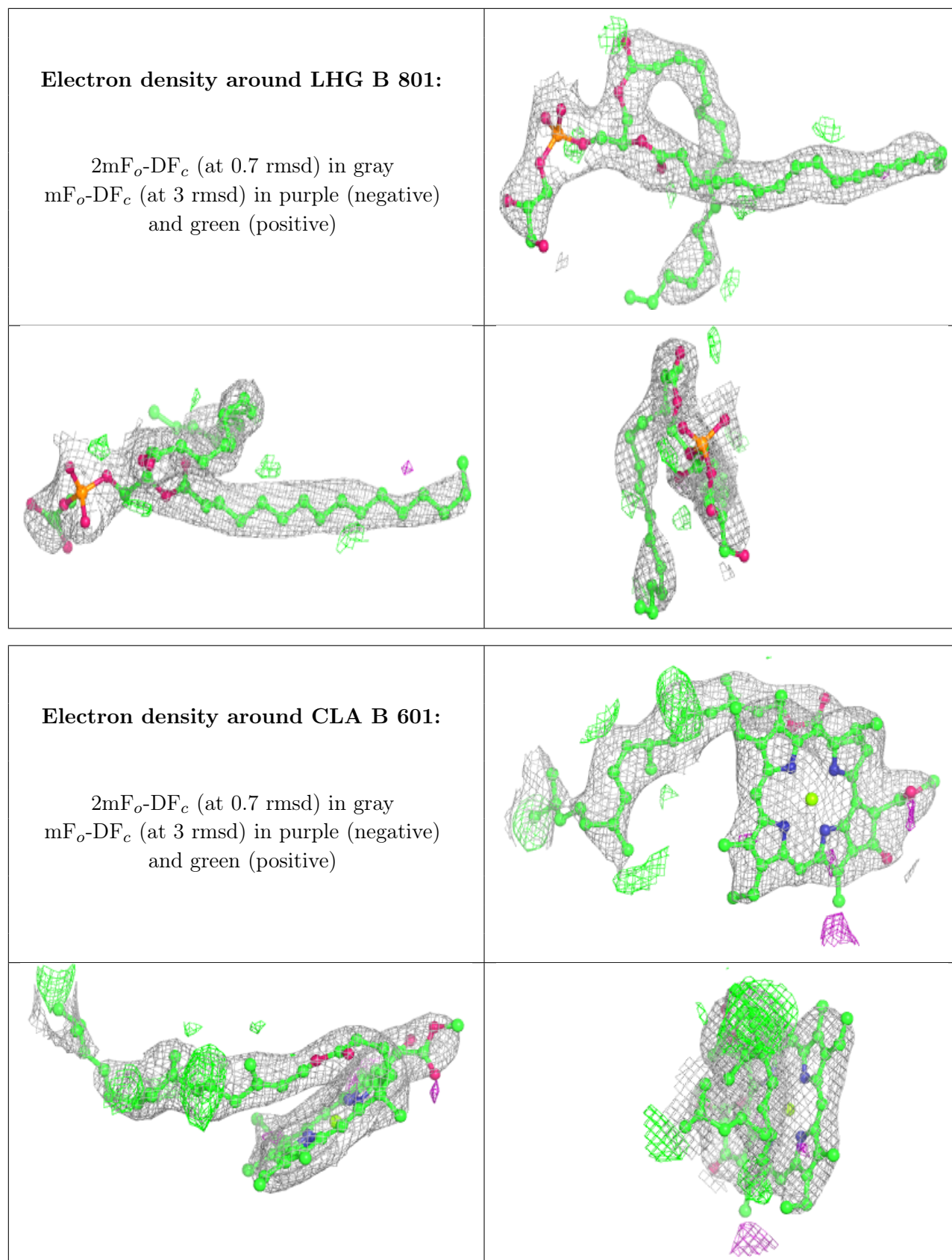
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around CLA C 604:

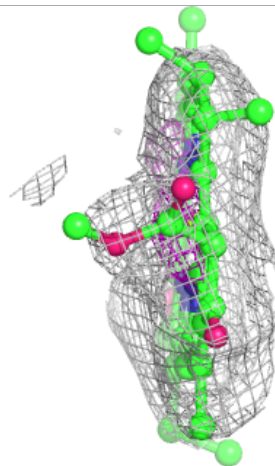
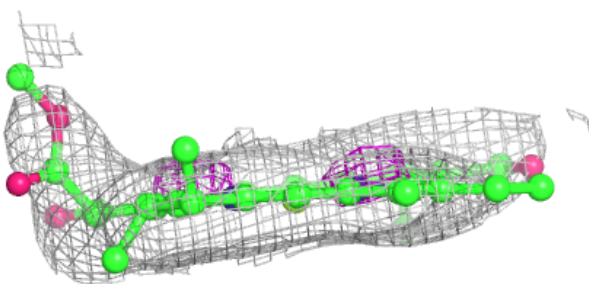
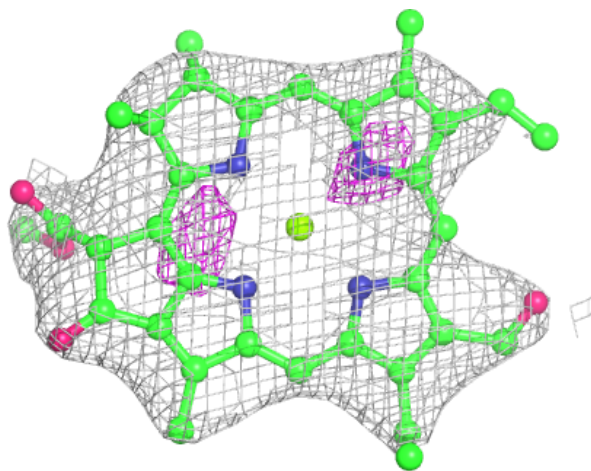
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

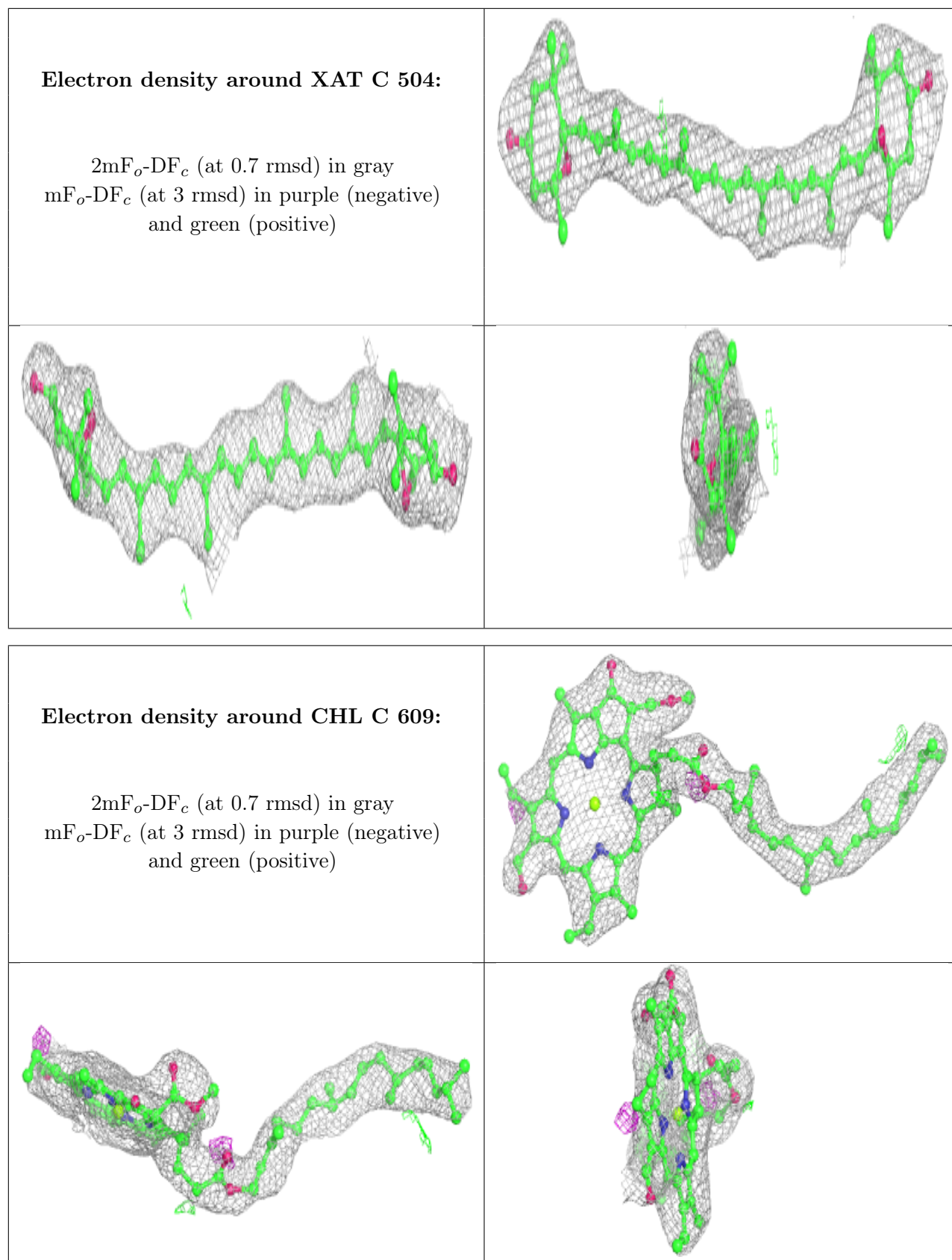




Electron density around CHL B 614:

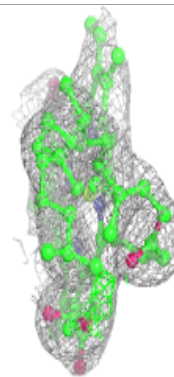
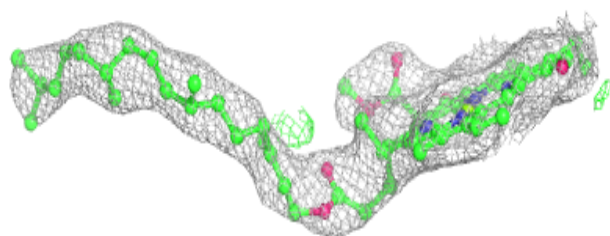
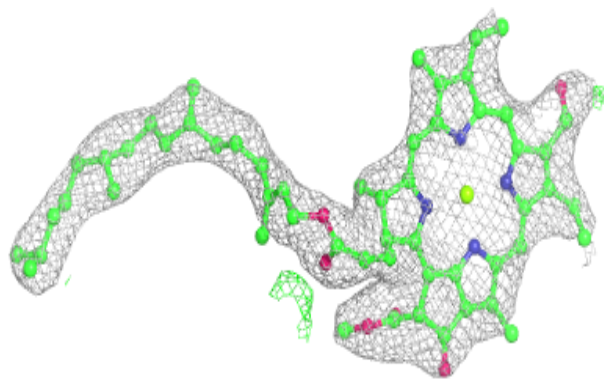
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



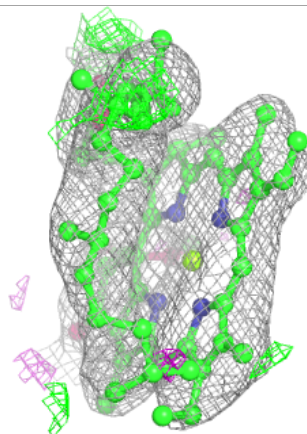
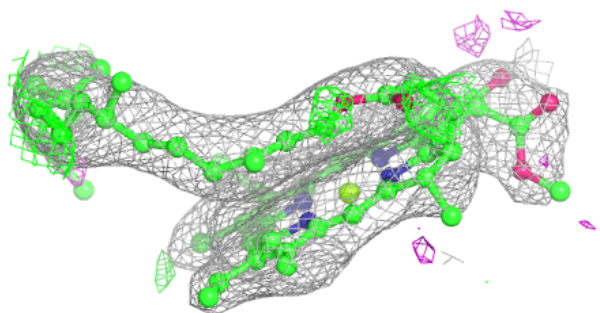
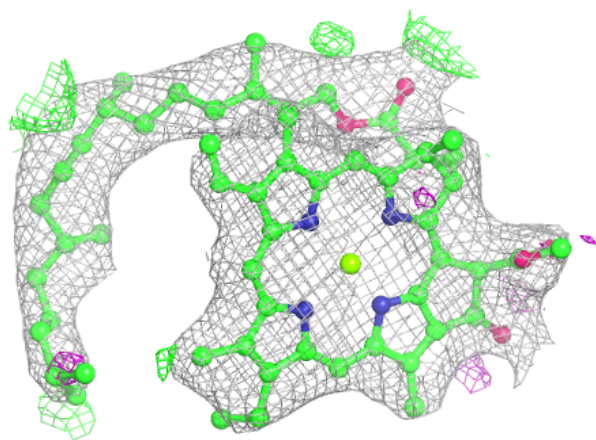


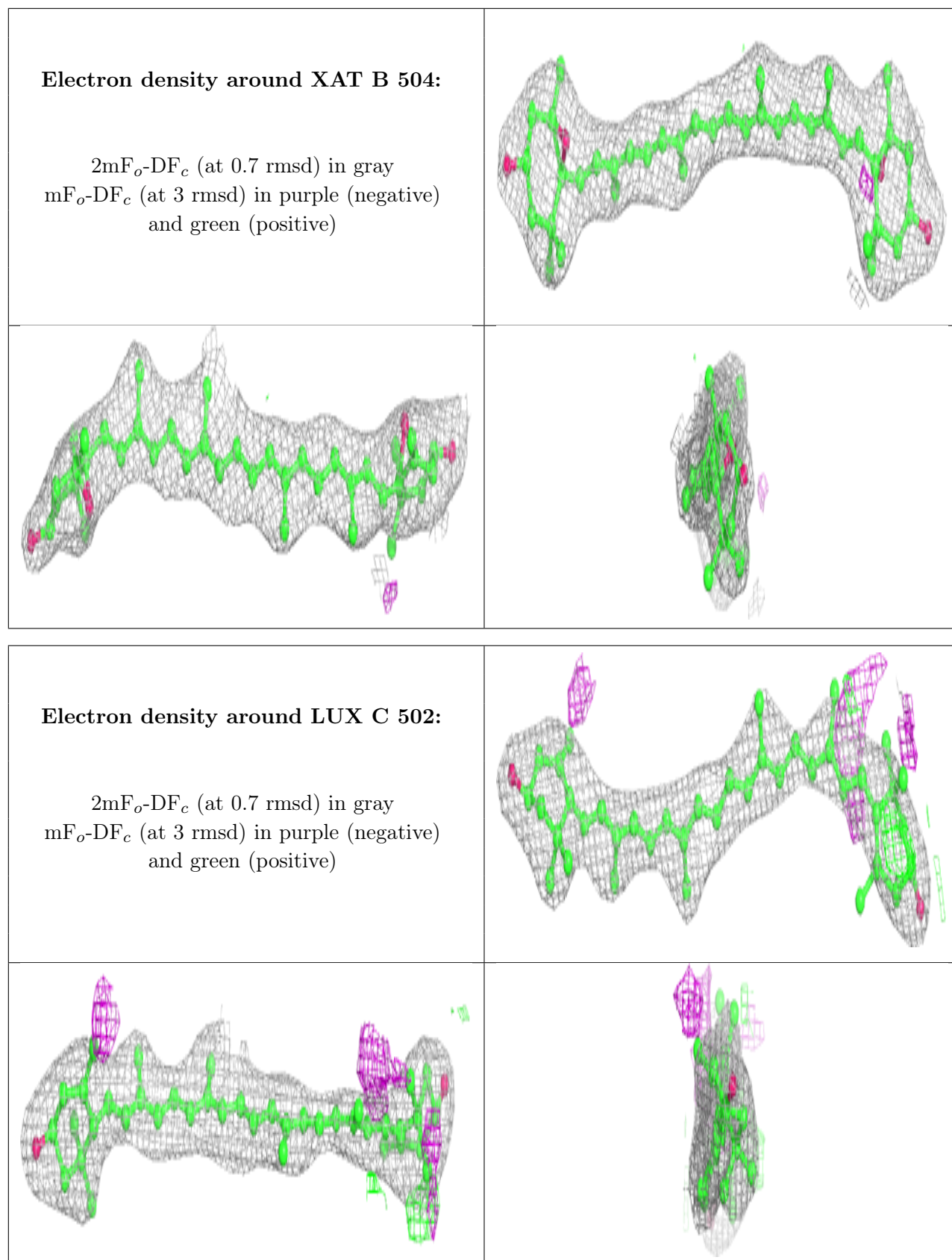
Electron density around CHL A 609:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around CLA B 604:**

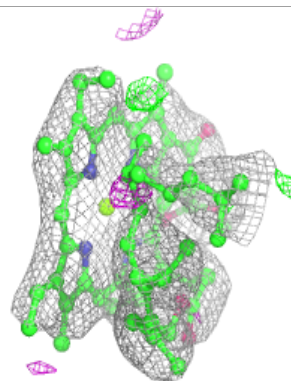
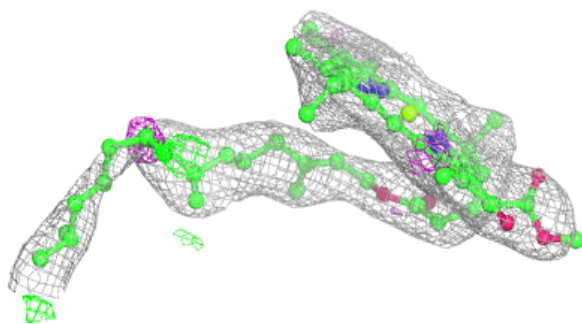
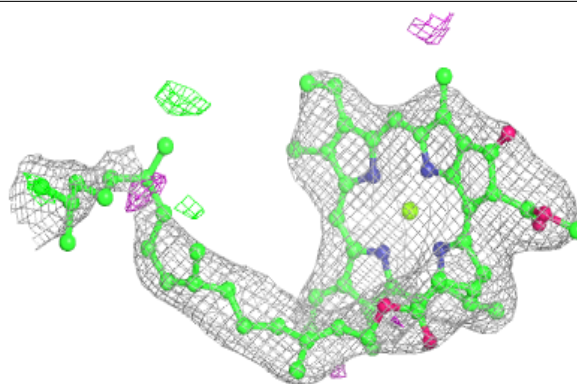
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



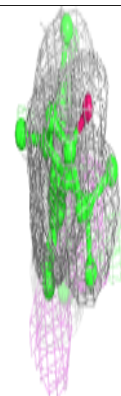
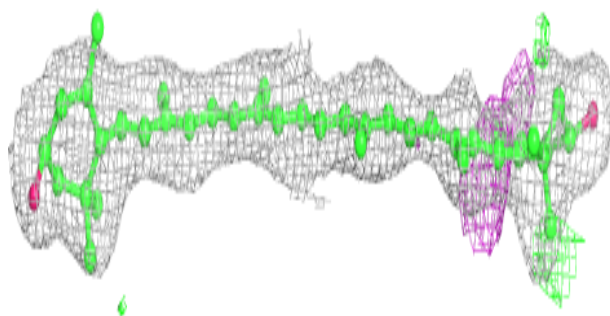
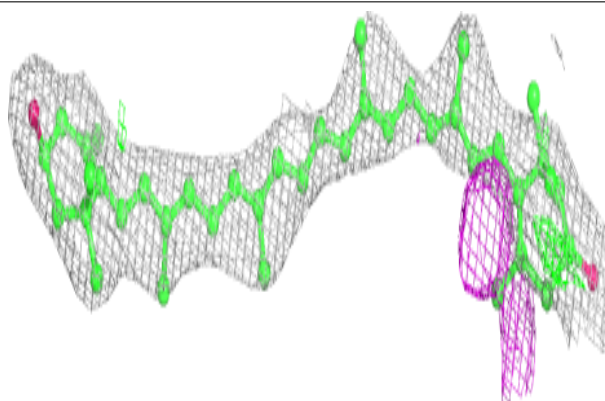


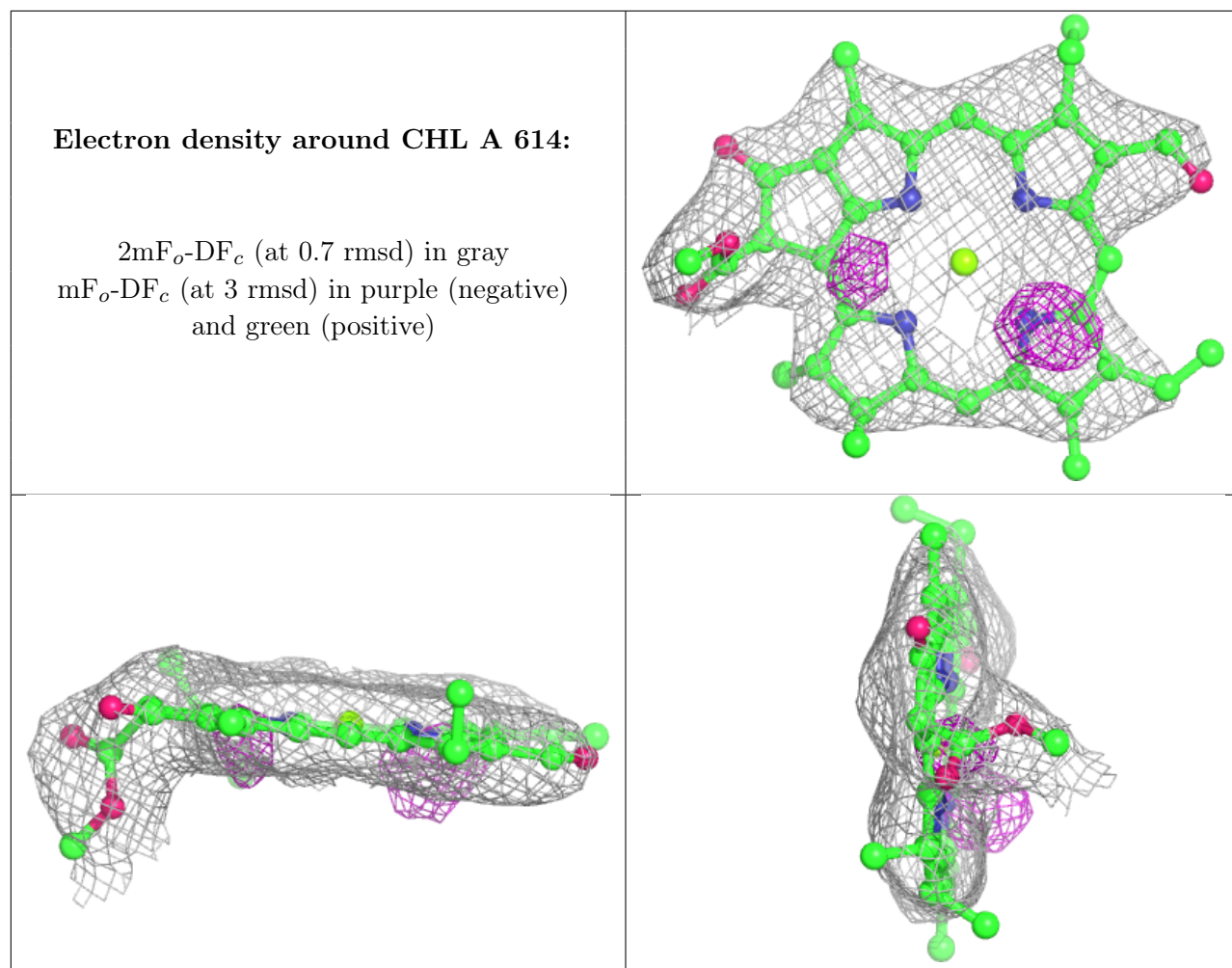
Electron density around CLA A 601:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

**Electron density around LUX C 501:**

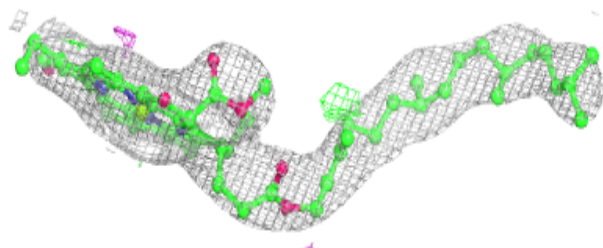
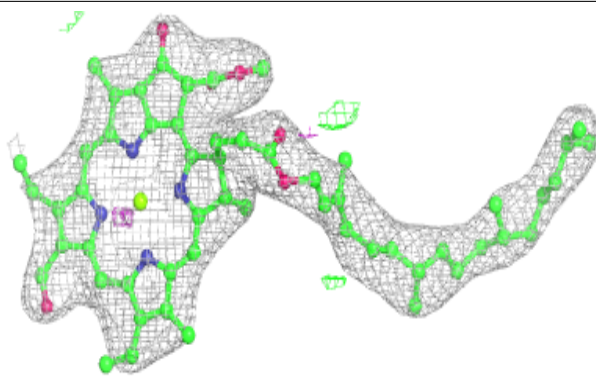
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



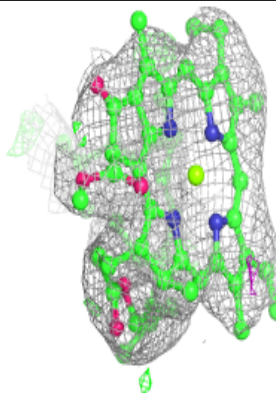
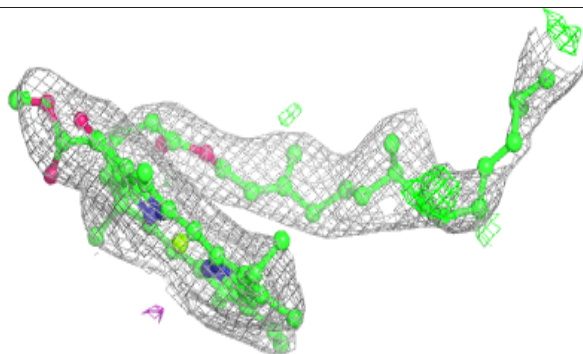
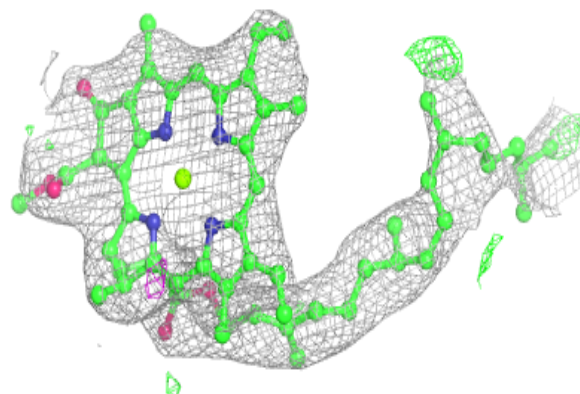


Electron density around CHL B 609:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

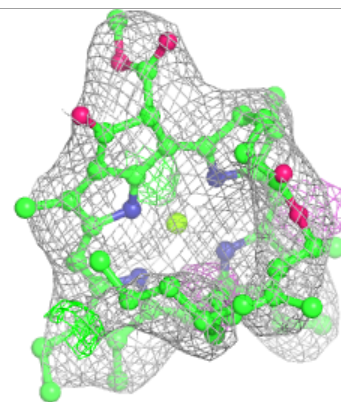
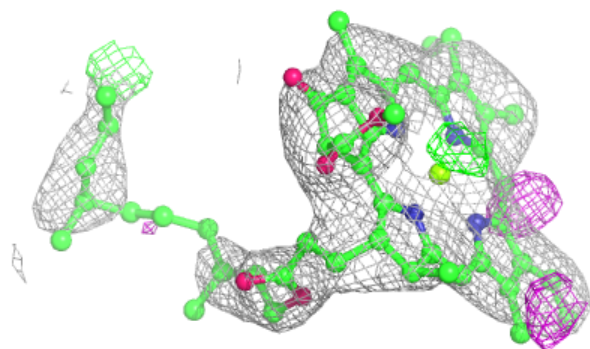
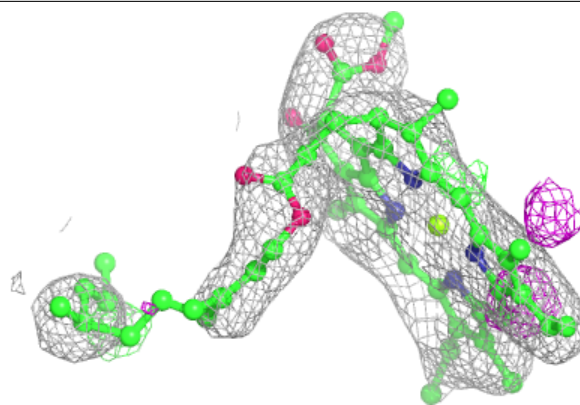
**Electron density around CLA C 601:**

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

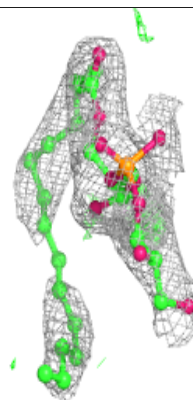
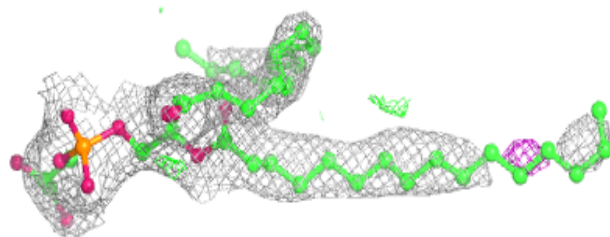
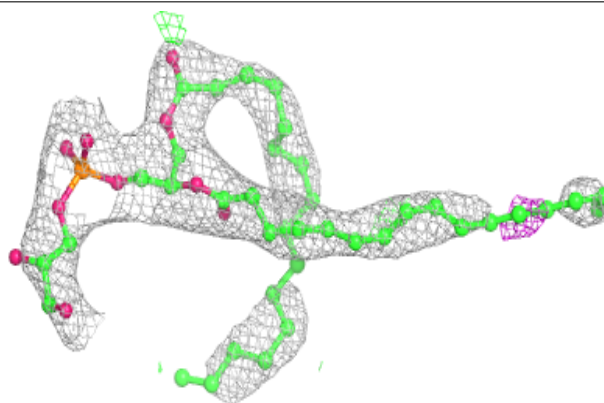


Electron density around CLA B 606:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

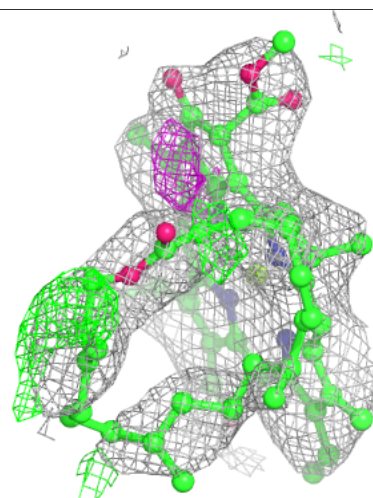
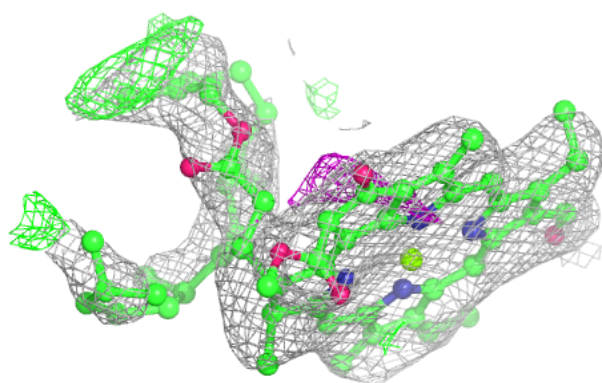
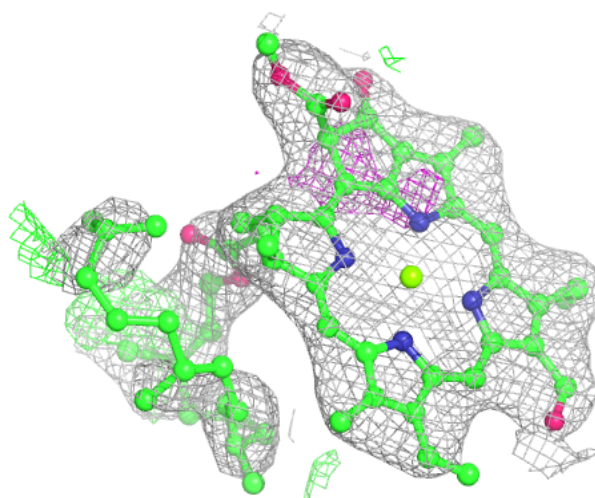
**Electron density around LHG A 801:**

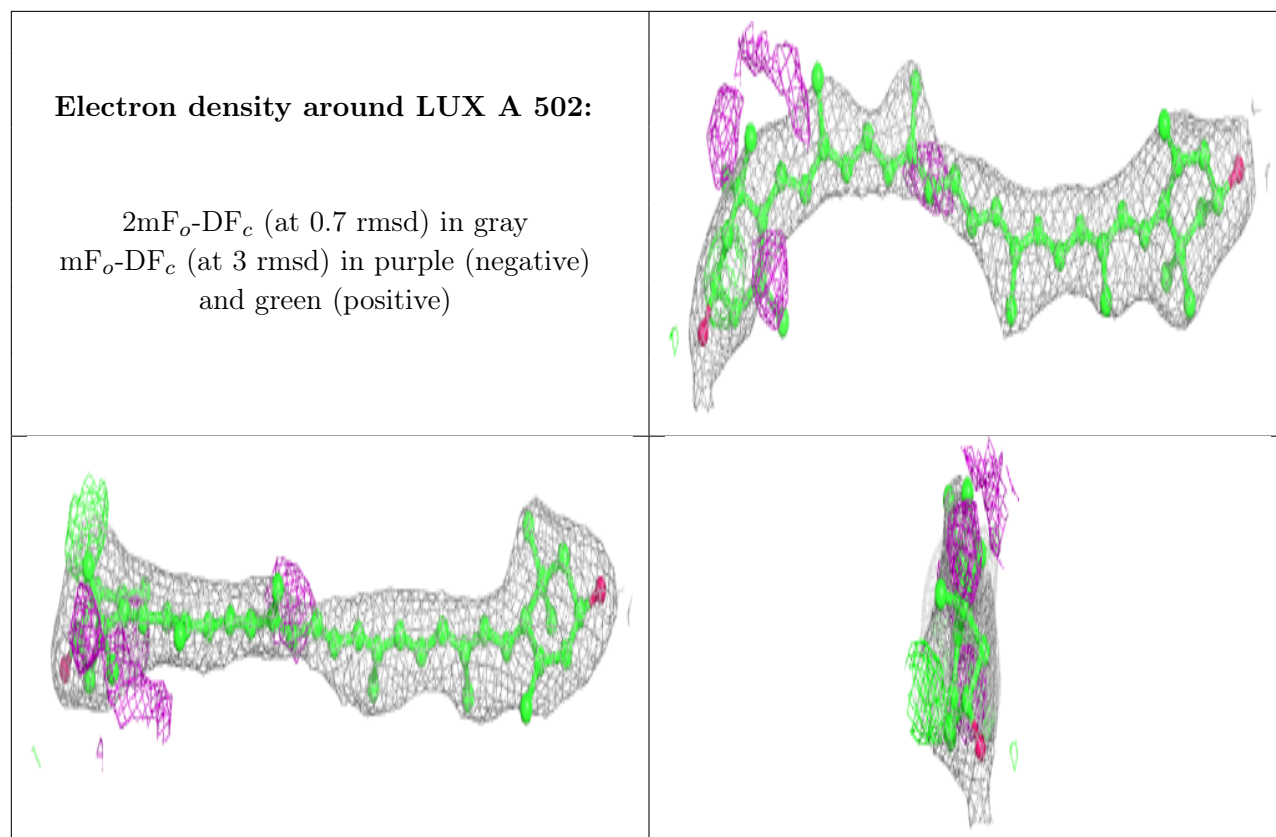
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around CHL A 611:

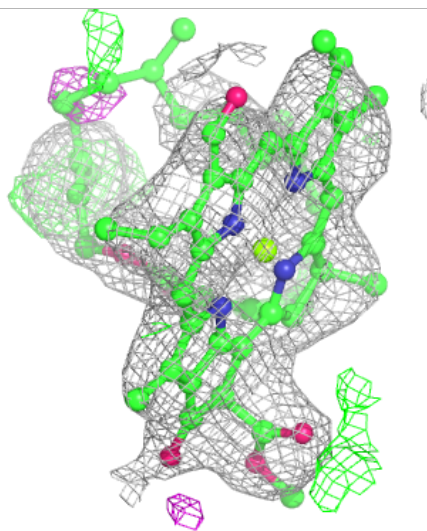
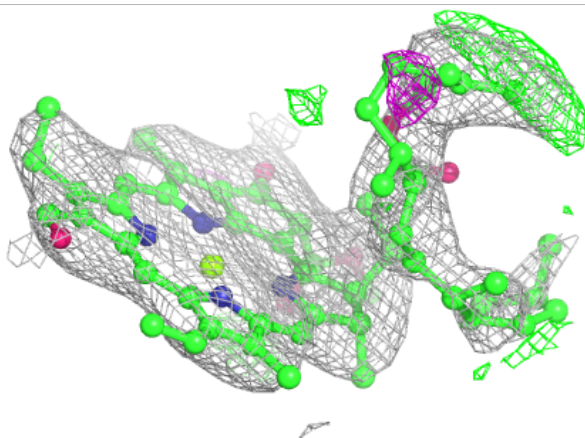
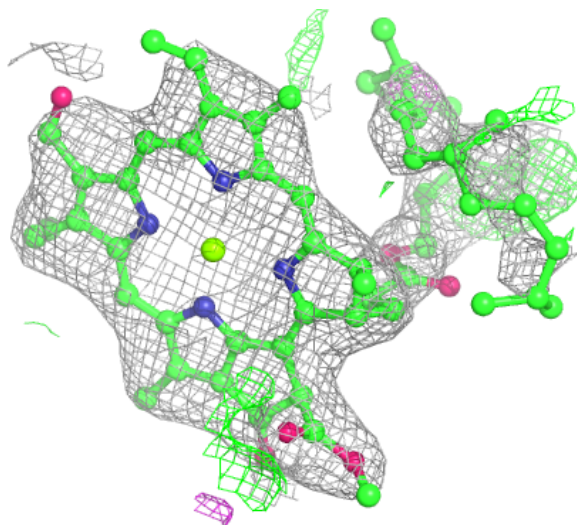
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





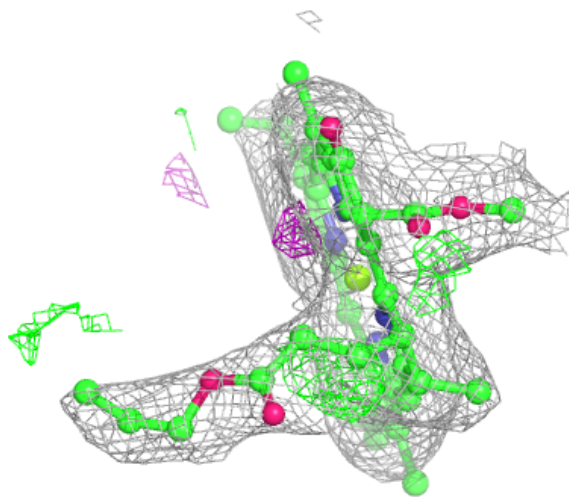
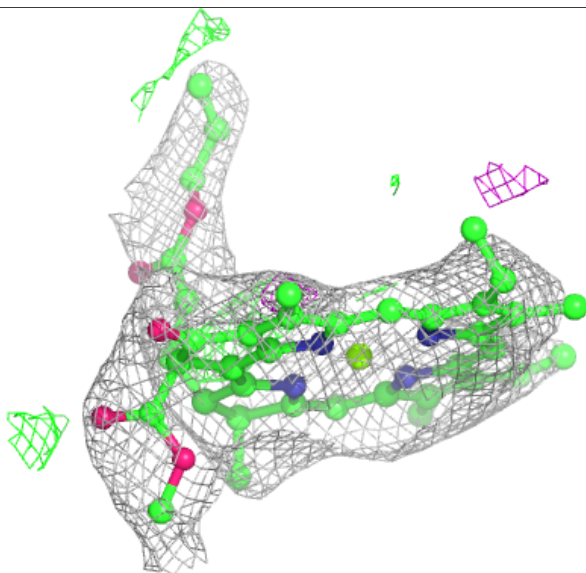
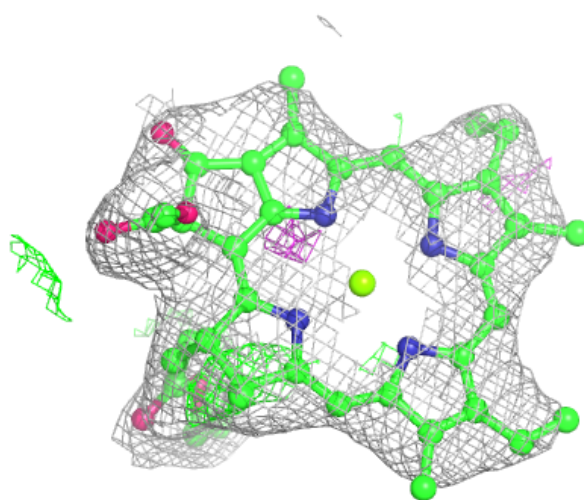
Electron density around CHL C 611:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



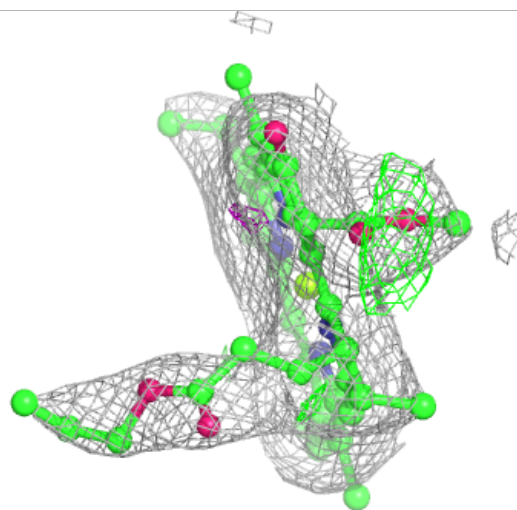
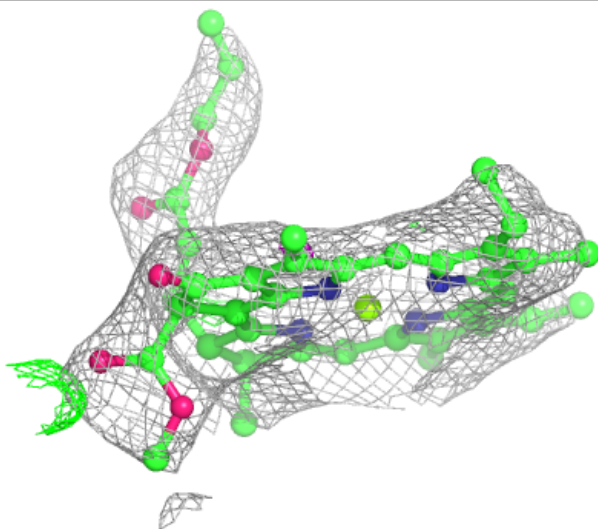
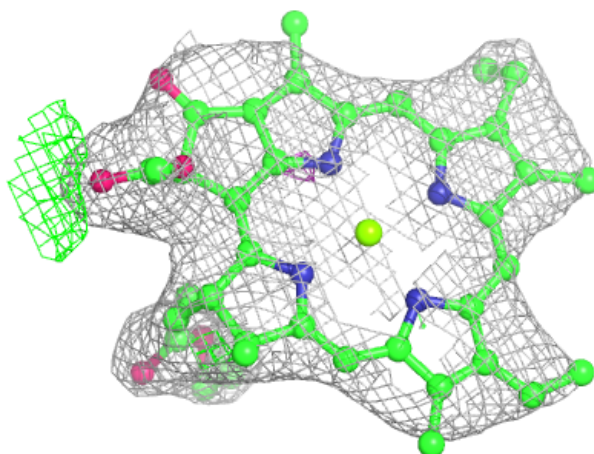
Electron density around CLA B 608:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



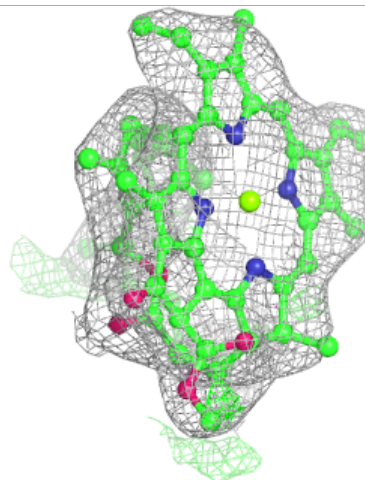
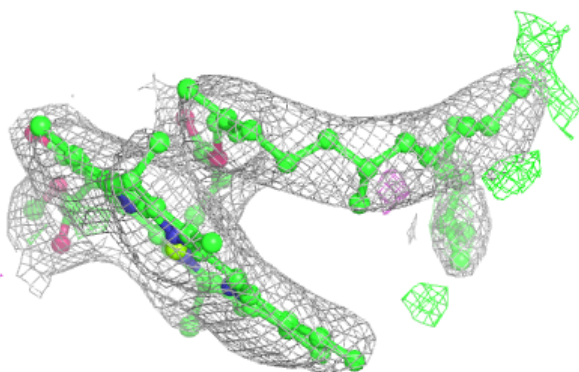
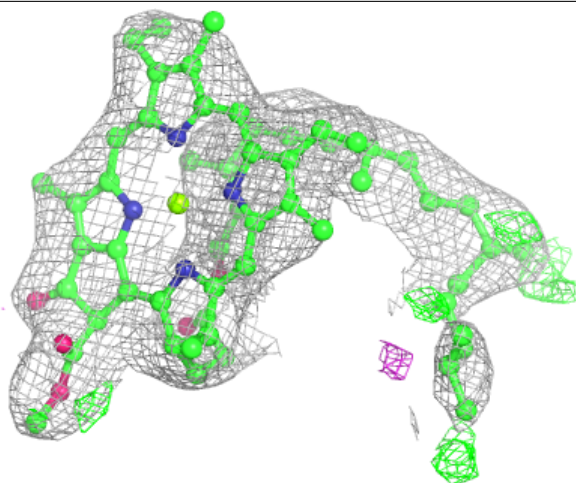
Electron density around CLA C 608:

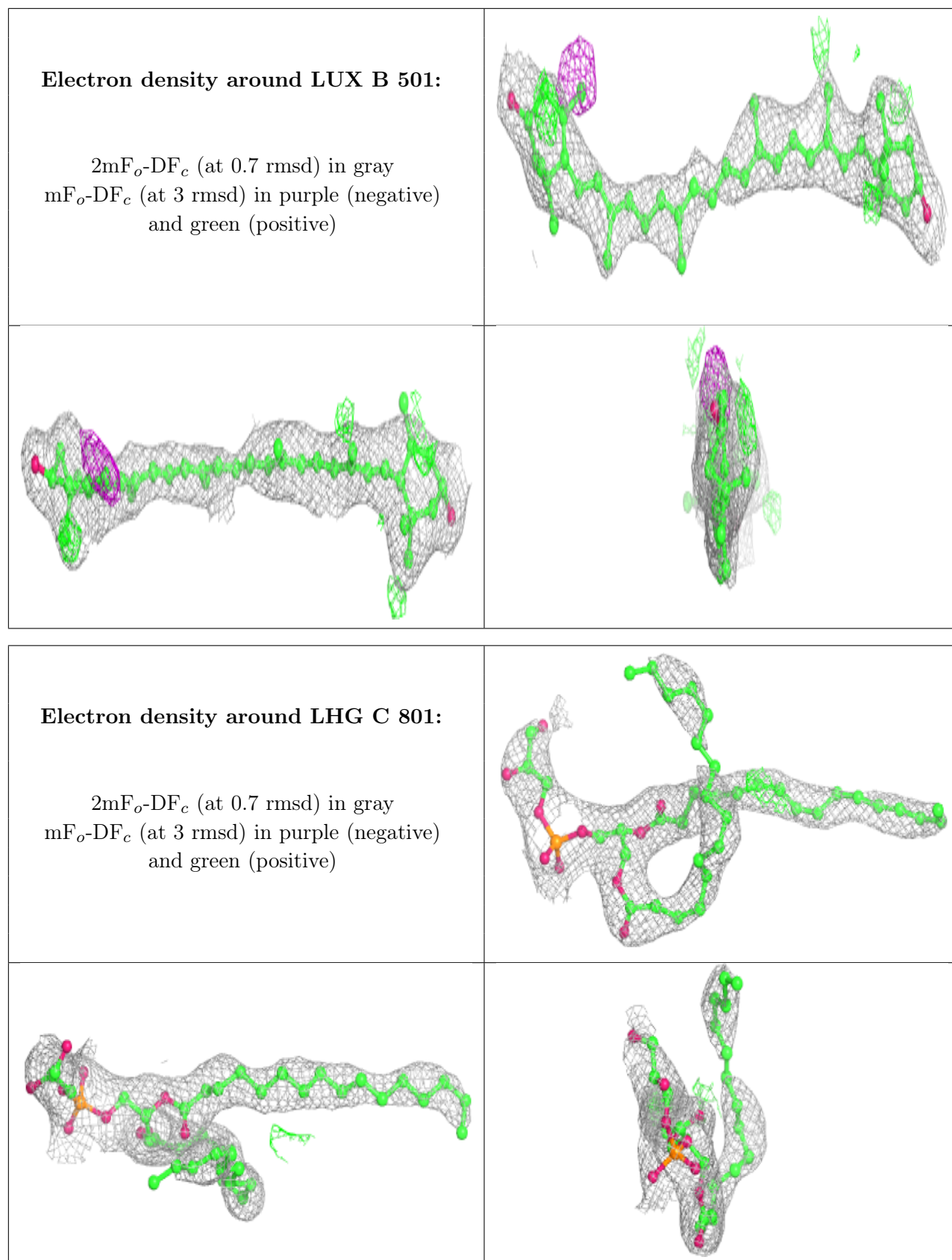
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around CLA B 603:

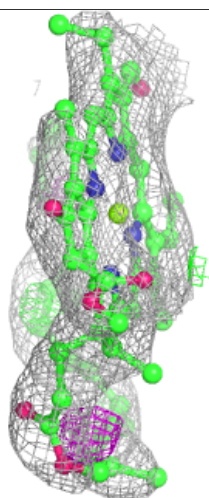
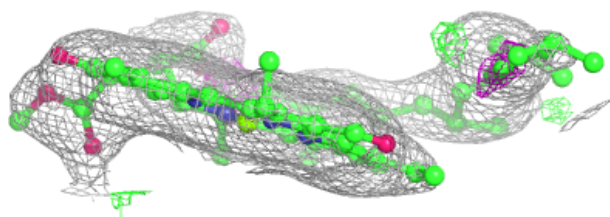
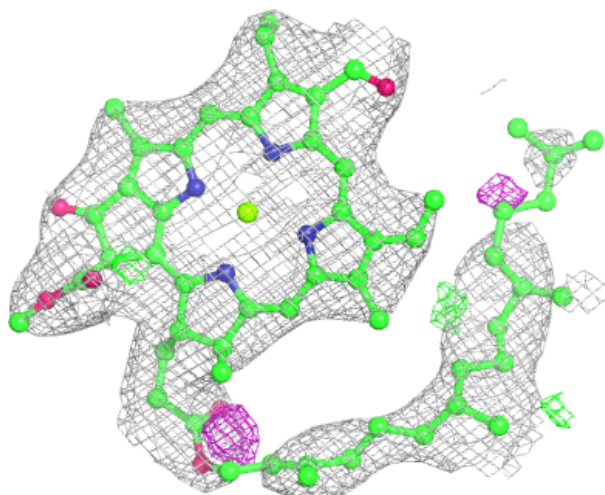
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





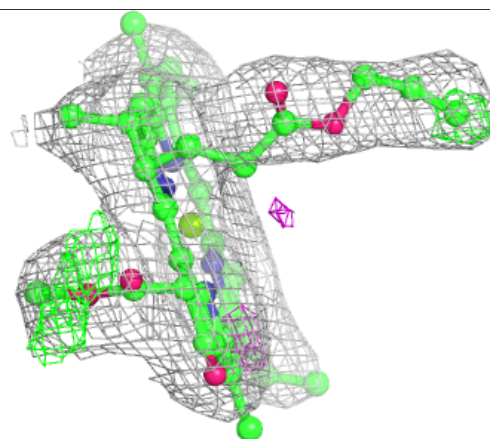
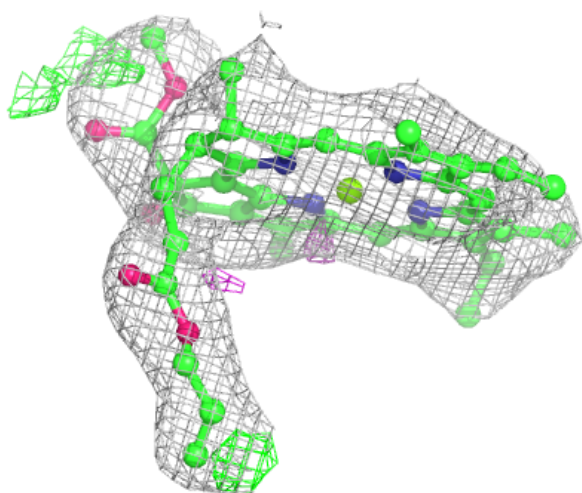
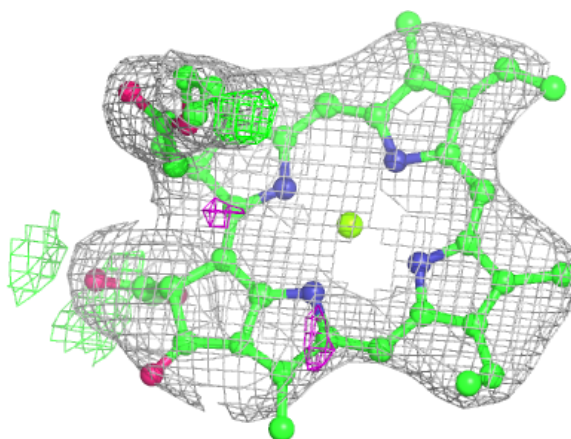
Electron density around CHL C 610:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



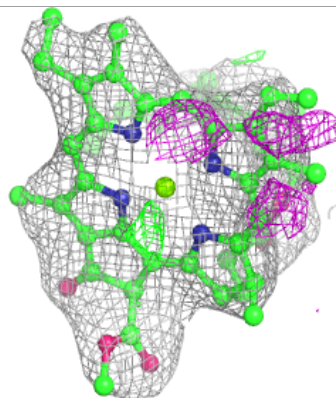
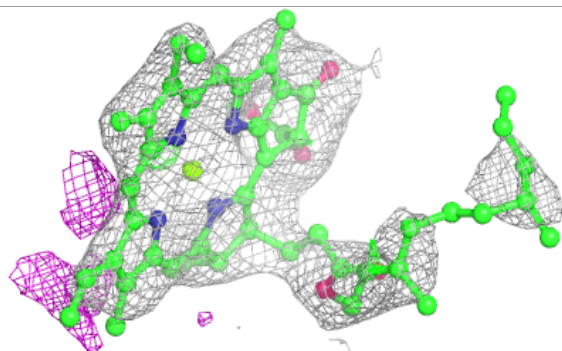
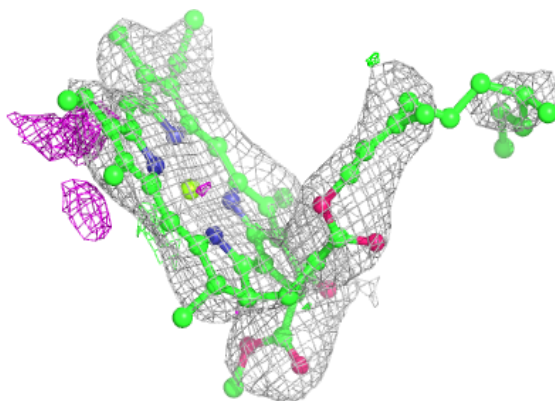
Electron density around CLA A 608:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

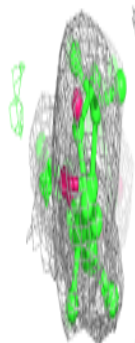
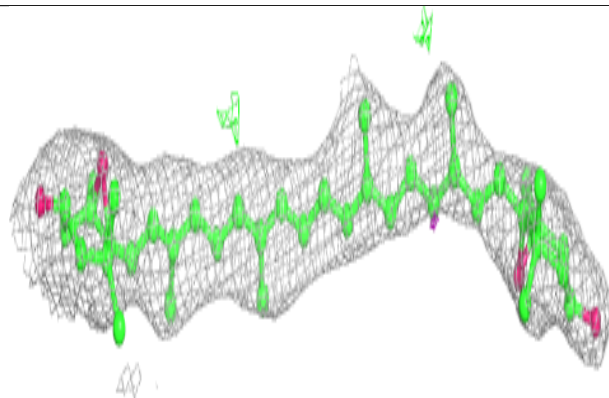
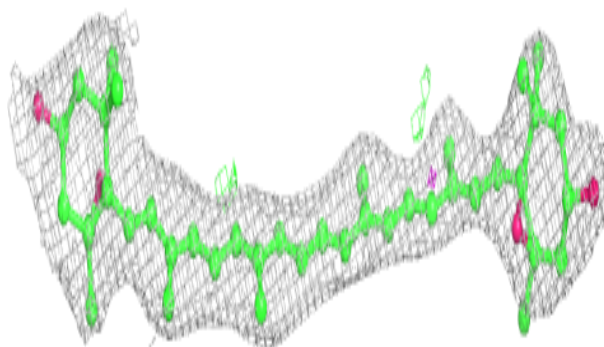


Electron density around CLA A 606:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

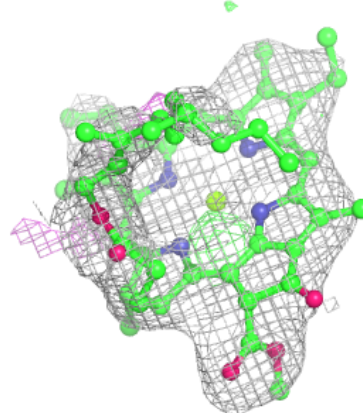
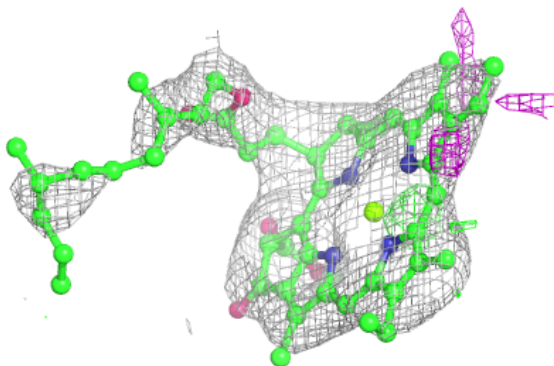
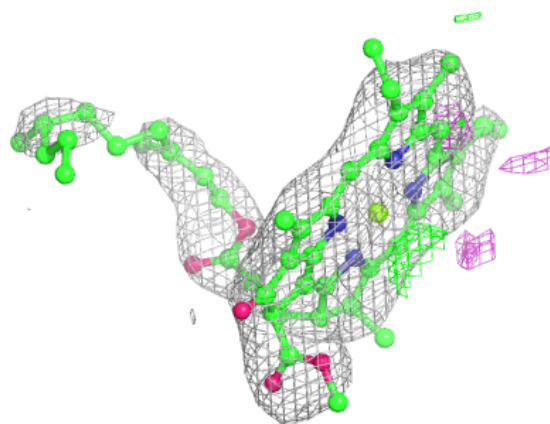
**Electron density around XAT A 504:**

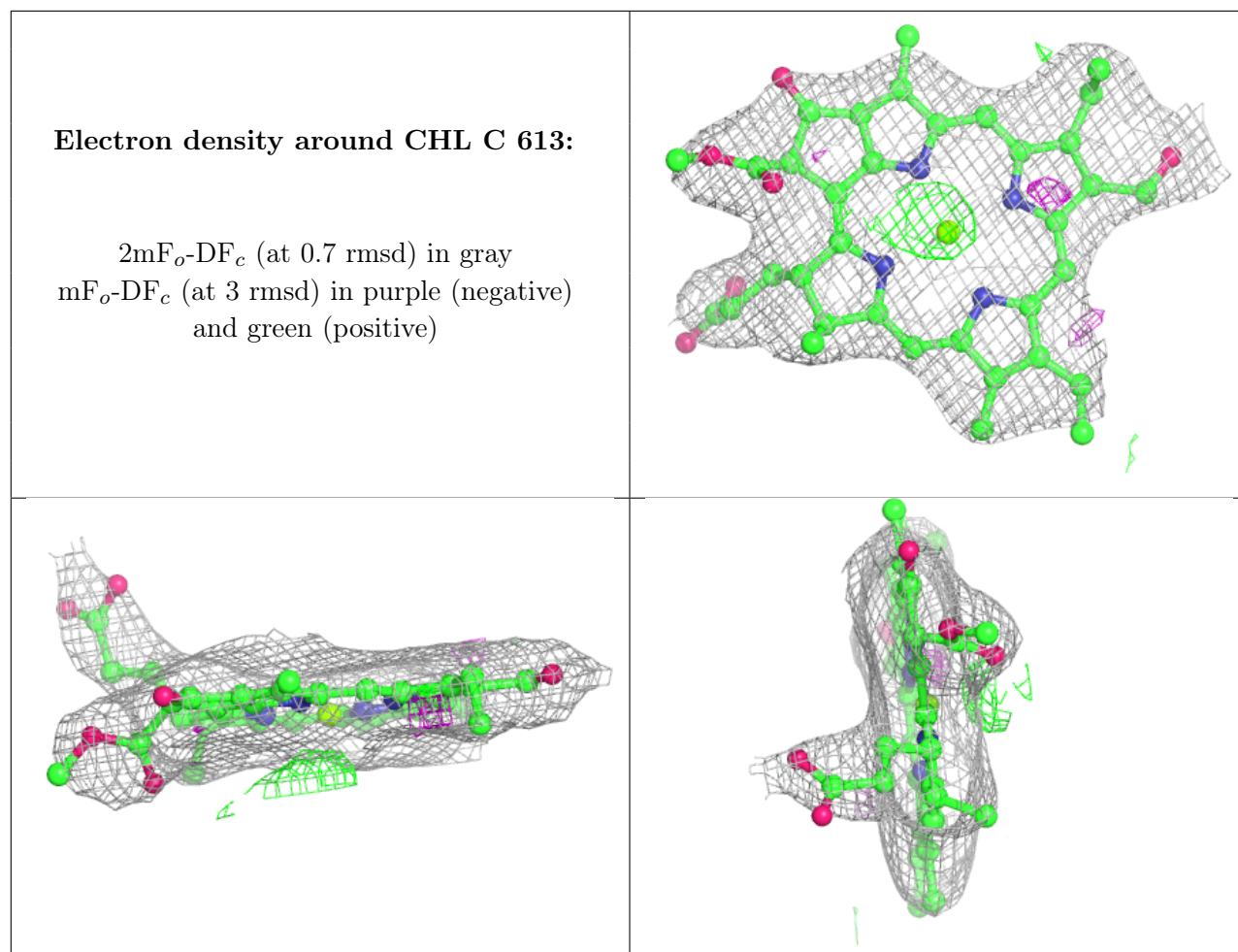
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

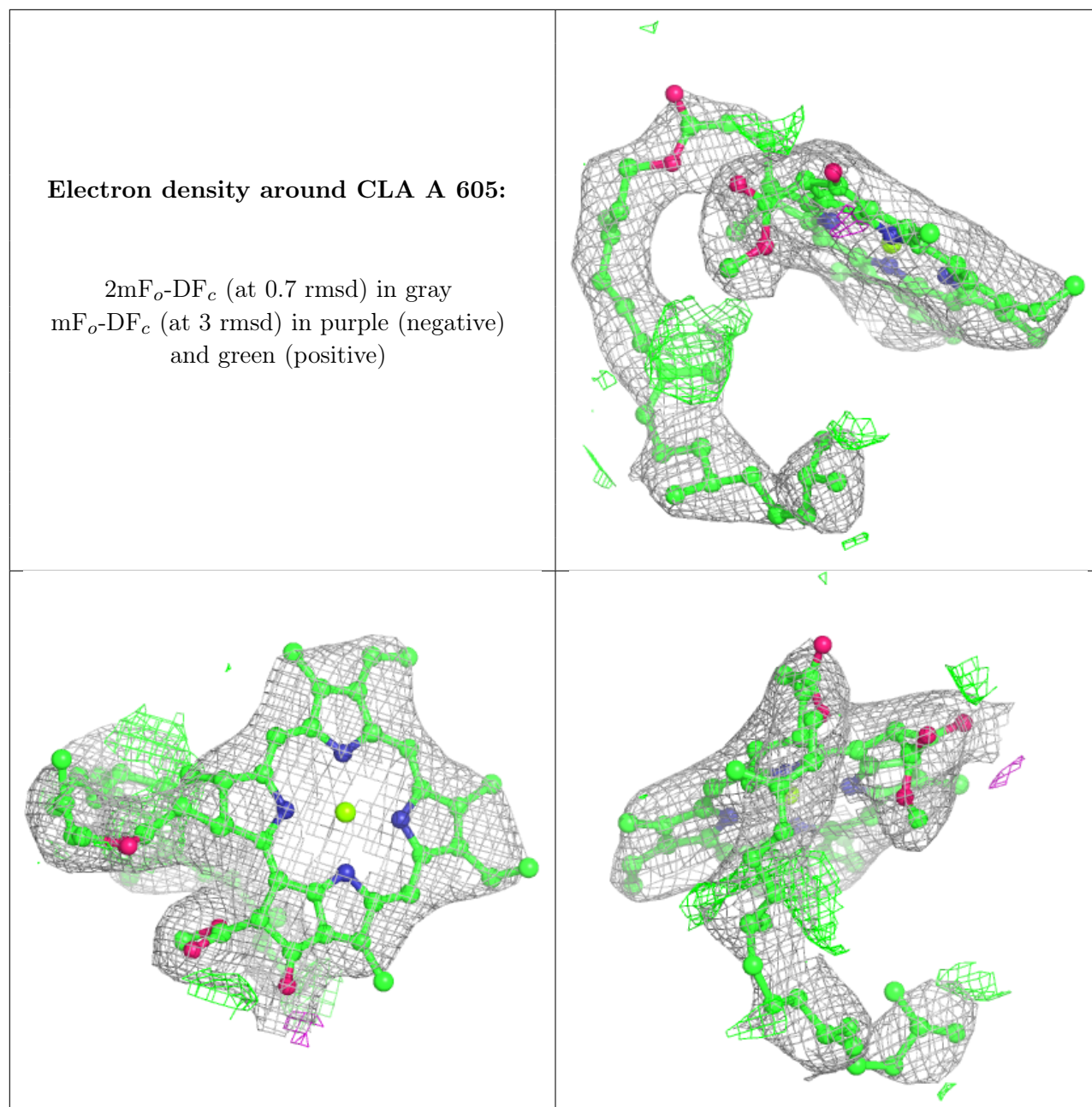


Electron density around CLA C 606:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

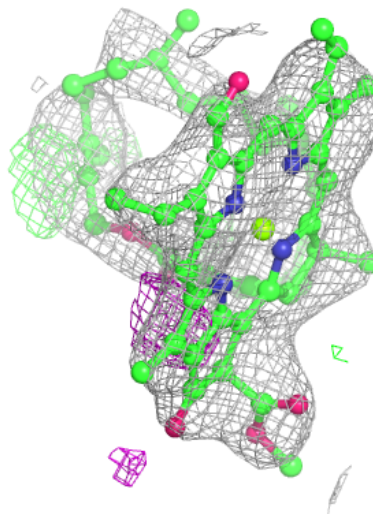
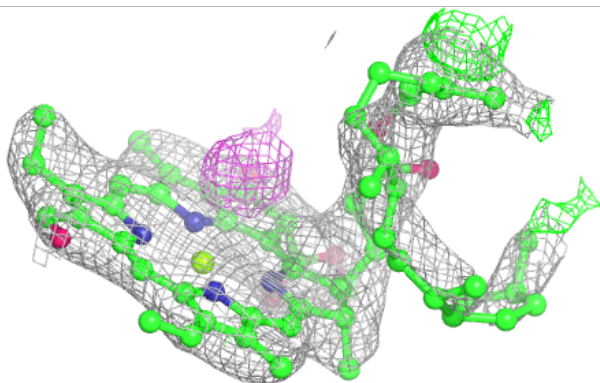
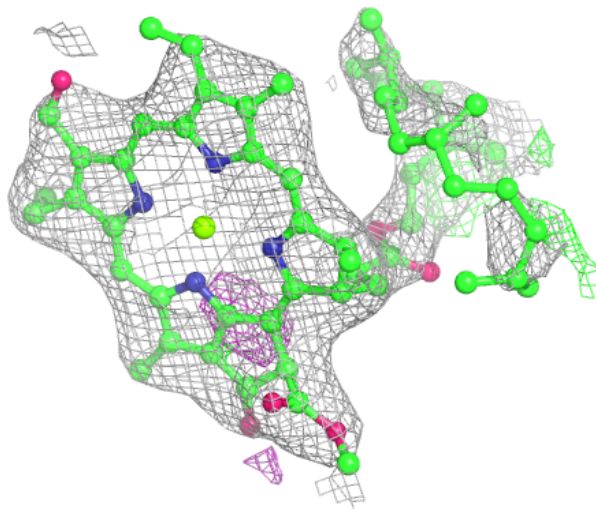


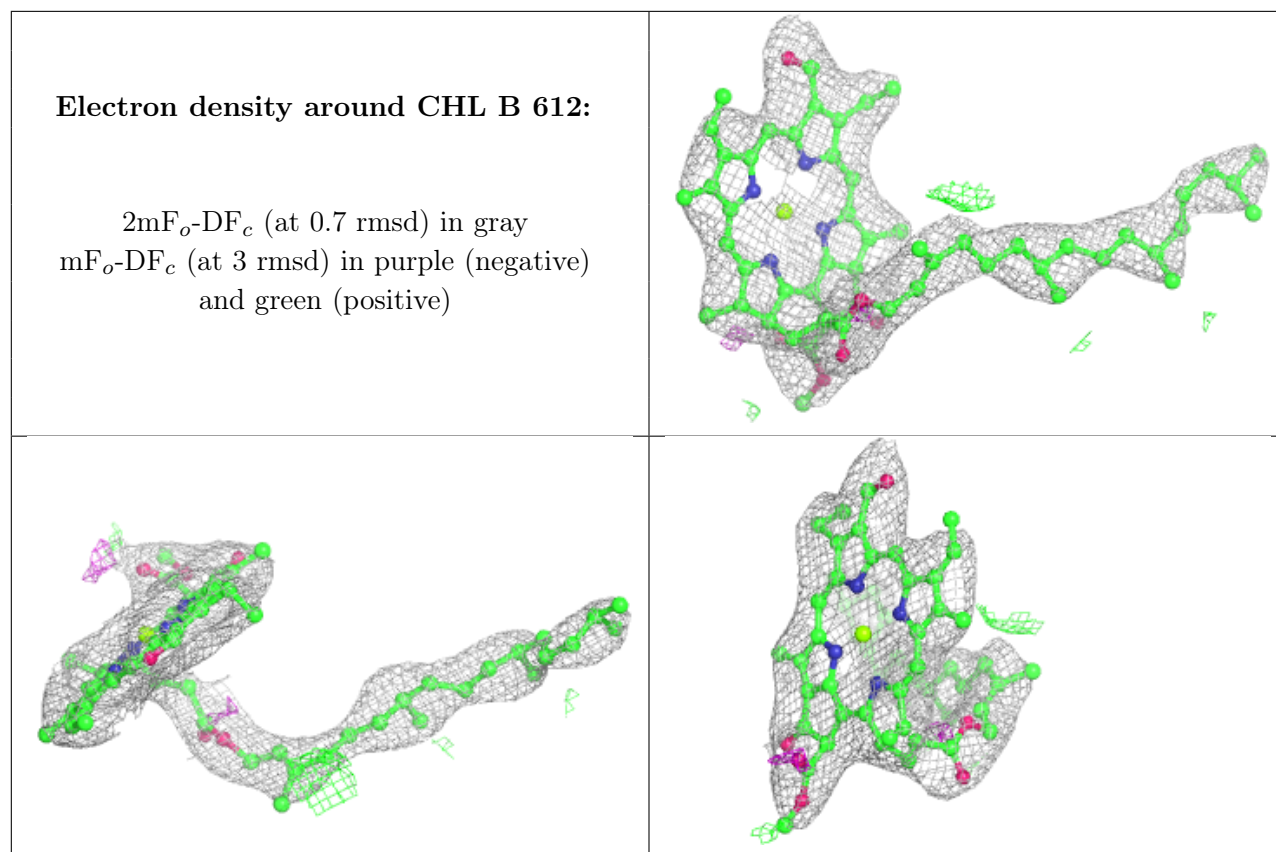


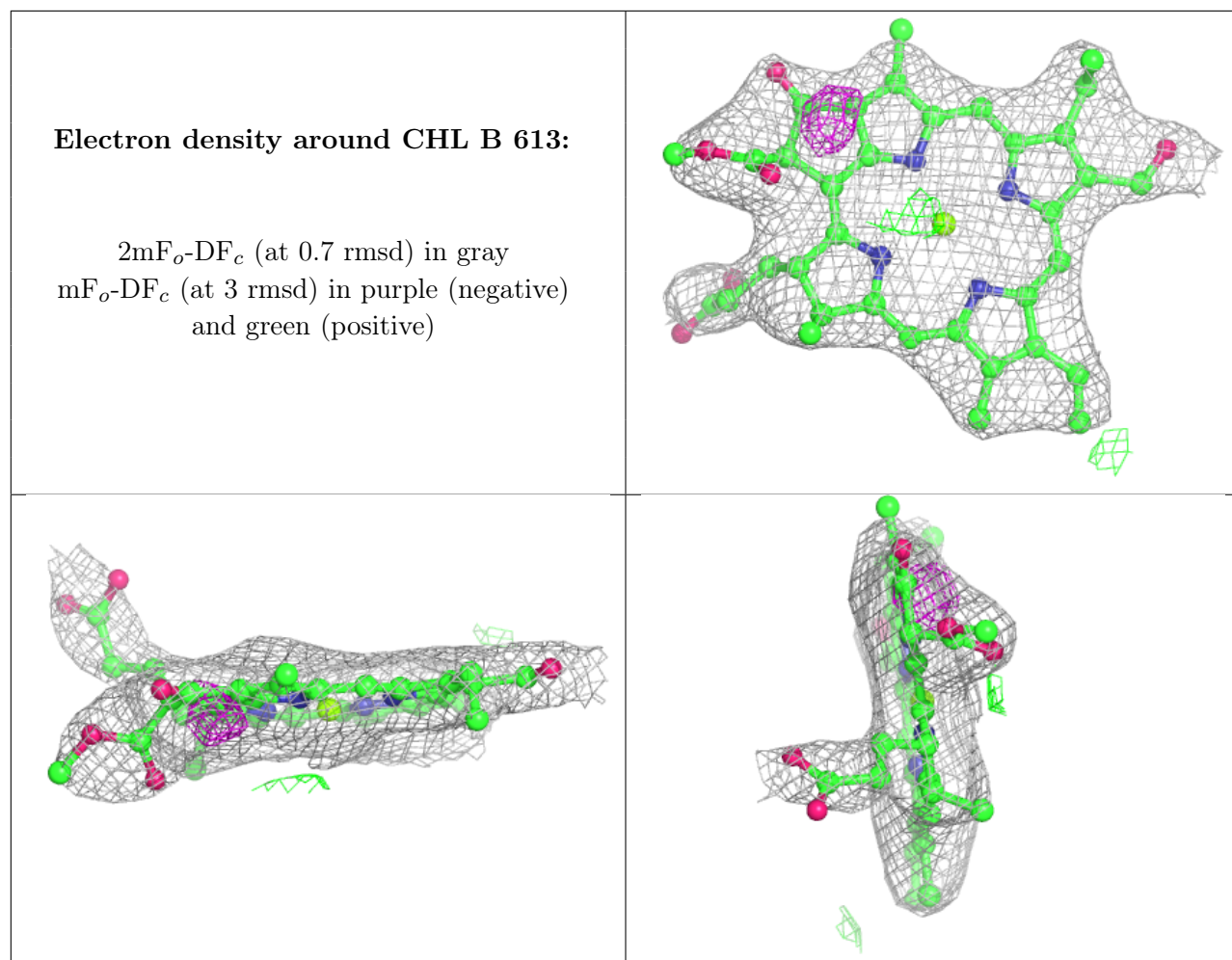


Electron density around CHL B 611:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

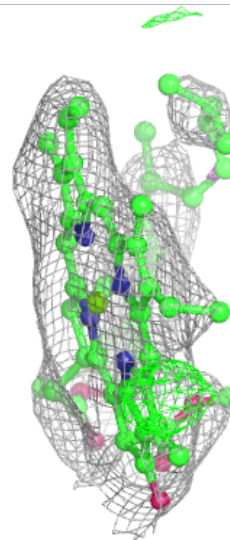
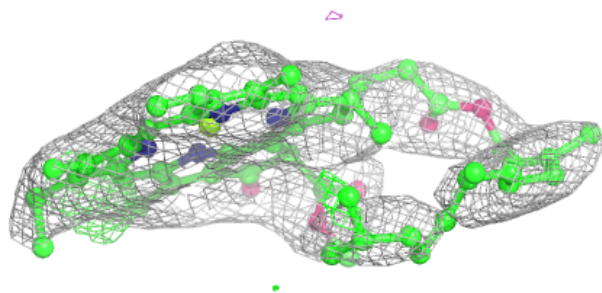
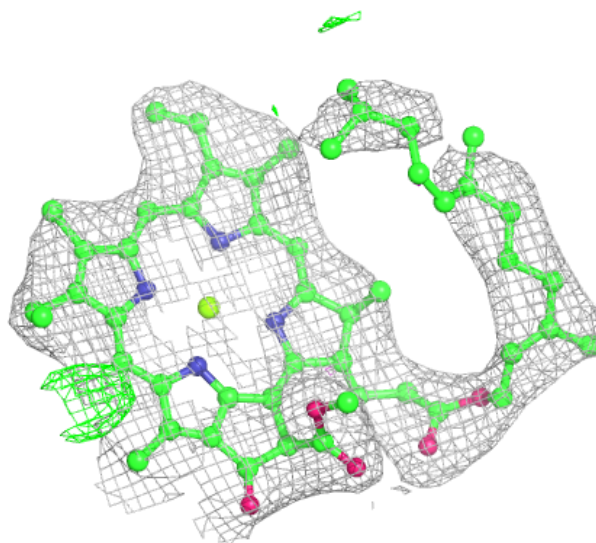


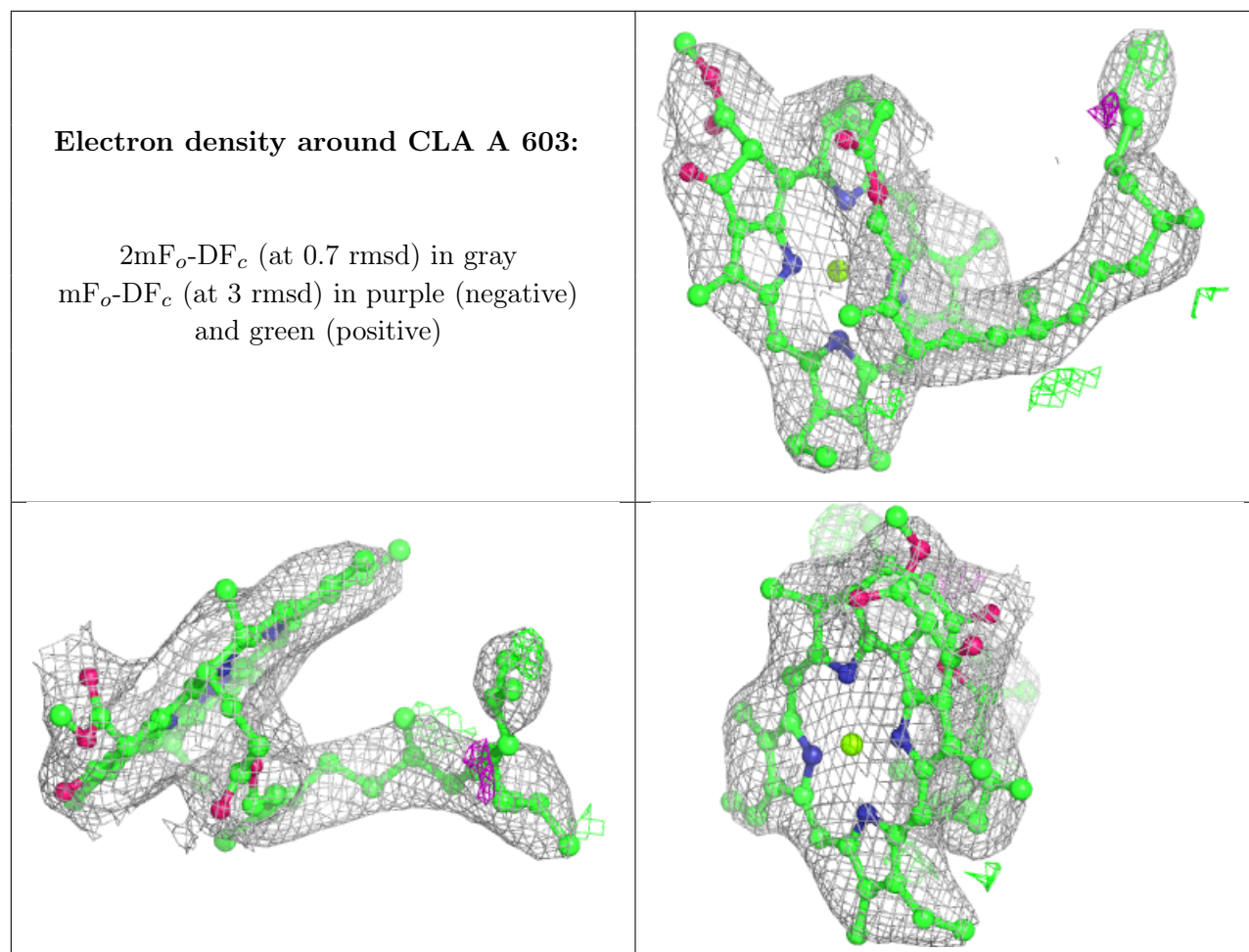




Electron density around CLA A 602:

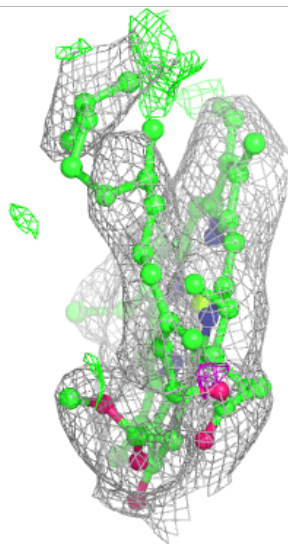
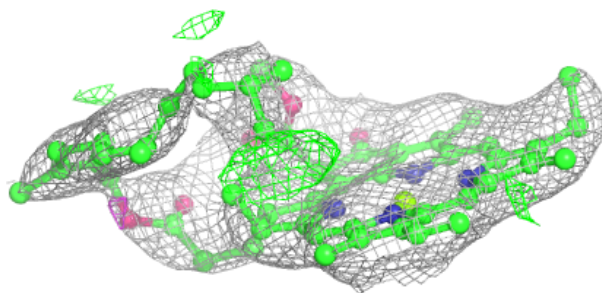
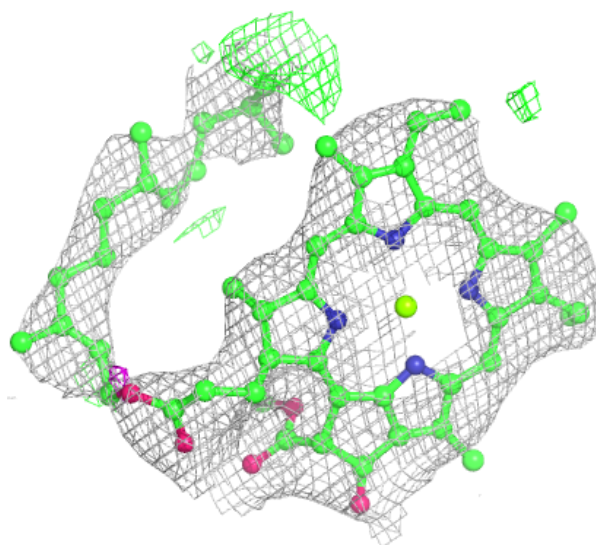
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





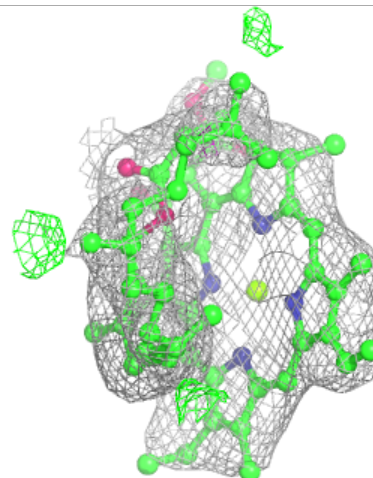
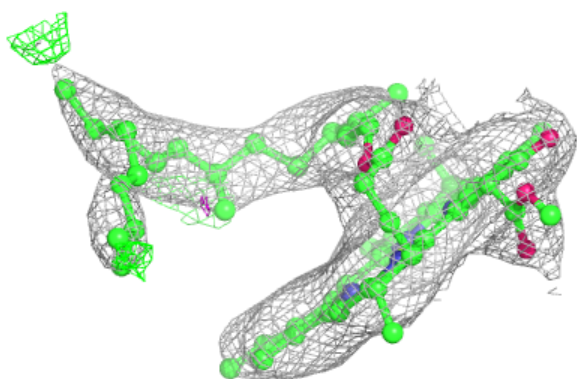
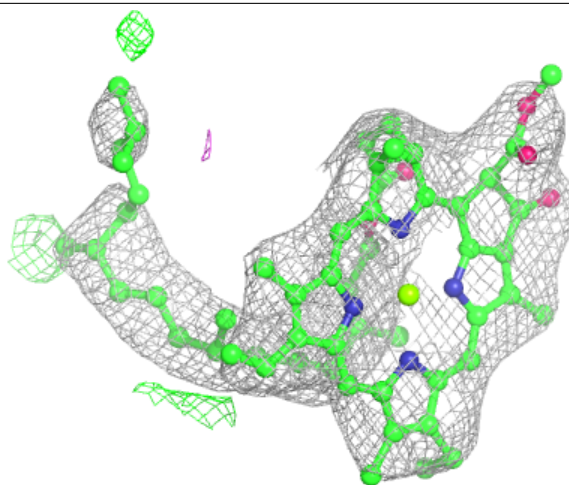
Electron density around CLA C 602:

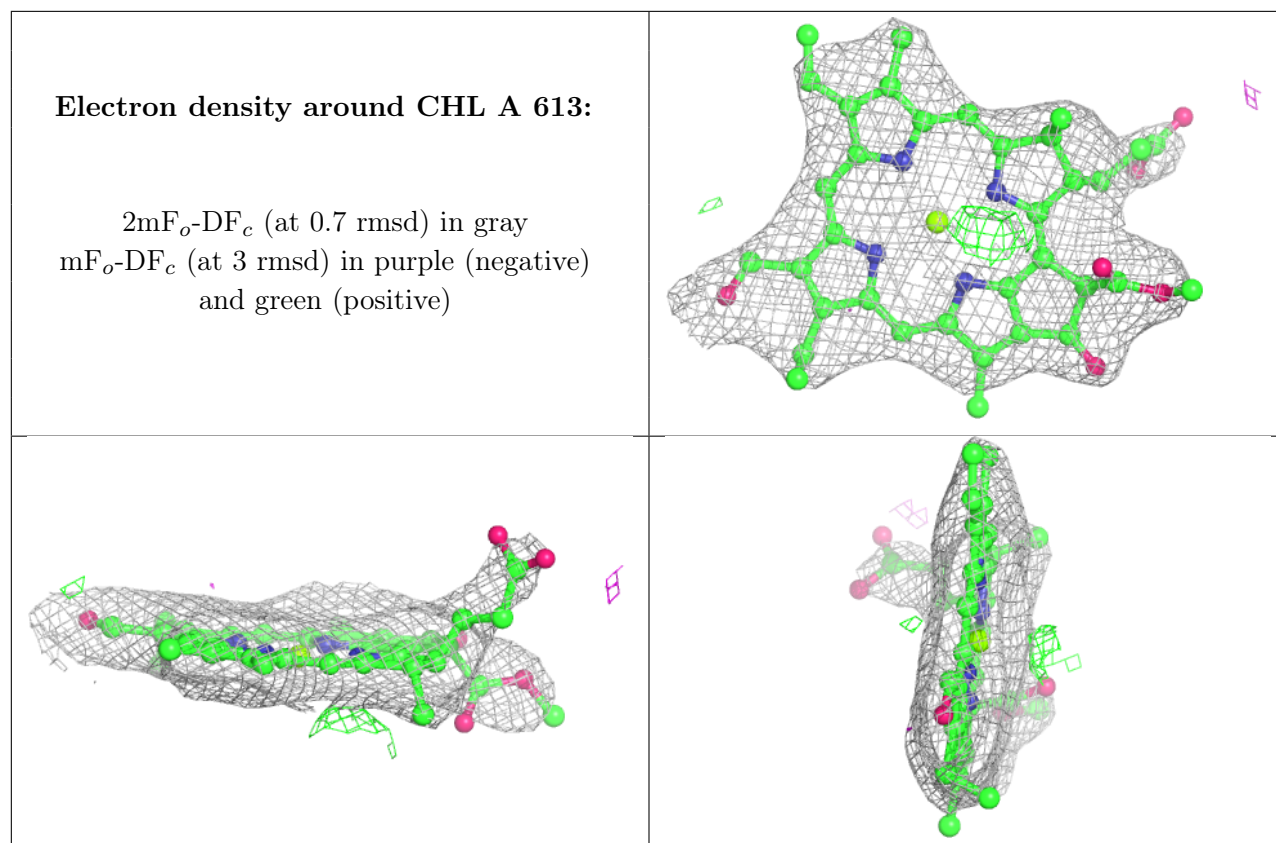
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



Electron density around CLA C 603:

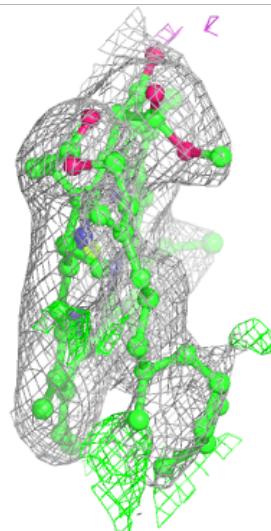
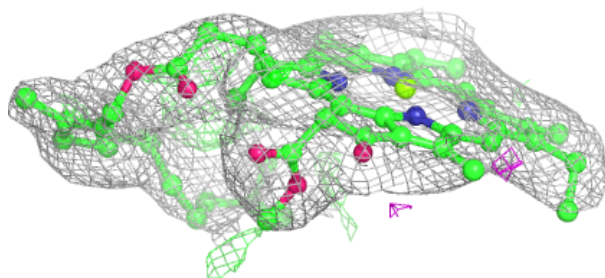
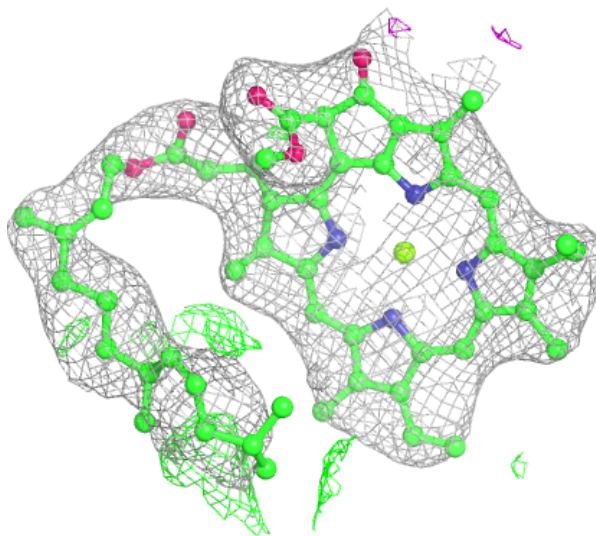
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)





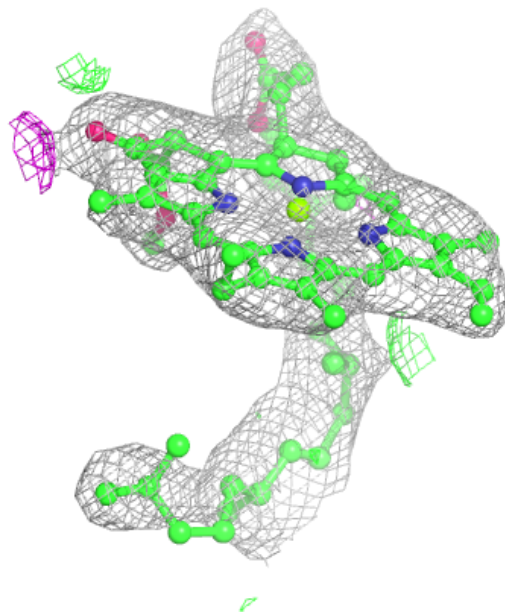
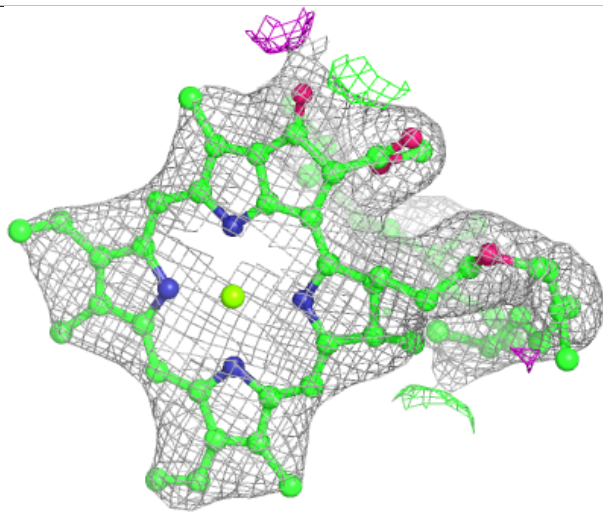
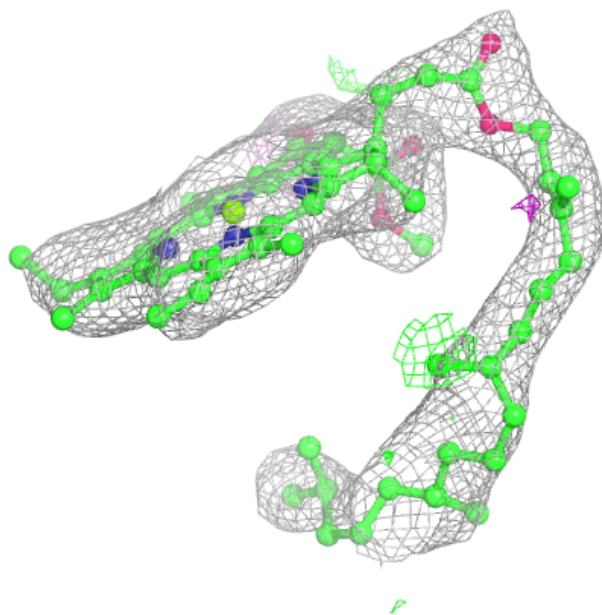
Electron density around CLA B 602:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)



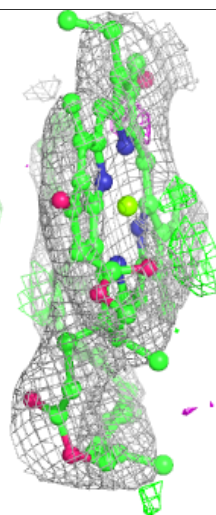
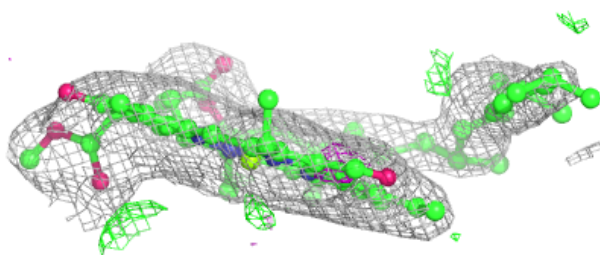
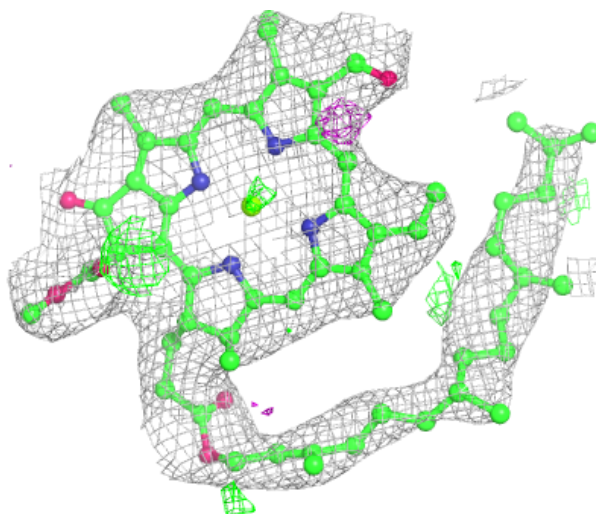
Electron density around CLA C 605:

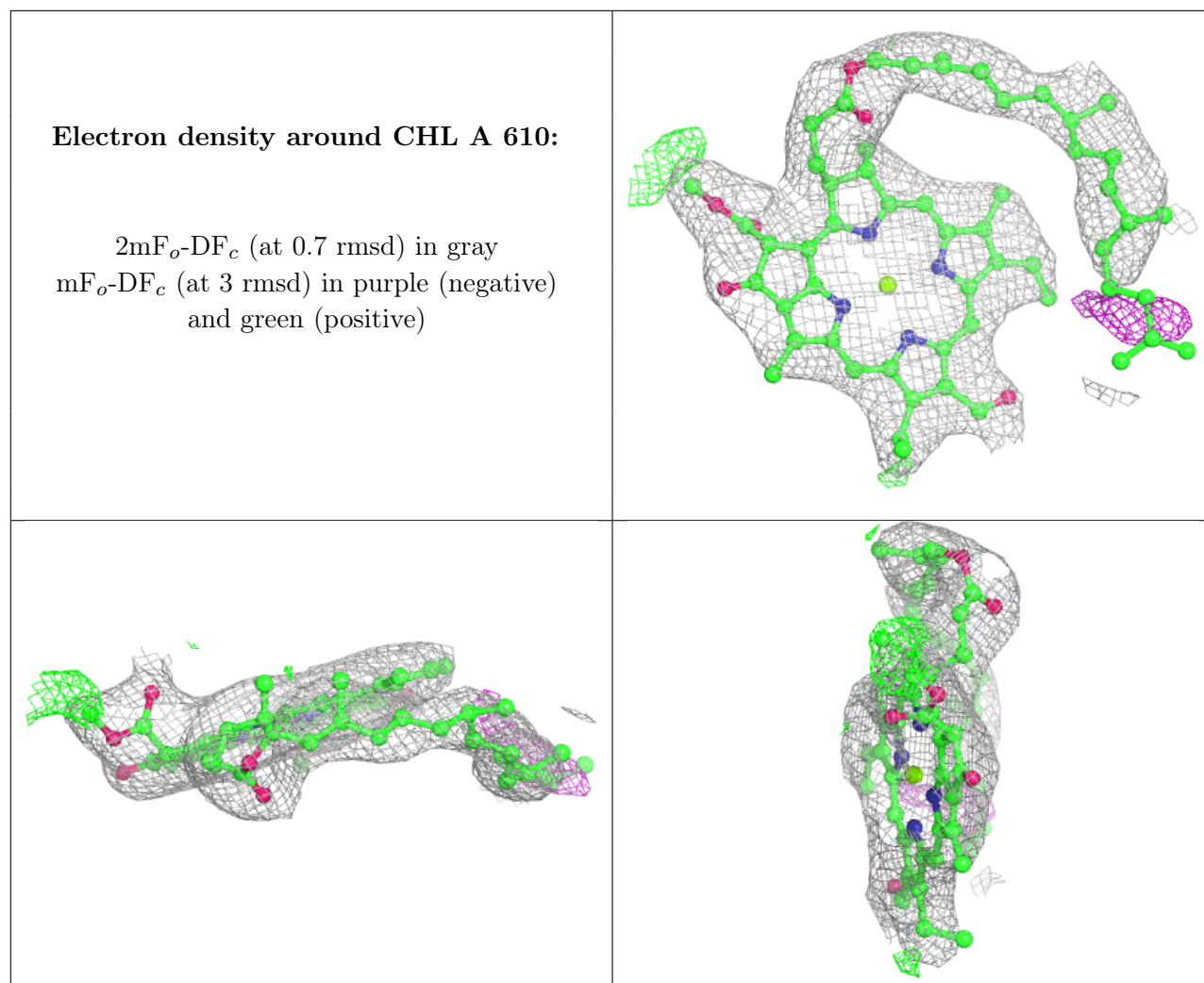
$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

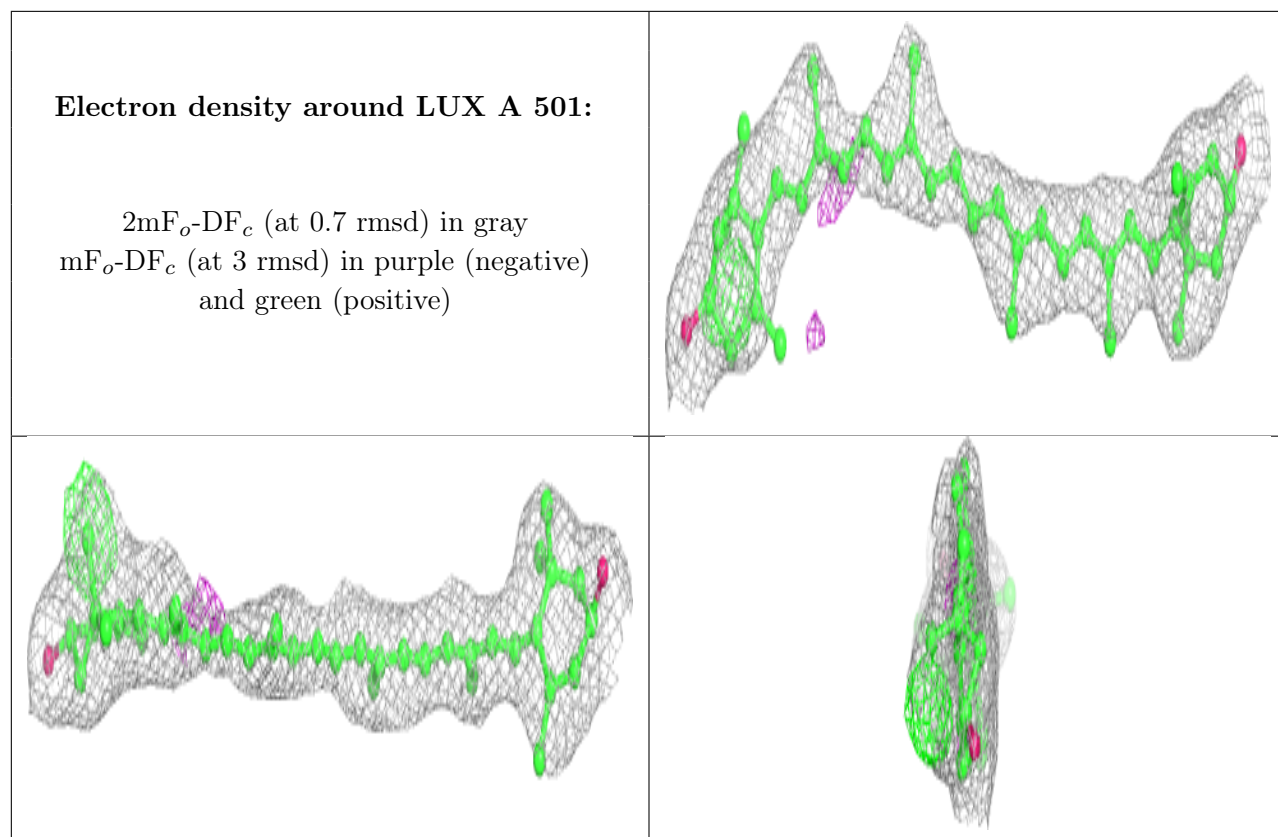


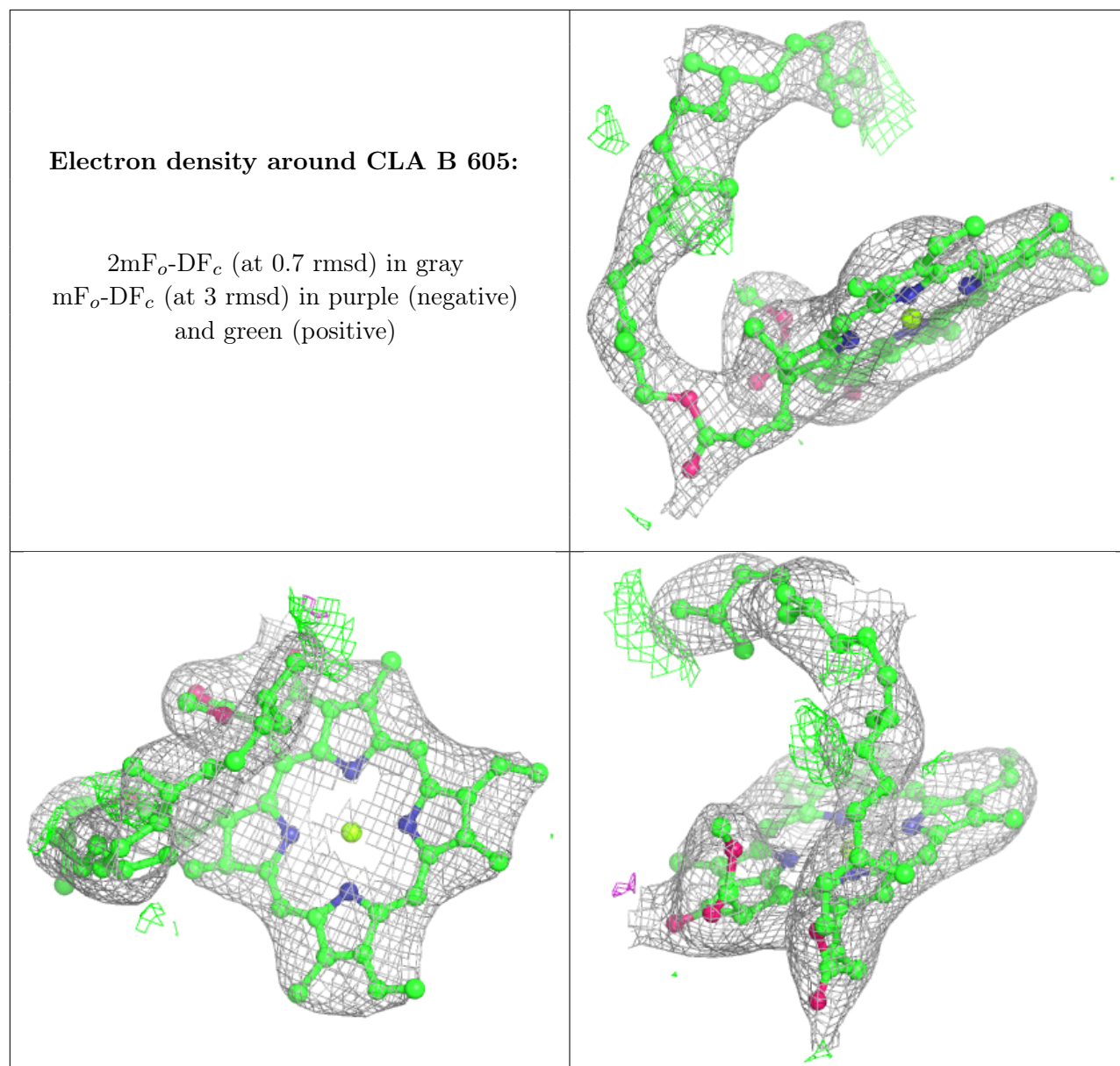
Electron density around CHL B 610:

$2mF_o-DF_c$ (at 0.7 rmsd) in gray
 mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)









6.5 Other polymers [i](#)

There are no such residues in this entry.