wwPDB EM Validation Summary Report (i)

Aug 13, 2026 – 09:04 PM JST

PDB ID	:	8HW1 / pdb_00008hw1
EMDB ID	:	EMD-35080
Title	:	Far-red light-harvesting complex of Antarctic alga Prasiola crispa
Authors	:	Kosugi, M.; Kawasaki, M.; Shibata, Y.; Moriya, T.; Adachi, N.; Senda, T.
Deposited on	:	2022-12-28
Resolution	:	3.13 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 (i) 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 (i)) were used in the production of this report:

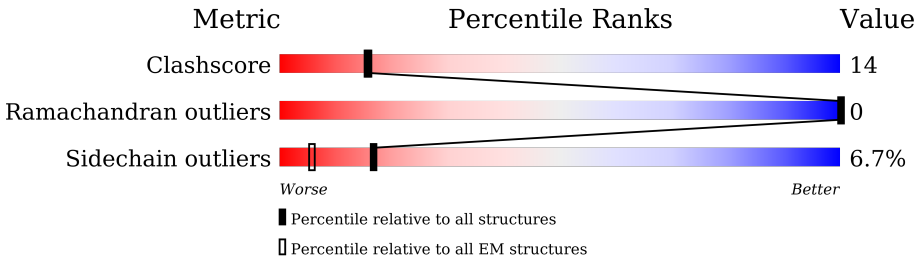
EMDB validation analysis	:	0.0.1.dev133
Mogul	:	1.8.5 (274361), CSD as541be (2020)
MolProbity	:	4-5-2 with Phenix2.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
EM percentile statistics	:	202505.v01 (Using 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)	:	2.50

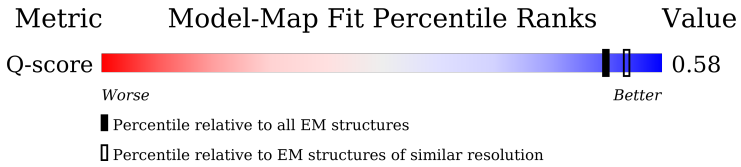
1 Overall quality at a glance (i)

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

The reported resolution of this entry is 3.13 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	14478 (2.63 - 3.63)

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 <=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	307	5% 62% 15% • 21%
1	B	307	6% 62% 14% • 21%
1	C	307	5% 61% 15% • 21%
1	D	307	6% 62% 14% • 21%

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Mol	Chain	Length	Quality of chain
1	E	307	6% 62% 15% 21%
1	F	307	5% 63% 14% 21%
1	G	307	5% 61% 15% 21%
1	H	307	6% 62% 15% 21%
1	I	307	6% 62% 14% 21%
1	J	307	6% 61% 15% 21%
1	K	307	5% 61% 16% 21%

2 Entry composition (i)

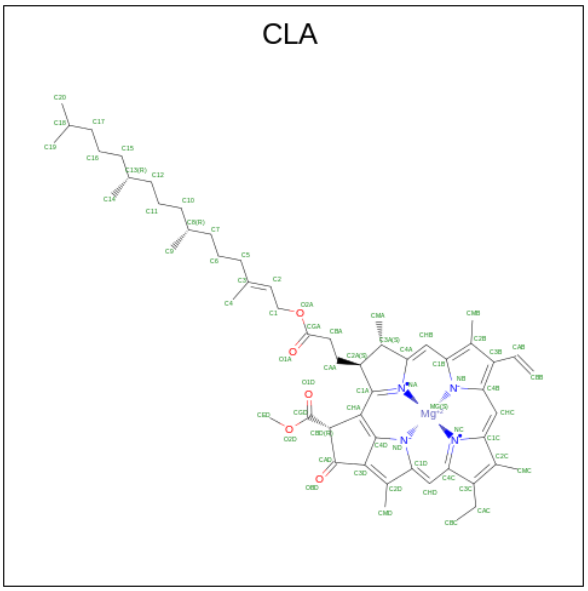
There are 4 unique types of molecules in this entry. The entry contains 29271 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Far-red light-harvesting complex.

Mol	Chain	Residues	Atoms					AltConf	Trace
1	A	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		
1	B	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		
1	C	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		
1	D	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		
1	E	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		
1	F	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		
1	G	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		
1	H	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		
1	I	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		
1	J	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		
1	K	241	Total	C	N	O	S	0	0
			1886	1214	321	346	5		

- Molecule 2 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
2	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
2	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
2	A	1	Total 56	C 46	Mg 1	N 4	O 5	0
2	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
2	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
2	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
2	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
2	A	1	Total 56	C 46	Mg 1	N 4	O 5	0
2	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
2	A	1	Total 65	C 55	Mg 1	N 4	O 5	0
2	A	1	Total 56	C 46	Mg 1	N 4	O 5	0
2	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
2	B	1	Total 65	C 55	Mg 1	N 4	O 5	0
2	B	1	Total 56	C 46	Mg 1	N 4	O 5	0

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Mol	Chain	Residues	Atoms					AltConf
2	B	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	B	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	B	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	B	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	B	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	B	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	B	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	B	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	C	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	D	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	D	1	Total	C	Mg	N	O	0
			65	55	1	4	5	

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Mol	Chain	Residues	Atoms					AltConf
2	D	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	D	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	D	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	D	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	D	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	D	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	D	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	D	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	D	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	E	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	E	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	E	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	E	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	E	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	E	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	E	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	E	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	E	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	E	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	F	1	Total	C	Mg	N	O	0
			65	55	1	4	5	

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Mol	Chain	Residues	Atoms					AltConf
2	F	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	F	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	F	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	F	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	F	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	F	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	F	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	F	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	F	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	F	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	G	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	G	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	G	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	G	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	G	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	G	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	G	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	G	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	G	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	G	1	Total	C	Mg	N	O	0
			56	46	1	4	5	

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Mol	Chain	Residues	Atoms					AltConf
2	H	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	H	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	H	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	H	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	H	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	H	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	H	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	H	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	H	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	H	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	H	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	I	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	I	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	I	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	I	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	I	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	I	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	I	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	I	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	I	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	I	1	Total	C	Mg	N	O	0
			65	55	1	4	5	

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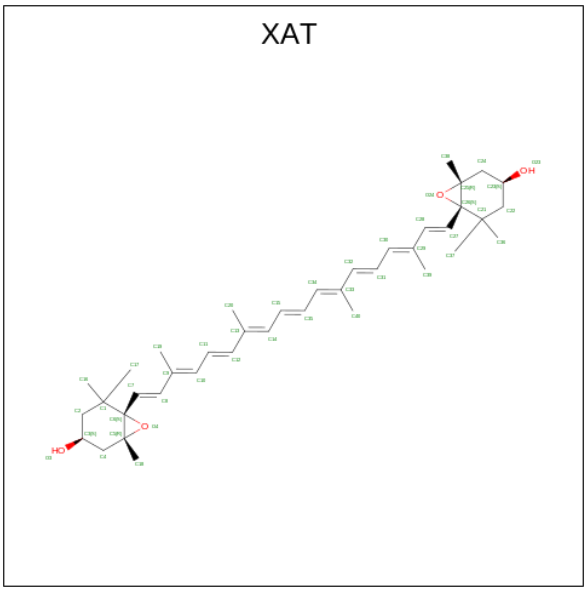
Mol	Chain	Residues	Atoms					AltConf
2	I	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	J	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	K	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	K	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	K	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	K	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	K	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	K	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	K	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	K	1	Total	C	Mg	N	O	0
			56	46	1	4	5	
2	K	1	Total	C	Mg	N	O	0
			65	55	1	4	5	

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Mol	Chain	Residues	Atoms					AltConf
2	K	1	Total	C	Mg	N	O	0
			65	55	1	4	5	
2	K	1	Total	C	Mg	N	O	0
			56	46	1	4	5	

- Molecule 3 is (3S,5R,6S,3'S,5'R,6'S)-5,6,5',6'-DIEPOXY-5,6,5',6'- TETRAHYDRO-BETA ,BETA-CAROTENE-3,3'-DIOL (CCD ID: XAT) (formula: $C_{40}H_{56}O_4$) (labeled as "Ligand of Interest" by depositor).



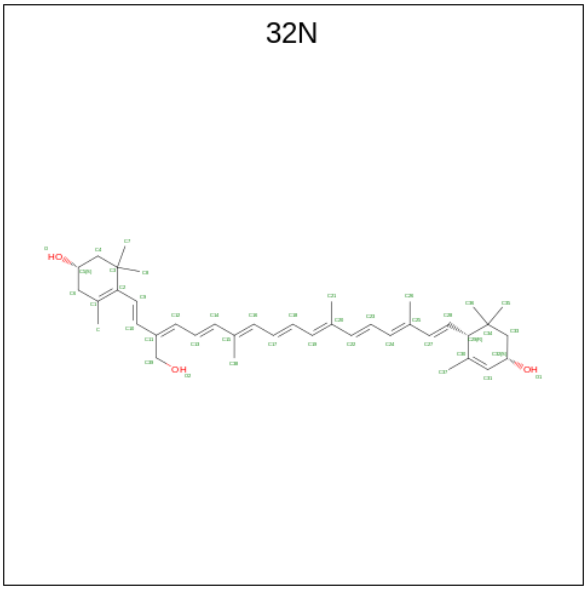
Mol	Chain	Residues	Atoms			AltConf
3	A	1	Total	C	O	0
			44	40	4	
3	B	1	Total	C	O	0
			44	40	4	
3	C	1	Total	C	O	0
			44	40	4	
3	D	1	Total	C	O	0
			44	40	4	
3	E	1	Total	C	O	0
			44	40	4	
3	F	1	Total	C	O	0
			44	40	4	
3	G	1	Total	C	O	0
			44	40	4	
3	H	1	Total	C	O	0
			44	40	4	

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Mol	Chain	Residues	Atoms	AltConf
3	I	1	Total C O 44 40 4	0
3	J	1	Total C O 44 40 4	0
3	K	1	Total C O 44 40 4	0

- Molecule 4 is (1S)-4-[(1E,3Z,5E,7E,9E,11E,13E,15E,17E)-3-(hydroxymethyl)-7,12,16-trimethyl-18-[(1R,4S)-2,6,6-trimethyl-4-oxidanyl-cyclohex-2-en-1-yl]octadeca-1,3,5,7,9,11,13,15,17-nonaenyl]-3,5,5-trimethyl-cyclohex-3-en-1-ol (CCD ID: 32N) (formula: $C_{40}H_{56}O_3$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms	AltConf
4	A	1	Total C O 43 40 3	0
4	B	1	Total C O 43 40 3	0
4	C	1	Total C O 43 40 3	0
4	D	1	Total C O 43 40 3	0
4	E	1	Total C O 43 40 3	0
4	F	1	Total C O 43 40 3	0
4	G	1	Total C O 43 40 3	0

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Mol	Chain	Residues	Atoms			AltConf
4	H	1	Total	C	O	0
			43	40	3	
4	I	1	Total	C	O	0
			43	40	3	
4	J	1	Total	C	O	0
			43	40	3	
4	K	1	Total	C	O	0
			43	40	3	

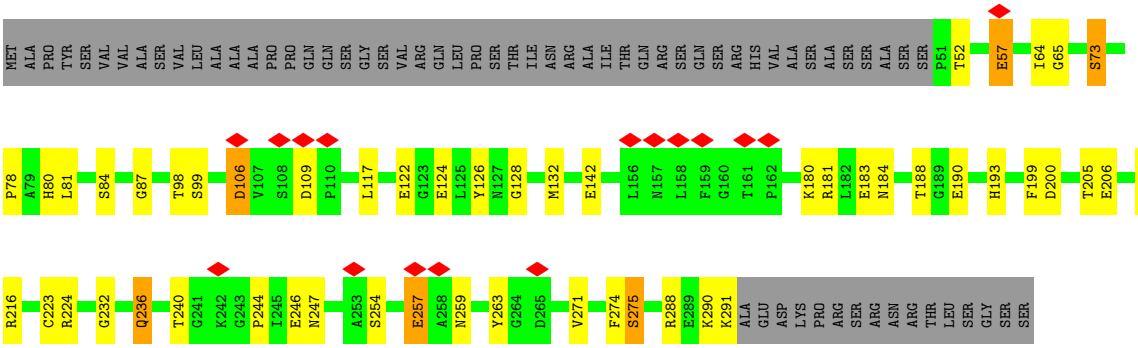
3 Residue-property plots (i)

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 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: Far-red light-harvesting complex

Chain A:

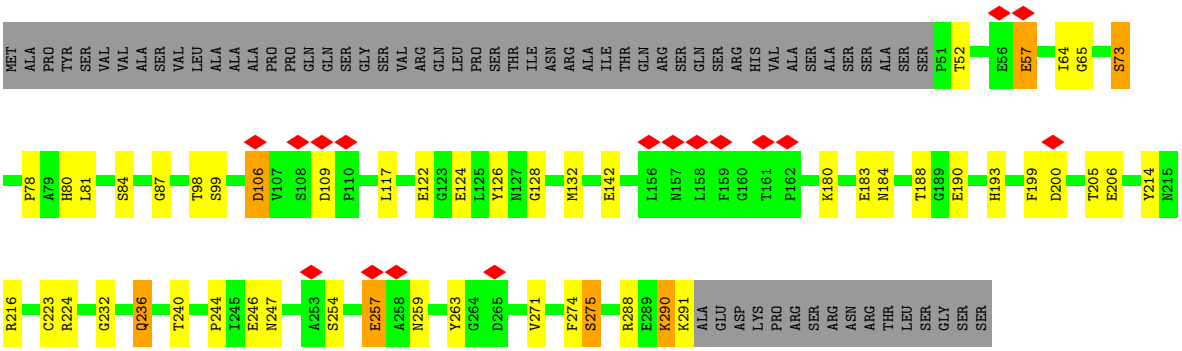




• Molecule 1: Far-red light-harvesting complex

Chain B:

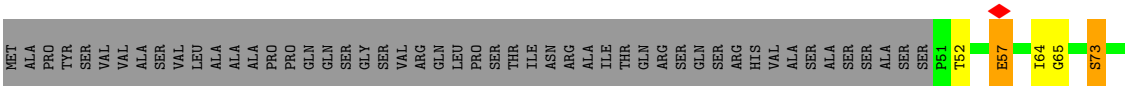


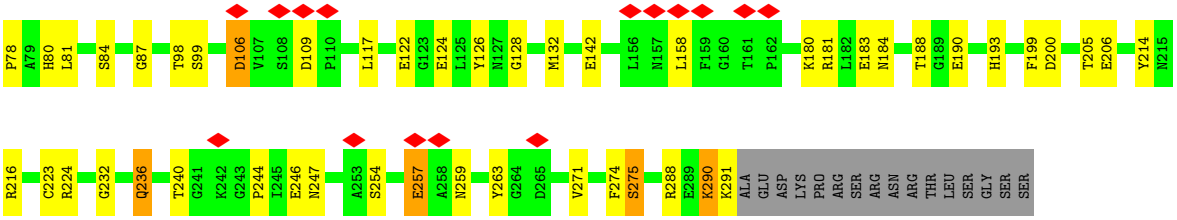


• Molecule 1: Far-red light-harvesting complex

Chain C:

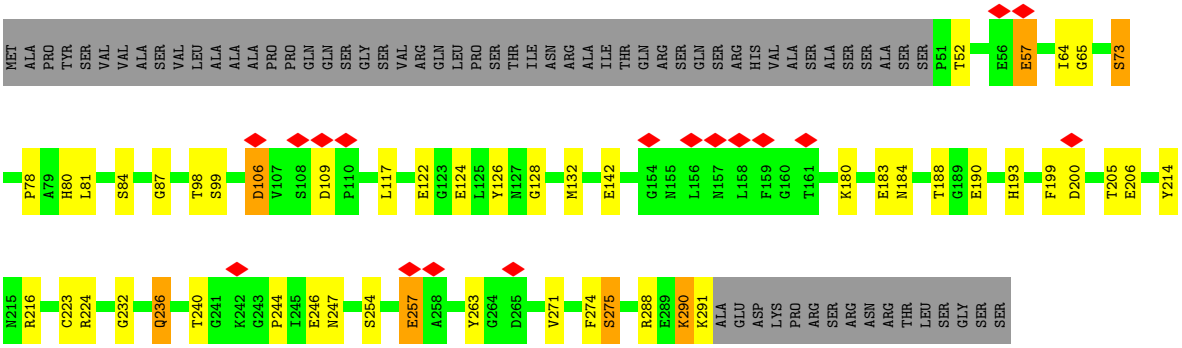






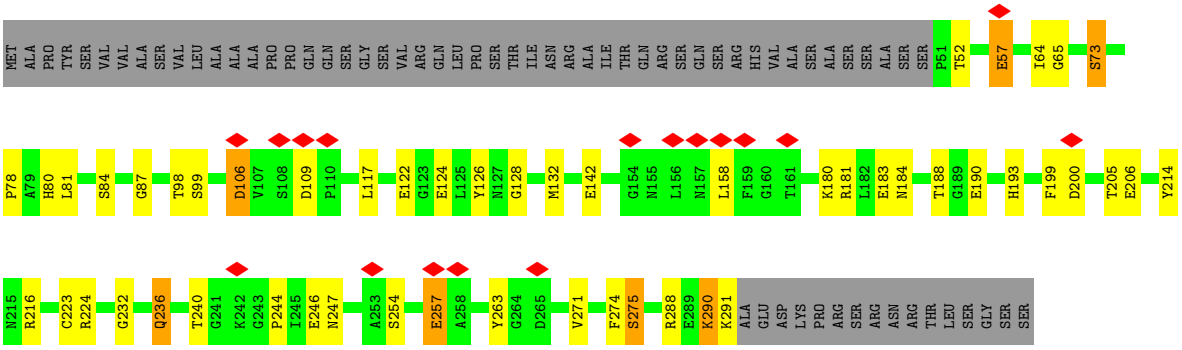
• Molecule 1: Far-red light-harvesting complex





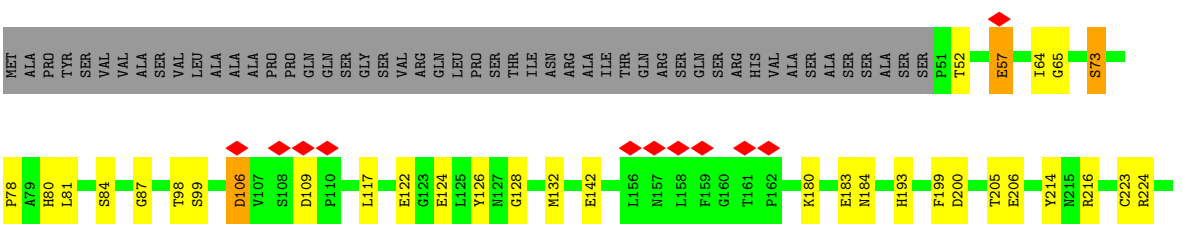
• Molecule 1: Far-red light-harvesting complex

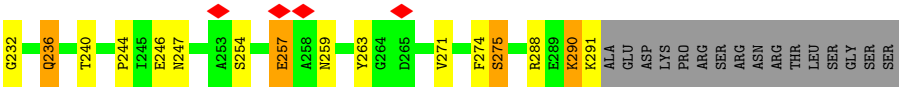




• Molecule 1: Far-red light-harvesting complex



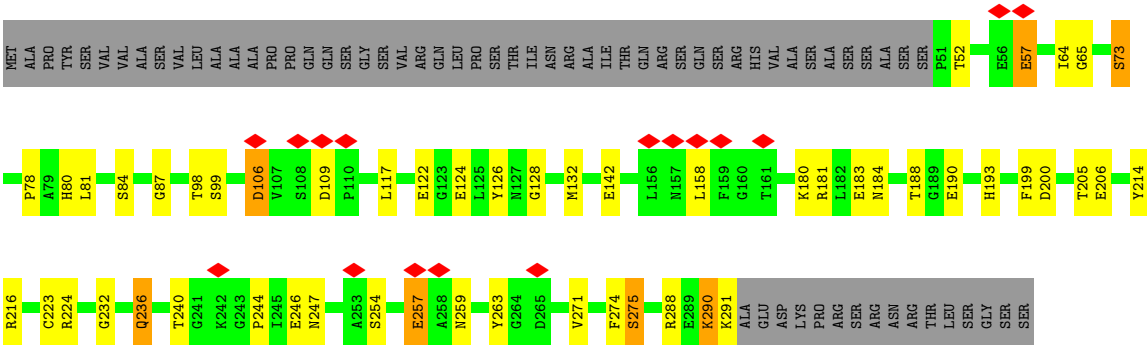




• Molecule 1: Far-red light-harvesting complex

Chain G:

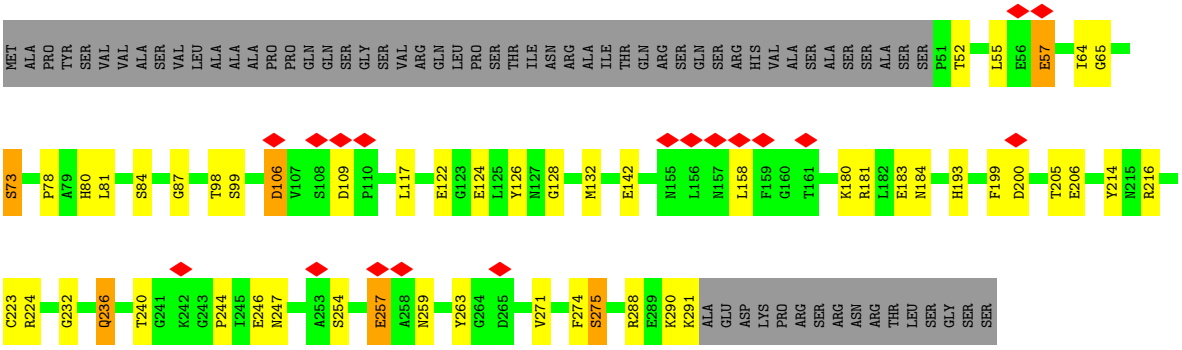




• Molecule 1: Far-red light-harvesting complex

Chain H:

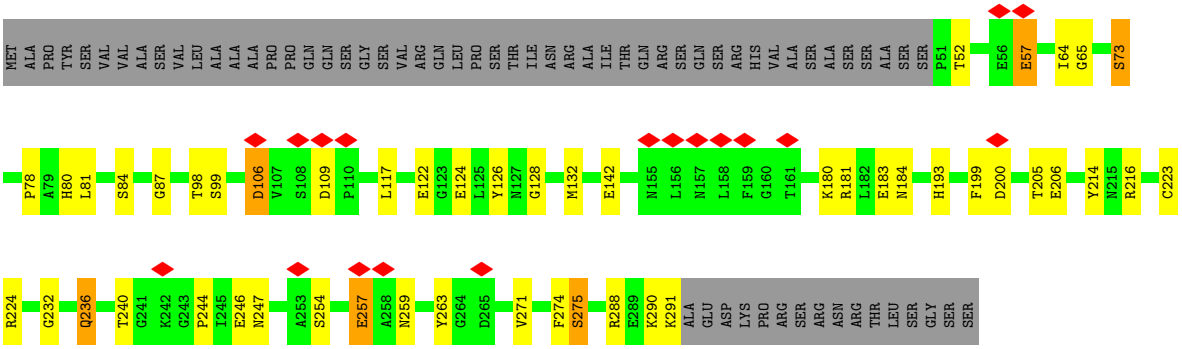




• Molecule 1: Far-red light-harvesting complex

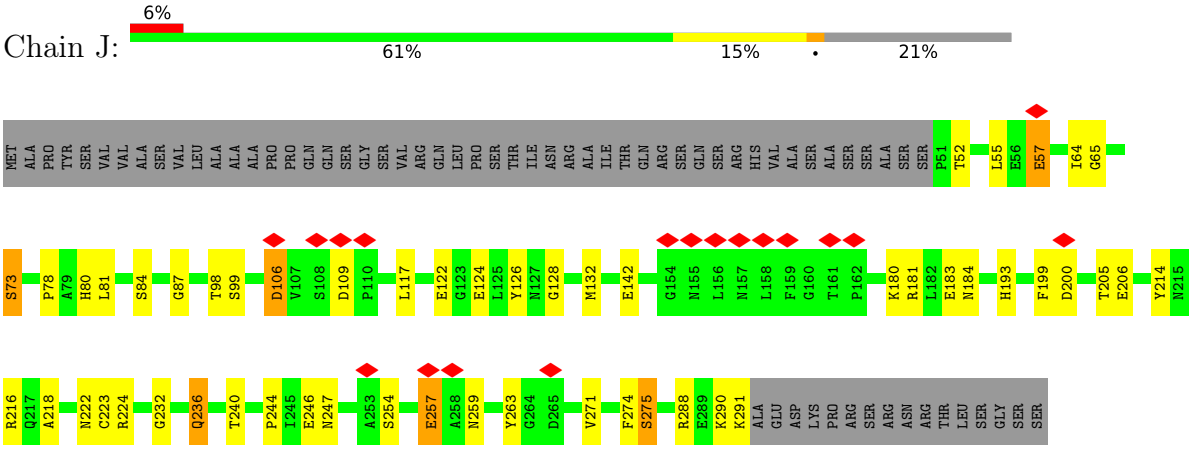
Chain I:





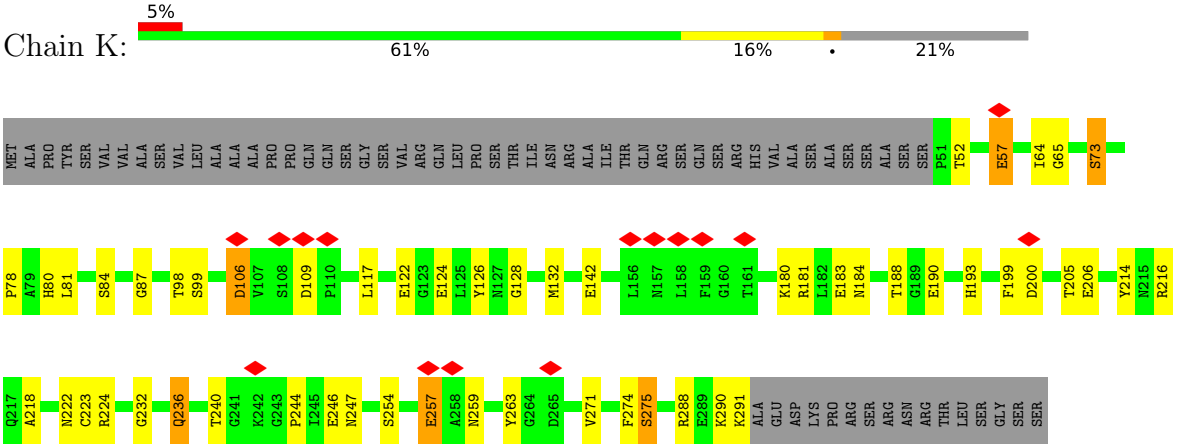
• Molecule 1: Far-red light-harvesting complex

Chain J:



• Molecule 1: Far-red light-harvesting complex

Chain K:



4 Experimental information (i)

Property	Value	Source
EM reconstruction method	SINGLE PARTICLE	Depositor
Imposed symmetry	POINT, C11	Depositor
Number of particles used	99510	Depositor
Resolution determination method	FSC 0.143 CUT-OFF	Depositor
CTF correction method	PHASE FLIPPING AND AMPLITUDE CORRECTION	Depositor
Microscope	TFS TALOS	Depositor
Voltage (kV)	200	Depositor
Electron dose ($e^-/Å^2$)	50	Depositor
Minimum defocus (nm)	1000	Depositor
Maximum defocus (nm)	3000	Depositor
Magnification	92000	Depositor
Image detector	FEI FALCON III (4k x 4k)	Depositor
Maximum map value	0.415	Depositor
Minimum map value	-0.229	Depositor
Average map value	0.000	Depositor
Map value standard deviation	0.015	Depositor
Recommended contour level	0.06	Depositor
Map size (Å)	282.5, 282.5, 282.5	wwPDB
Map dimensions	250, 250, 250	wwPDB
Map angles (°)	90.0, 90.0, 90.0	wwPDB
Pixel spacing (Å)	1.13, 1.13, 1.13	Depositor

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: XAT, 32N, CLA

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	0.39	0/1944	0.60	1/2650 (0.0%)
1	B	0.39	0/1944	0.60	1/2650 (0.0%)
1	C	0.38	0/1944	0.60	1/2650 (0.0%)
1	D	0.39	0/1944	0.60	1/2650 (0.0%)
1	E	0.39	0/1944	0.60	1/2650 (0.0%)
1	F	0.39	0/1944	0.60	1/2650 (0.0%)
1	G	0.39	0/1944	0.60	1/2650 (0.0%)
1	H	0.38	0/1944	0.60	1/2650 (0.0%)
1	I	0.39	0/1944	0.60	1/2650 (0.0%)
1	J	0.38	0/1944	0.60	1/2650 (0.0%)
1	K	0.39	0/1944	0.60	1/2650 (0.0%)
All	All	0.39	0/21384	0.60	11/29150 (0.0%)

There are no bond length outliers.

The worst 5 of 11 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	G	199	PHE	N-CA-C	5.78	118.24	111.02
1	C	199	PHE	N-CA-C	5.76	118.22	111.02
1	E	199	PHE	N-CA-C	5.76	118.22	111.02
1	A	199	PHE	N-CA-C	5.76	118.22	111.02
1	F	199	PHE	N-CA-C	5.76	118.22	111.02

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	1886	0	1811	47	0
1	B	1886	0	1811	48	0
1	C	1886	0	1811	49	0
1	D	1886	0	1811	46	0
1	E	1886	0	1811	48	0
1	F	1886	0	1811	47	0
1	G	1886	0	1811	50	0
1	H	1886	0	1811	49	0
1	I	1886	0	1811	47	0
1	J	1886	0	1811	49	0
1	K	1886	0	1811	48	0
2	A	688	0	729	43	0
2	B	688	0	729	40	0
2	C	688	0	729	40	0
2	D	688	0	729	40	0
2	E	688	0	729	40	0
2	F	688	0	729	40	0
2	G	688	0	729	41	0
2	H	688	0	729	40	0
2	I	688	0	729	43	0
2	J	688	0	729	43	0
2	K	688	0	729	39	0
3	A	44	0	56	4	0
3	B	44	0	56	4	0
3	C	44	0	56	4	0
3	D	44	0	56	4	0
3	E	44	0	56	4	0
3	F	44	0	56	4	0
3	G	44	0	56	4	0
3	H	44	0	56	4	0
3	I	44	0	56	4	0
3	J	44	0	56	4	0
3	K	44	0	56	4	0
4	A	43	0	0	0	0
4	B	43	0	0	0	0
4	C	43	0	0	0	0
4	D	43	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	E	43	0	0	0	0
4	F	43	0	0	0	0
4	G	43	0	0	0	0
4	H	43	0	0	0	0
4	I	43	0	0	0	0
4	J	43	0	0	0	0
4	K	43	0	0	0	0
All	All	29271	0	28556	837	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 14.

The worst 5 of 837 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
2:A:601:CLA:H141	2:A:601:CLA:H171	1.15	1.14
2:B:601:CLA:H141	2:B:601:CLA:H171	1.15	1.14
2:K:601:CLA:H141	2:K:601:CLA:H171	1.15	1.13
2:C:601:CLA:H141	2:C:601:CLA:H171	1.15	1.13
2:J:601:CLA:H141	2:J:601:CLA:H171	1.15	1.11

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	239/307 (78%)	212 (89%)	27 (11%)	0	100	100
1	B	239/307 (78%)	212 (89%)	27 (11%)	0	100	100
1	C	239/307 (78%)	212 (89%)	27 (11%)	0	100	100
1	D	239/307 (78%)	212 (89%)	27 (11%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	E	239/307 (78%)	212 (89%)	27 (11%)	0	100	100
1	F	239/307 (78%)	212 (89%)	27 (11%)	0	100	100
1	G	239/307 (78%)	212 (89%)	27 (11%)	0	100	100
1	H	239/307 (78%)	212 (89%)	27 (11%)	0	100	100
1	I	239/307 (78%)	212 (89%)	27 (11%)	0	100	100
1	J	239/307 (78%)	212 (89%)	27 (11%)	0	100	100
1	K	239/307 (78%)	212 (89%)	27 (11%)	0	100	100
All	All	2629/3377 (78%)	2332 (89%)	297 (11%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	193/247 (78%)	180 (93%)	13 (7%)	15	39
1	B	193/247 (78%)	180 (93%)	13 (7%)	15	39
1	C	193/247 (78%)	180 (93%)	13 (7%)	15	39
1	D	193/247 (78%)	180 (93%)	13 (7%)	15	39
1	E	193/247 (78%)	180 (93%)	13 (7%)	15	39
1	F	193/247 (78%)	180 (93%)	13 (7%)	15	39
1	G	193/247 (78%)	180 (93%)	13 (7%)	15	39
1	H	193/247 (78%)	180 (93%)	13 (7%)	15	39
1	I	193/247 (78%)	180 (93%)	13 (7%)	15	39
1	J	193/247 (78%)	180 (93%)	13 (7%)	15	39
1	K	193/247 (78%)	180 (93%)	13 (7%)	15	39
All	All	2123/2717 (78%)	1980 (93%)	143 (7%)	17	39

5 of 143 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	I	275	SER
1	J	84	SER
1	K	57	GLU
1	D	236	GLN
1	D	200	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

143 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
2	CLA	E	603	1	60,64,73	1.77	9 (15%)	72,102,113	1.40	9 (12%)
2	CLA	D	606	-	69,73,73	1.67	10 (14%)	83,113,113	1.35	12 (14%)
2	CLA	D	611	1	60,64,73	1.69	9 (15%)	72,102,113	1.31	9 (12%)
2	CLA	F	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)
2	CLA	E	605	1	69,73,73	1.61	9 (13%)	83,113,113	1.33	10 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	CLA	H	609	1	69,73,73	1.66	9 (13%)	83,113,113	1.25	10 (12%)
2	CLA	F	605	1	69,73,73	1.60	9 (13%)	83,113,113	1.34	10 (12%)
4	32N	C	613	-	42,44,44	0.27	0	50,61,61	0.60	1 (2%)
2	CLA	G	608	-	60,64,73	1.84	10 (16%)	72,102,113	1.43	12 (16%)
2	CLA	K	605	1	69,73,73	1.61	9 (13%)	83,113,113	1.33	10 (12%)
2	CLA	J	603	1	60,64,73	1.77	9 (15%)	72,102,113	1.40	9 (12%)
3	XAT	D	612	-	39,47,47	0.21	0	54,74,74	1.00	4 (7%)
2	CLA	I	605	1	69,73,73	1.61	9 (13%)	83,113,113	1.33	10 (12%)
2	CLA	A	608	-	60,64,73	1.83	10 (16%)	72,102,113	1.43	12 (16%)
2	CLA	F	601	-	69,73,73	1.62	10 (14%)	83,113,113	1.18	8 (9%)
2	CLA	A	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)
2	CLA	K	604	-	69,73,73	1.62	9 (13%)	83,113,113	1.17	8 (9%)
2	CLA	I	604	-	69,73,73	1.63	9 (13%)	83,113,113	1.17	8 (9%)
2	CLA	D	610	-	69,73,73	1.66	10 (14%)	83,113,113	1.38	10 (12%)
2	CLA	K	601	-	69,73,73	1.62	10 (14%)	83,113,113	1.17	9 (10%)
2	CLA	B	608	-	60,64,73	1.83	10 (16%)	72,102,113	1.43	12 (16%)
2	CLA	C	605	1	69,73,73	1.60	9 (13%)	83,113,113	1.33	10 (12%)
2	CLA	B	601	-	69,73,73	1.62	10 (14%)	83,113,113	1.18	8 (9%)
2	CLA	K	611	1	60,64,73	1.69	9 (15%)	72,102,113	1.31	9 (12%)
2	CLA	D	605	1	69,73,73	1.61	9 (13%)	83,113,113	1.33	10 (12%)
2	CLA	B	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)
2	CLA	J	604	-	69,73,73	1.63	9 (13%)	83,113,113	1.16	8 (9%)
3	XAT	E	612	-	39,47,47	0.21	0	54,74,74	1.00	4 (7%)
3	XAT	C	612	-	39,47,47	0.21	0	54,74,74	1.00	4 (7%)
2	CLA	B	610	-	69,73,73	1.66	10 (14%)	83,113,113	1.39	10 (12%)
2	CLA	C	603	1	60,64,73	1.77	9 (15%)	72,102,113	1.40	9 (12%)
2	CLA	E	601	-	69,73,73	1.62	10 (14%)	83,113,113	1.18	9 (10%)
2	CLA	C	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
2	CLA	E	611	1	60,64,73	1.69	9 (15%)	72,102,113	1.31	9 (12%)
2	CLA	H	603	1	60,64,73	1.77	9 (15%)	72,102,113	1.40	9 (12%)
2	CLA	I	610	-	69,73,73	1.66	10 (14%)	83,113,113	1.39	10 (12%)
2	CLA	F	609	1	69,73,73	1.66	9 (13%)	83,113,113	1.25	10 (12%)
2	CLA	D	603	1	60,64,73	1.77	9 (15%)	72,102,113	1.40	9 (12%)
4	32N	H	613	-	42,44,44	0.28	0	50,61,61	0.60	1 (2%)
4	32N	J	613	-	42,44,44	0.27	0	50,61,61	0.60	1 (2%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	CLA	G	606	-	69,73,73	1.67	10 (14%)	83,113,113	1.36	12 (14%)
2	CLA	J	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)
4	32N	F	613	-	42,44,44	0.27	0	50,61,61	0.60	1 (2%)
2	CLA	C	608	-	60,64,73	1.83	10 (16%)	72,102,113	1.43	12 (16%)
4	32N	A	613	-	42,44,44	0.27	0	50,61,61	0.60	1 (2%)
2	CLA	G	610	-	69,73,73	1.66	10 (14%)	83,113,113	1.38	10 (12%)
2	CLA	E	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
2	CLA	G	609	1	69,73,73	1.66	9 (13%)	83,113,113	1.25	10 (12%)
2	CLA	C	611	1	60,64,73	1.69	9 (15%)	72,102,113	1.31	9 (12%)
2	CLA	B	609	1	69,73,73	1.66	9 (13%)	83,113,113	1.25	10 (12%)
2	CLA	J	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
2	CLA	K	606	-	69,73,73	1.67	10 (14%)	83,113,113	1.36	12 (14%)
2	CLA	H	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)
4	32N	B	613	-	42,44,44	0.28	0	50,61,61	0.60	1 (2%)
2	CLA	G	603	1	60,64,73	1.78	9 (15%)	72,102,113	1.40	9 (12%)
2	CLA	H	605	1	69,73,73	1.60	9 (13%)	83,113,113	1.33	10 (12%)
2	CLA	J	605	1	69,73,73	1.61	9 (13%)	83,113,113	1.33	10 (12%)
2	CLA	C	606	-	69,73,73	1.67	10 (14%)	83,113,113	1.35	12 (14%)
2	CLA	K	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
2	CLA	D	609	1	69,73,73	1.66	9 (13%)	83,113,113	1.25	10 (12%)
2	CLA	I	601	-	69,73,73	1.62	10 (14%)	83,113,113	1.18	9 (10%)
2	CLA	A	604	-	69,73,73	1.63	9 (13%)	83,113,113	1.17	8 (9%)
2	CLA	B	603	1	60,64,73	1.77	9 (15%)	72,102,113	1.40	9 (12%)
2	CLA	H	601	-	69,73,73	1.62	10 (14%)	83,113,113	1.17	9 (10%)
2	CLA	F	604	-	69,73,73	1.63	9 (13%)	83,113,113	1.16	8 (9%)
2	CLA	J	611	1	60,64,73	1.69	9 (15%)	72,102,113	1.31	9 (12%)
2	CLA	F	611	1	60,64,73	1.70	9 (15%)	72,102,113	1.31	9 (12%)
2	CLA	K	609	1	69,73,73	1.66	8 (11%)	83,113,113	1.26	10 (12%)
2	CLA	I	608	-	60,64,73	1.83	10 (16%)	72,102,113	1.43	12 (16%)
2	CLA	F	608	-	60,64,73	1.83	10 (16%)	72,102,113	1.42	12 (16%)
3	XAT	G	612	-	39,47,47	0.21	0	54,74,74	1.00	4 (7%)
2	CLA	A	605	1	69,73,73	1.60	9 (13%)	83,113,113	1.33	10 (12%)
2	CLA	E	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)
2	CLA	H	611	1	60,64,73	1.69	9 (15%)	72,102,113	1.31	9 (12%)
4	32N	G	613	-	42,44,44	0.27	0	50,61,61	0.60	1 (2%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	CLA	B	605	1	69,73,73	1.61	9 (13%)	83,113,113	1.33	10 (12%)
2	CLA	C	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)
2	CLA	G	611	1	60,64,73	1.69	9 (15%)	72,102,113	1.30	9 (12%)
2	CLA	H	606	-	69,73,73	1.67	10 (14%)	83,113,113	1.35	12 (14%)
2	CLA	G	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)
2	CLA	A	611	1	60,64,73	1.69	9 (15%)	72,102,113	1.31	9 (12%)
3	XAT	J	612	-	39,47,47	0.21	0	54,74,74	1.00	5 (9%)
2	CLA	G	604	-	69,73,73	1.63	9 (13%)	83,113,113	1.17	8 (9%)
2	CLA	D	601	-	69,73,73	1.62	10 (14%)	83,113,113	1.18	9 (10%)
2	CLA	B	611	1	60,64,73	1.69	9 (15%)	72,102,113	1.31	9 (12%)
3	XAT	F	612	-	39,47,47	0.22	0	54,74,74	1.00	4 (7%)
2	CLA	F	606	-	69,73,73	1.67	10 (14%)	83,113,113	1.35	12 (14%)
3	XAT	K	612	-	39,47,47	0.21	0	54,74,74	1.00	4 (7%)
3	XAT	I	612	-	39,47,47	0.22	0	54,74,74	0.99	4 (7%)
2	CLA	D	604	-	69,73,73	1.63	9 (13%)	83,113,113	1.17	8 (9%)
2	CLA	G	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
4	32N	K	613	-	42,44,44	0.27	0	50,61,61	0.60	1 (2%)
3	XAT	A	612	-	39,47,47	0.21	0	54,74,74	1.00	4 (7%)
2	CLA	H	608	-	60,64,73	1.83	10 (16%)	72,102,113	1.43	12 (16%)
2	CLA	I	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)
2	CLA	J	601	-	69,73,73	1.63	10 (14%)	83,113,113	1.18	8 (9%)
2	CLA	F	610	-	69,73,73	1.66	10 (14%)	83,113,113	1.38	10 (12%)
2	CLA	K	610	-	69,73,73	1.67	10 (14%)	83,113,113	1.38	10 (12%)
2	CLA	E	606	-	69,73,73	1.67	10 (14%)	83,113,113	1.35	12 (14%)
2	CLA	H	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
2	CLA	A	601	-	69,73,73	1.62	10 (14%)	83,113,113	1.18	9 (10%)
2	CLA	E	608	-	60,64,73	1.83	10 (16%)	72,102,113	1.43	12 (16%)
4	32N	D	613	-	42,44,44	0.27	0	50,61,61	0.60	1 (2%)
2	CLA	C	609	1	69,73,73	1.65	9 (13%)	83,113,113	1.25	10 (12%)
2	CLA	E	610	-	69,73,73	1.66	10 (14%)	83,113,113	1.38	10 (12%)
2	CLA	A	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
2	CLA	B	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
2	CLA	J	609	1	69,73,73	1.65	9 (13%)	83,113,113	1.24	10 (12%)
2	CLA	K	608	-	60,64,73	1.83	10 (16%)	72,102,113	1.43	12 (16%)
2	CLA	K	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	CLA	I	603	1	60,64,73	1.78	9 (15%)	72,102,113	1.40	9 (12%)
2	CLA	A	609	1	69,73,73	1.66	9 (13%)	83,113,113	1.25	10 (12%)
2	CLA	F	603	1	60,64,73	1.78	9 (15%)	72,102,113	1.40	9 (12%)
2	CLA	K	603	1	60,64,73	1.77	9 (15%)	72,102,113	1.40	9 (12%)
2	CLA	B	604	-	69,73,73	1.62	9 (13%)	83,113,113	1.17	8 (9%)
2	CLA	C	610	-	69,73,73	1.65	10 (14%)	83,113,113	1.38	10 (12%)
2	CLA	D	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
2	CLA	C	601	-	69,73,73	1.62	10 (14%)	83,113,113	1.17	9 (10%)
2	CLA	I	606	-	69,73,73	1.67	10 (14%)	83,113,113	1.36	12 (14%)
4	32N	I	613	-	42,44,44	0.27	0	50,61,61	0.60	1 (2%)
2	CLA	J	610	-	69,73,73	1.66	10 (14%)	83,113,113	1.38	10 (12%)
2	CLA	D	608	-	60,64,73	1.83	10 (16%)	72,102,113	1.43	12 (16%)
2	CLA	J	606	-	69,73,73	1.67	10 (14%)	83,113,113	1.35	12 (14%)
3	XAT	H	612	-	39,47,47	0.21	0	54,74,74	0.99	4 (7%)
2	CLA	A	603	1	60,64,73	1.77	9 (15%)	72,102,113	1.40	9 (12%)
2	CLA	E	609	1	69,73,73	1.66	9 (13%)	83,113,113	1.25	10 (12%)
2	CLA	I	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
2	CLA	E	604	-	69,73,73	1.62	9 (13%)	83,113,113	1.16	8 (9%)
2	CLA	I	611	1	60,64,73	1.69	9 (15%)	72,102,113	1.31	9 (12%)
2	CLA	A	606	-	69,73,73	1.67	10 (14%)	83,113,113	1.35	12 (14%)
2	CLA	G	605	1	69,73,73	1.60	9 (13%)	83,113,113	1.33	10 (12%)
2	CLA	B	606	-	69,73,73	1.68	10 (14%)	83,113,113	1.35	12 (14%)
2	CLA	I	609	1	69,73,73	1.65	9 (13%)	83,113,113	1.25	10 (12%)
2	CLA	J	608	-	60,64,73	1.83	10 (16%)	72,102,113	1.43	12 (16%)
4	32N	E	613	-	42,44,44	0.27	0	50,61,61	0.60	1 (2%)
2	CLA	C	604	-	69,73,73	1.63	9 (13%)	83,113,113	1.17	8 (9%)
2	CLA	D	602	1	69,73,73	1.61	9 (13%)	83,113,113	1.30	10 (12%)
2	CLA	G	601	-	69,73,73	1.62	10 (14%)	83,113,113	1.18	8 (9%)
3	XAT	B	612	-	39,47,47	0.21	0	54,74,74	0.99	4 (7%)
2	CLA	H	604	-	69,73,73	1.63	9 (13%)	83,113,113	1.17	8 (9%)
2	CLA	H	610	-	69,73,73	1.66	10 (14%)	83,113,113	1.39	10 (12%)
2	CLA	F	607	1	69,73,73	1.56	8 (11%)	83,113,113	1.20	8 (9%)
2	CLA	A	610	-	69,73,73	1.66	10 (14%)	83,113,113	1.38	10 (12%)

In the following table, the Chirals column lists the number of chiral outliers, the number of chiral

centers analysed, the number of these observed in the model and the number defined in the Chemical Component Dictionary. Similar counts are reported in the Torsion and Rings columns. '-' means no outliers of that kind were identified.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CLA	E	603	1	-	10/29/105/115	-
2	CLA	D	606	-	-	10/39/115/115	-
2	CLA	D	611	1	-	8/29/105/115	-
2	CLA	F	602	1	-	9/39/115/115	-
2	CLA	E	605	1	-	18/39/115/115	-
2	CLA	H	609	1	-	3/39/115/115	-
2	CLA	F	605	1	-	18/39/115/115	-
4	32N	C	613	-	-	3/31/69/69	0/2/2/2
2	CLA	G	608	-	-	4/29/105/115	-
2	CLA	K	605	1	-	18/39/115/115	-
2	CLA	J	603	1	-	10/29/105/115	-
3	XAT	D	612	-	-	0/31/93/93	0/4/4/4
2	CLA	I	605	1	-	18/39/115/115	-
2	CLA	A	608	-	-	4/29/105/115	-
2	CLA	F	601	-	-	13/39/115/115	-
2	CLA	A	602	1	-	9/39/115/115	-
2	CLA	K	604	-	-	7/39/115/115	-
2	CLA	I	604	-	-	7/39/115/115	-
2	CLA	D	610	-	-	13/39/115/115	-
2	CLA	K	601	-	-	13/39/115/115	-
2	CLA	B	608	-	-	4/29/105/115	-
2	CLA	C	605	1	-	18/39/115/115	-
2	CLA	B	601	-	-	13/39/115/115	-
2	CLA	K	611	1	-	8/29/105/115	-
2	CLA	D	605	1	-	18/39/115/115	-
2	CLA	B	602	1	-	9/39/115/115	-
2	CLA	J	604	-	-	7/39/115/115	-
3	XAT	E	612	-	-	0/31/93/93	0/4/4/4
3	XAT	C	612	-	-	0/31/93/93	0/4/4/4
2	CLA	B	610	-	-	13/39/115/115	-
2	CLA	C	603	1	-	11/29/105/115	-
2	CLA	E	601	-	-	13/39/115/115	-
2	CLA	C	607	1	-	14/39/115/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CLA	E	611	1	-	8/29/105/115	-
2	CLA	H	603	1	-	10/29/105/115	-
2	CLA	I	610	-	-	13/39/115/115	-
2	CLA	F	609	1	-	3/39/115/115	-
2	CLA	D	603	1	-	10/29/105/115	-
4	32N	H	613	-	-	3/31/69/69	0/2/2/2
4	32N	J	613	-	-	3/31/69/69	0/2/2/2
2	CLA	G	606	-	-	10/39/115/115	-
2	CLA	J	602	1	-	9/39/115/115	-
4	32N	F	613	-	-	3/31/69/69	0/2/2/2
2	CLA	C	608	-	-	4/29/105/115	-
4	32N	A	613	-	-	3/31/69/69	0/2/2/2
2	CLA	G	610	-	-	13/39/115/115	-
2	CLA	E	607	1	-	14/39/115/115	-
2	CLA	G	609	1	-	3/39/115/115	-
2	CLA	C	611	1	-	8/29/105/115	-
2	CLA	B	609	1	-	3/39/115/115	-
2	CLA	J	607	1	-	14/39/115/115	-
2	CLA	K	606	-	-	10/39/115/115	-
2	CLA	H	602	1	-	9/39/115/115	-
4	32N	B	613	-	-	3/31/69/69	0/2/2/2
2	CLA	G	603	1	-	10/29/105/115	-
2	CLA	H	605	1	-	18/39/115/115	-
2	CLA	J	605	1	-	18/39/115/115	-
2	CLA	C	606	-	-	10/39/115/115	-
2	CLA	K	607	1	-	14/39/115/115	-
2	CLA	D	609	1	-	3/39/115/115	-
2	CLA	I	601	-	-	13/39/115/115	-
2	CLA	A	604	-	-	6/39/115/115	-
2	CLA	B	603	1	-	10/29/105/115	-
2	CLA	H	601	-	-	13/39/115/115	-
2	CLA	F	604	-	-	6/39/115/115	-
2	CLA	J	611	1	-	8/29/105/115	-
2	CLA	F	611	1	-	8/29/105/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CLA	K	609	1	-	3/39/115/115	-
2	CLA	I	608	-	-	4/29/105/115	-
2	CLA	F	608	-	-	4/29/105/115	-
3	XAT	G	612	-	-	0/31/93/93	0/4/4/4
2	CLA	A	605	1	-	18/39/115/115	-
2	CLA	E	602	1	-	9/39/115/115	-
2	CLA	H	611	1	-	8/29/105/115	-
4	32N	G	613	-	-	3/31/69/69	0/2/2/2
2	CLA	B	605	1	-	18/39/115/115	-
2	CLA	C	602	1	-	9/39/115/115	-
2	CLA	G	611	1	-	8/29/105/115	-
2	CLA	H	606	-	-	10/39/115/115	-
2	CLA	G	602	1	-	9/39/115/115	-
2	CLA	A	611	1	-	8/29/105/115	-
3	XAT	J	612	-	-	0/31/93/93	0/4/4/4
2	CLA	G	604	-	-	6/39/115/115	-
2	CLA	D	601	-	-	13/39/115/115	-
2	CLA	B	611	1	-	8/29/105/115	-
3	XAT	F	612	-	-	0/31/93/93	0/4/4/4
2	CLA	F	606	-	-	10/39/115/115	-
3	XAT	K	612	-	-	0/31/93/93	0/4/4/4
3	XAT	I	612	-	-	0/31/93/93	0/4/4/4
2	CLA	D	604	-	-	6/39/115/115	-
2	CLA	G	607	1	-	14/39/115/115	-
4	32N	K	613	-	-	3/31/69/69	0/2/2/2
3	XAT	A	612	-	-	0/31/93/93	0/4/4/4
2	CLA	H	608	-	-	4/29/105/115	-
2	CLA	I	602	1	-	9/39/115/115	-
2	CLA	J	601	-	-	13/39/115/115	-
2	CLA	F	610	-	-	13/39/115/115	-
2	CLA	K	610	-	-	13/39/115/115	-
2	CLA	E	606	-	-	10/39/115/115	-
2	CLA	H	607	1	-	14/39/115/115	-
2	CLA	A	601	-	-	13/39/115/115	-
2	CLA	E	608	-	-	4/29/105/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	32N	D	613	-	-	3/31/69/69	0/2/2/2
2	CLA	C	609	1	-	3/39/115/115	-
2	CLA	E	610	-	-	13/39/115/115	-
2	CLA	A	607	1	-	14/39/115/115	-
2	CLA	B	607	1	-	14/39/115/115	-
2	CLA	J	609	1	-	3/39/115/115	-
2	CLA	K	608	-	-	4/29/105/115	-
2	CLA	K	602	1	-	9/39/115/115	-
2	CLA	I	603	1	-	10/29/105/115	-
2	CLA	A	609	1	-	3/39/115/115	-
2	CLA	F	603	1	-	10/29/105/115	-
2	CLA	K	603	1	-	10/29/105/115	-
2	CLA	B	604	-	-	7/39/115/115	-
2	CLA	C	610	-	-	13/39/115/115	-
2	CLA	D	607	1	-	14/39/115/115	-
2	CLA	C	601	-	-	13/39/115/115	-
2	CLA	I	606	-	-	10/39/115/115	-
4	32N	I	613	-	-	3/31/69/69	0/2/2/2
2	CLA	J	610	-	-	13/39/115/115	-
2	CLA	D	608	-	-	4/29/105/115	-
2	CLA	J	606	-	-	10/39/115/115	-
3	XAT	H	612	-	-	0/31/93/93	0/4/4/4
2	CLA	A	603	1	-	10/29/105/115	-
2	CLA	E	609	1	-	3/39/115/115	-
2	CLA	I	607	1	-	14/39/115/115	-
2	CLA	E	604	-	-	7/39/115/115	-
2	CLA	I	611	1	-	8/29/105/115	-
2	CLA	A	606	-	-	10/39/115/115	-
2	CLA	G	605	1	-	18/39/115/115	-
2	CLA	B	606	-	-	10/39/115/115	-
2	CLA	I	609	1	-	3/39/115/115	-
2	CLA	J	608	-	-	4/29/105/115	-
4	32N	E	613	-	-	3/31/69/69	0/2/2/2
2	CLA	C	604	-	-	7/39/115/115	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	CLA	D	602	1	-	9/39/115/115	-
2	CLA	G	601	-	-	13/39/115/115	-
3	XAT	B	612	-	-	0/31/93/93	0/4/4/4
2	CLA	H	604	-	-	7/39/115/115	-
2	CLA	H	610	-	-	13/39/115/115	-
2	CLA	F	607	1	-	14/39/115/115	-
2	CLA	A	610	-	-	13/39/115/115	-

The worst 5 of 1121 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	609	CLA	C4B-NB	7.56	1.47	1.37
2	K	609	CLA	C4B-NB	7.55	1.47	1.37
2	B	609	CLA	C4B-NB	7.54	1.47	1.37
2	E	609	CLA	C4B-NB	7.54	1.47	1.37
2	G	609	CLA	C4B-NB	7.54	1.47	1.37

The worst 5 of 1229 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	K	606	CLA	C4A-NA-C1A	-6.57	103.75	106.71
2	D	606	CLA	C4A-NA-C1A	-6.54	103.77	106.71
2	G	606	CLA	C4A-NA-C1A	-6.53	103.77	106.71
2	I	606	CLA	C4A-NA-C1A	-6.52	103.77	106.71
2	F	606	CLA	C4A-NA-C1A	-6.52	103.78	106.71

There are no chirality outliers.

5 of 1229 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	602	CLA	C1A-C2A-CAA-CBA
2	A	602	CLA	C3A-C2A-CAA-CBA
2	A	605	CLA	C1A-C2A-CAA-CBA
2	A	605	CLA	C2-C3-C5-C6
2	A	605	CLA	C4-C3-C5-C6

There are no ring outliers.

113 monomers are involved in 438 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	603	CLA	4	0
2	D	606	CLA	7	0
2	D	611	CLA	1	0
2	F	602	CLA	4	0
2	E	605	CLA	3	0
2	H	609	CLA	3	0
2	F	605	CLA	2	0
2	G	608	CLA	2	0
2	K	605	CLA	3	0
2	J	603	CLA	4	0
3	D	612	XAT	4	0
2	I	605	CLA	3	0
2	A	608	CLA	2	0
2	F	601	CLA	7	0
2	A	602	CLA	4	0
2	K	601	CLA	6	0
2	B	608	CLA	2	0
2	C	605	CLA	3	0
2	B	601	CLA	7	0
2	K	611	CLA	1	0
2	D	605	CLA	2	0
2	B	602	CLA	4	0
3	E	612	XAT	4	0
3	C	612	XAT	4	0
2	C	603	CLA	3	0
2	E	601	CLA	6	0
2	C	607	CLA	12	0
2	E	611	CLA	1	0
2	H	603	CLA	3	0
2	I	610	CLA	1	0
2	F	609	CLA	3	0
2	D	603	CLA	4	0
2	G	606	CLA	7	0
2	J	602	CLA	4	0
2	C	608	CLA	2	0
2	E	607	CLA	11	0
2	G	609	CLA	3	0
2	C	611	CLA	1	0
2	B	609	CLA	3	0
2	J	607	CLA	11	0
2	K	606	CLA	6	0
2	H	602	CLA	5	0
2	G	603	CLA	3	0

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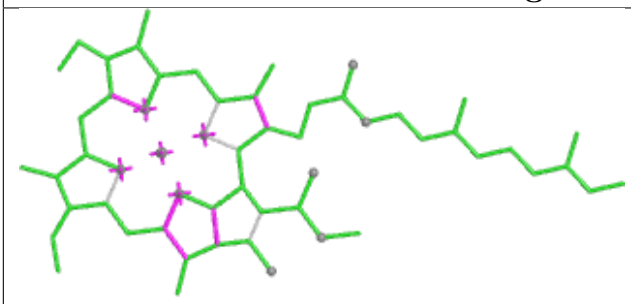
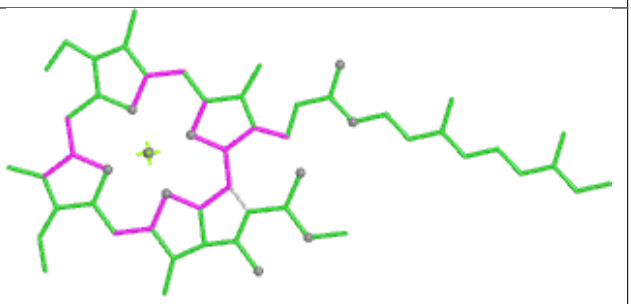
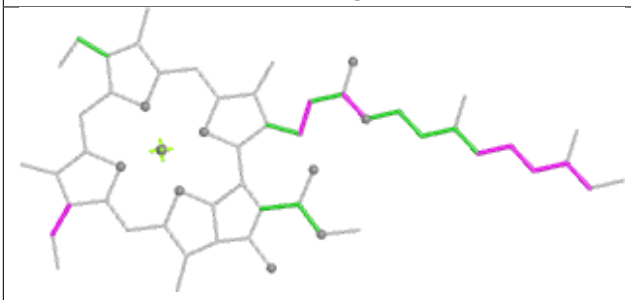
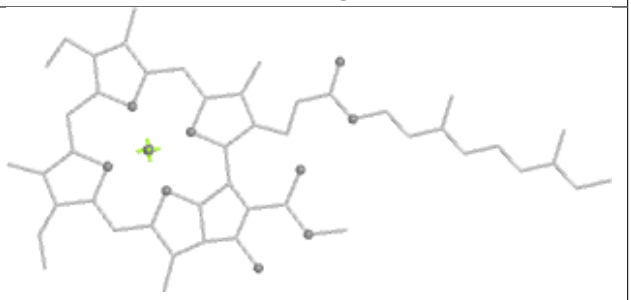
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	H	605	CLA	3	0
2	J	605	CLA	3	0
2	C	606	CLA	7	0
2	K	607	CLA	12	0
2	D	609	CLA	3	0
2	I	601	CLA	7	0
2	B	603	CLA	4	0
2	H	601	CLA	6	0
2	J	611	CLA	1	0
2	F	611	CLA	1	0
2	K	609	CLA	3	0
2	I	608	CLA	2	0
2	F	608	CLA	2	0
3	G	612	XAT	4	0
2	A	605	CLA	3	0
2	E	602	CLA	4	0
2	H	611	CLA	1	0
2	B	605	CLA	2	0
2	C	602	CLA	5	0
2	G	611	CLA	1	0
2	H	606	CLA	7	0
2	G	602	CLA	4	0
2	A	611	CLA	1	0
3	J	612	XAT	4	0
2	D	601	CLA	7	0
2	B	611	CLA	1	0
3	F	612	XAT	4	0
2	F	606	CLA	7	0
3	K	612	XAT	4	0
3	I	612	XAT	4	0
2	G	607	CLA	12	0
3	A	612	XAT	4	0
2	H	608	CLA	2	0
2	I	602	CLA	4	0
2	J	601	CLA	9	0
2	E	606	CLA	7	0
2	H	607	CLA	12	0
2	A	601	CLA	9	0
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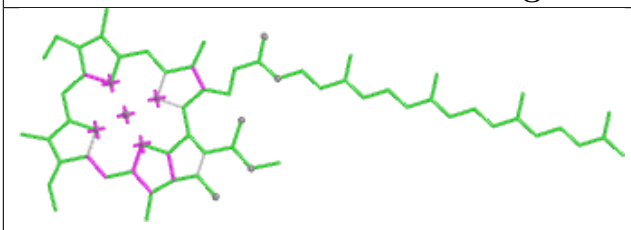
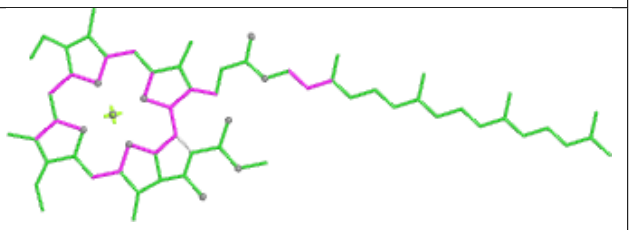
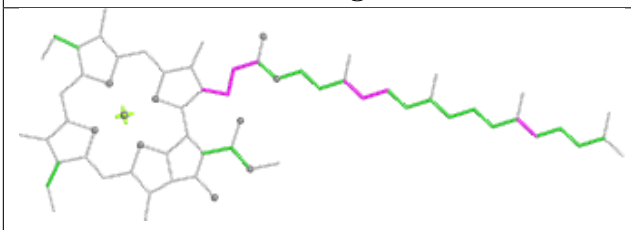
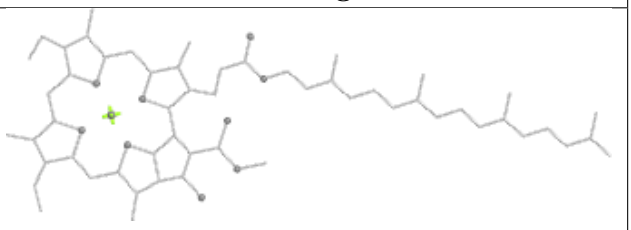
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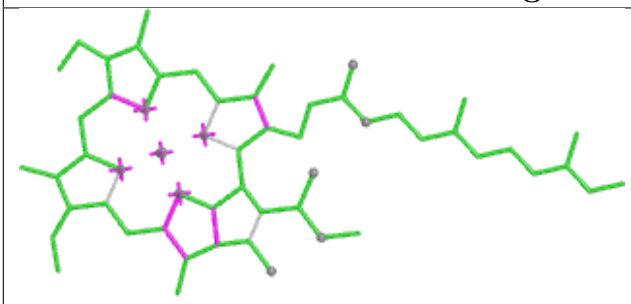
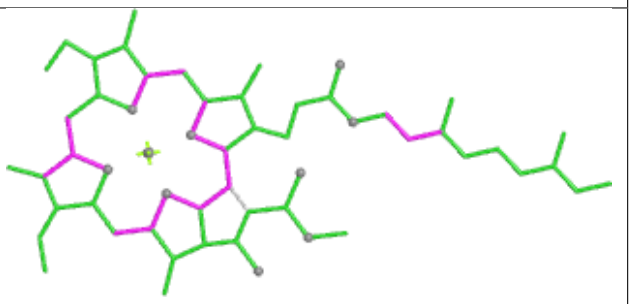
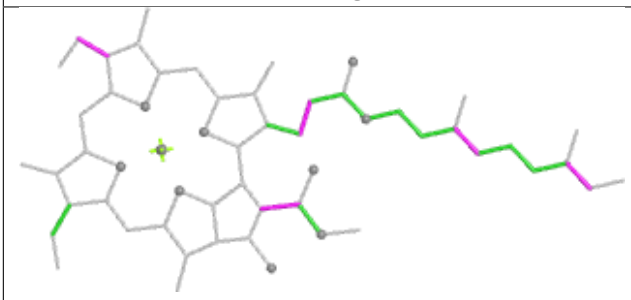
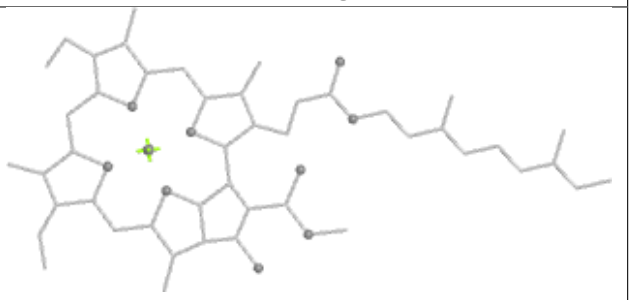
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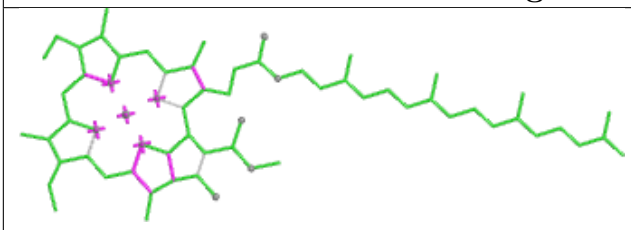
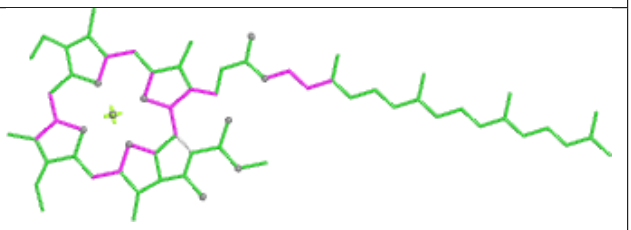
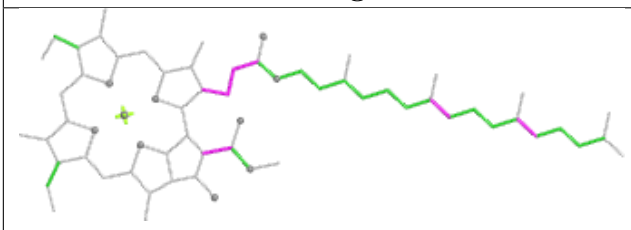
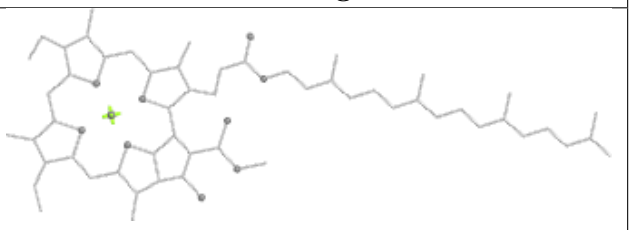
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2	K	608	CLA	2	0
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2	I	603	CLA	5	0
2	A	609	CLA	3	0
2	F	603	CLA	4	0
2	K	603	CLA	3	0
2	D	607	CLA	12	0
2	C	601	CLA	6	0
2	I	606	CLA	7	0
2	J	610	CLA	1	0
2	D	608	CLA	2	0
2	J	606	CLA	7	0
3	H	612	XAT	4	0
2	A	603	CLA	4	0
2	E	609	CLA	3	0
2	I	607	CLA	12	0
2	I	611	CLA	1	0
2	A	606	CLA	6	0
2	G	605	CLA	3	0
2	B	606	CLA	7	0
2	I	609	CLA	3	0
2	J	608	CLA	2	0
2	D	602	CLA	4	0
2	G	601	CLA	8	0
3	B	612	XAT	4	0
2	F	607	CLA	12	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 then 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.

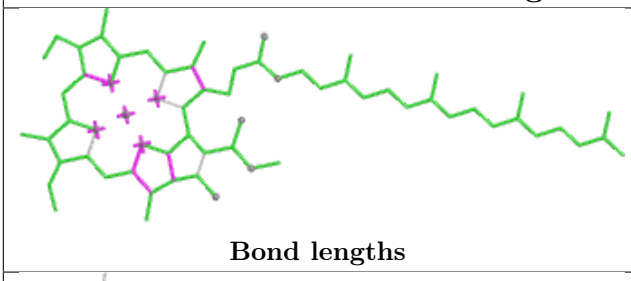
Ligand CLA E 603	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA D 606	
	
Bond lengths	Bond angles
	
Torsions	Rings

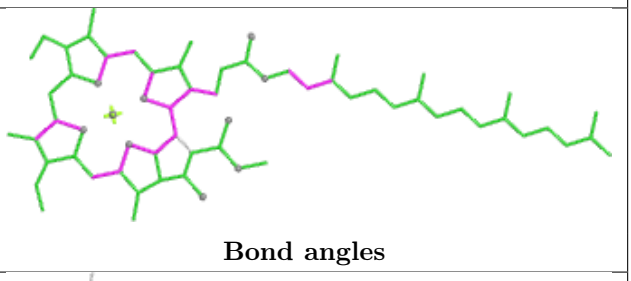
Ligand CLA D 611	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA F 602	
	
Bond lengths	Bond angles
	
Torsions	Rings

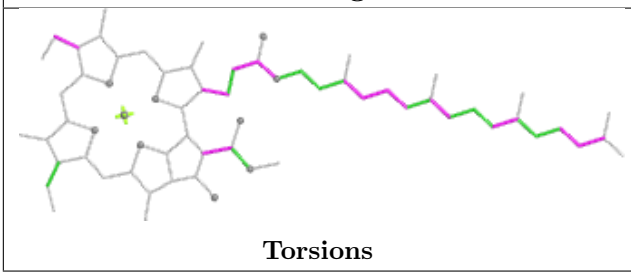
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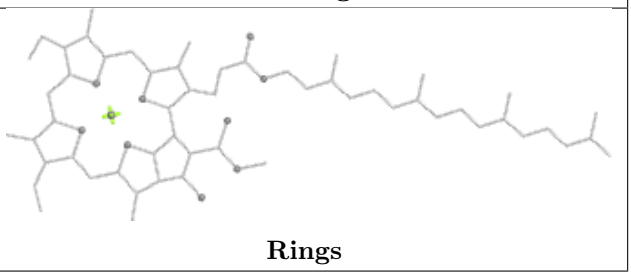
Bond lengths



Bond angles

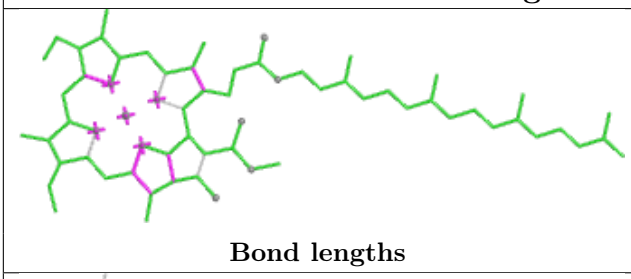


Torsions

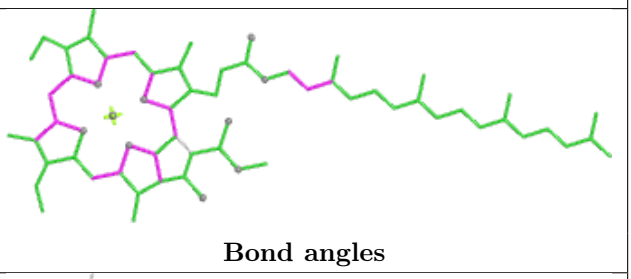


Rings

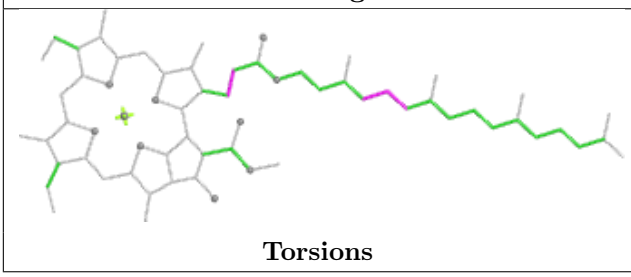
Ligand CLA H 609



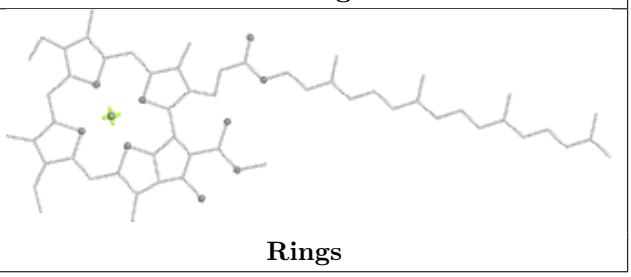
Bond lengths



Bond angles

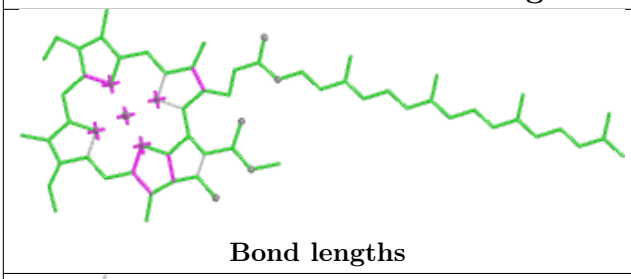


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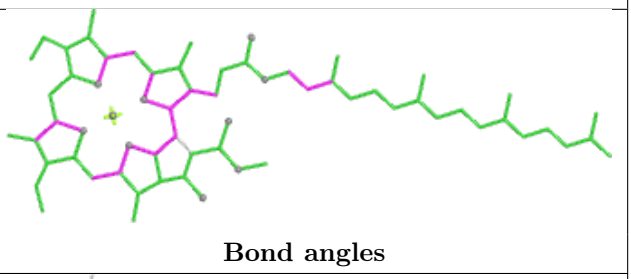


Rings

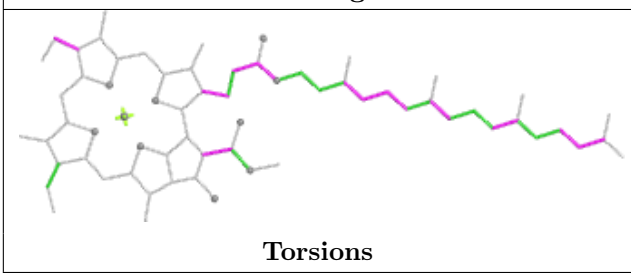
Ligand CLA F 605



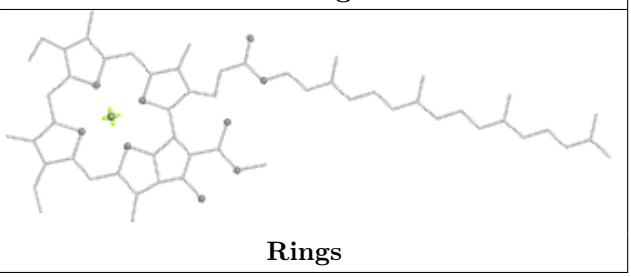
Bond lengths



Bond angles

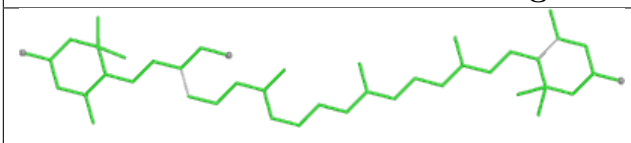


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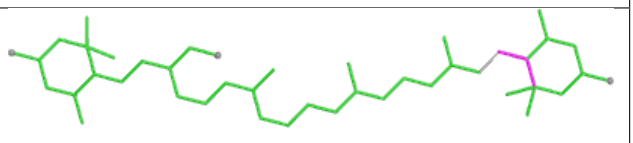


Rings

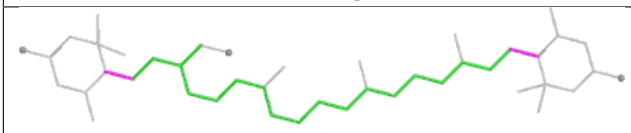
Ligand 32N C 613



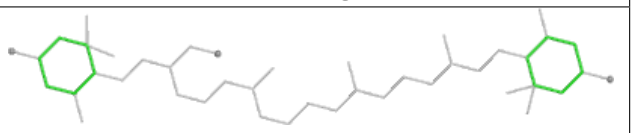
Bond lengths



Bond angles

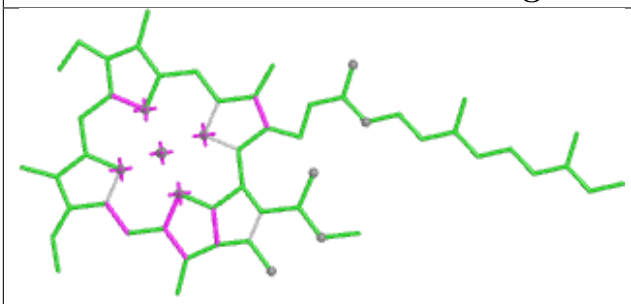


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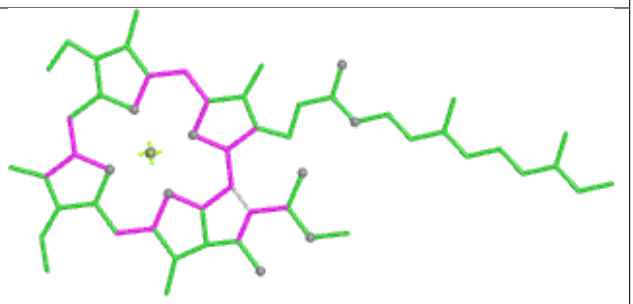


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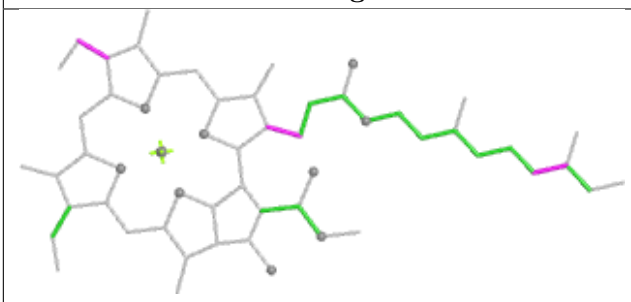
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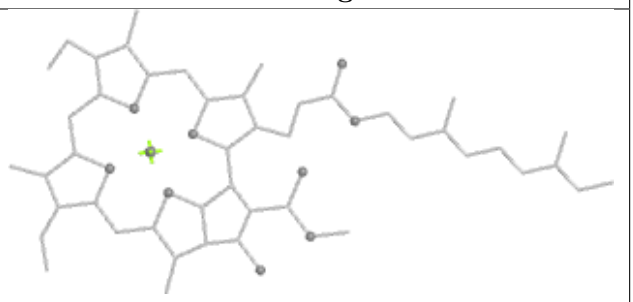
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Bond angles

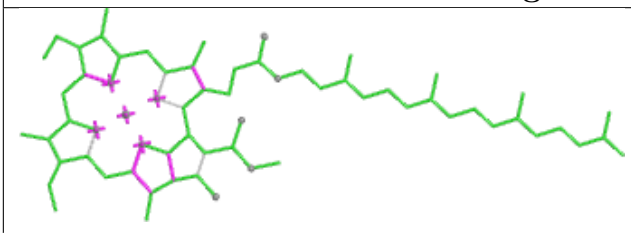


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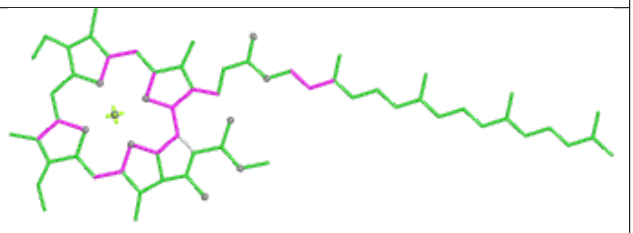


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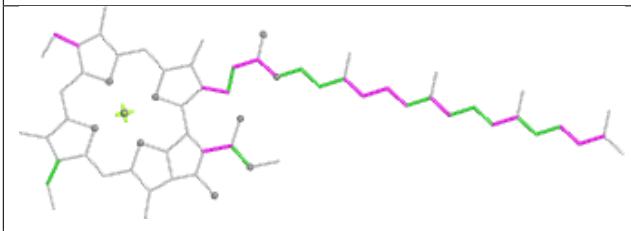
Ligand CLA K 605



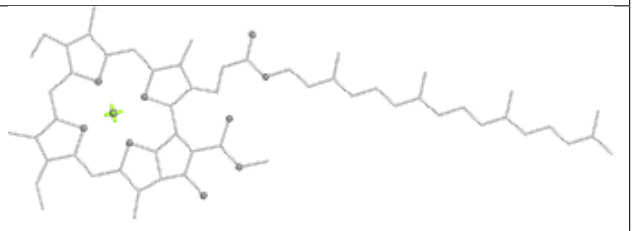
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Bond angles

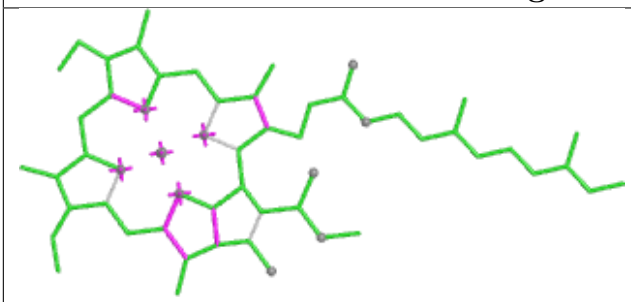


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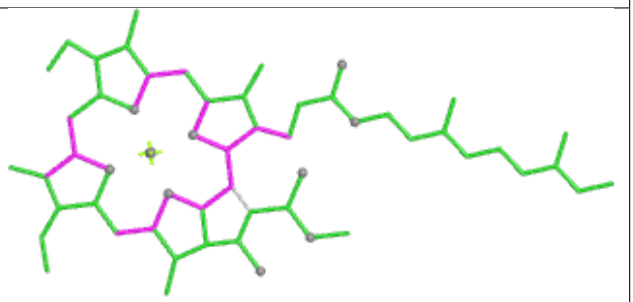


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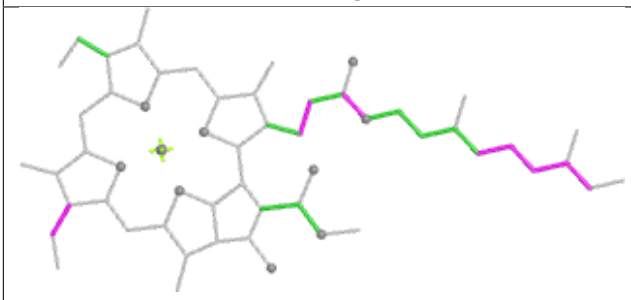
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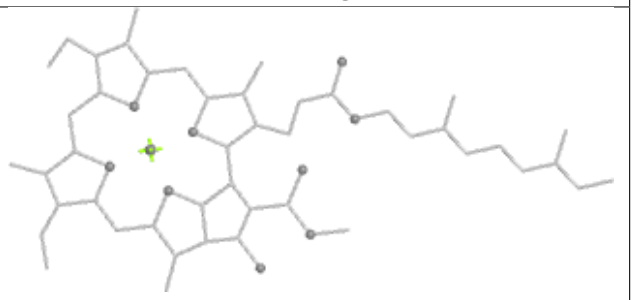
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Bond angles

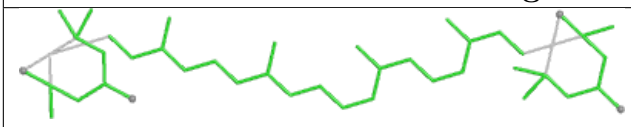


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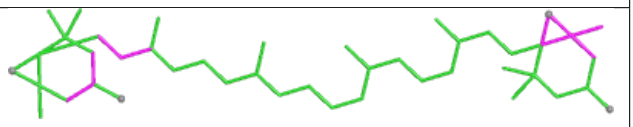


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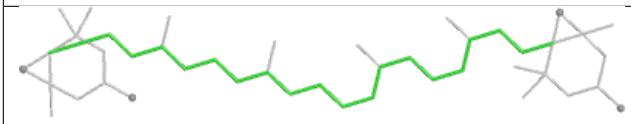
Ligand XAT D 612



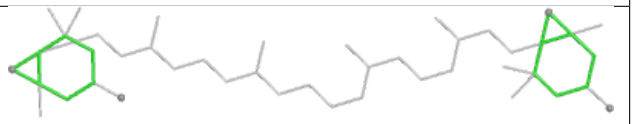
Bond lengths



Bond angles

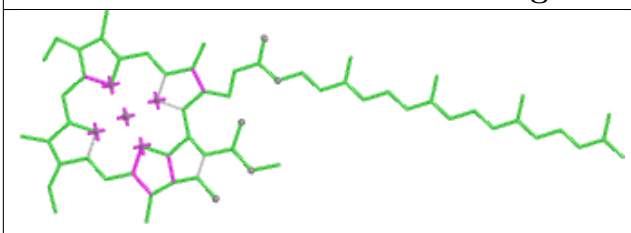


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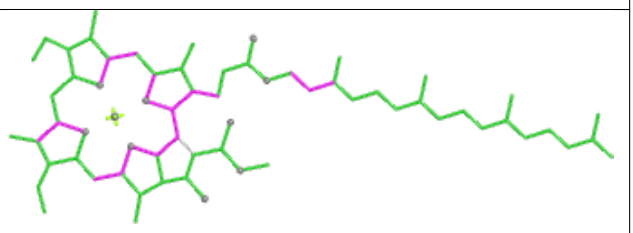


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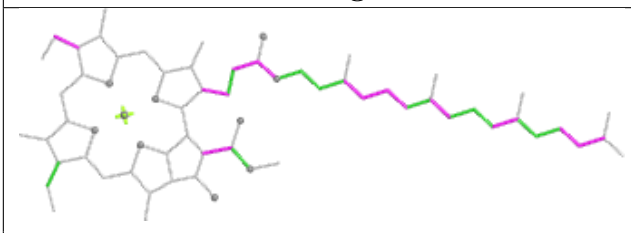
Ligand CLA I 605



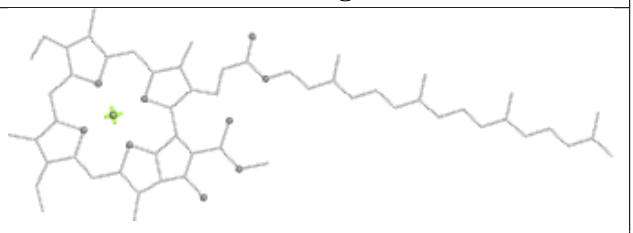
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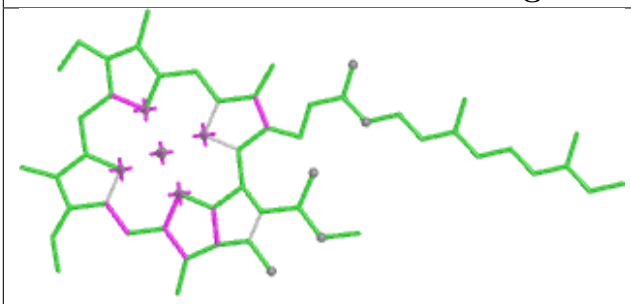
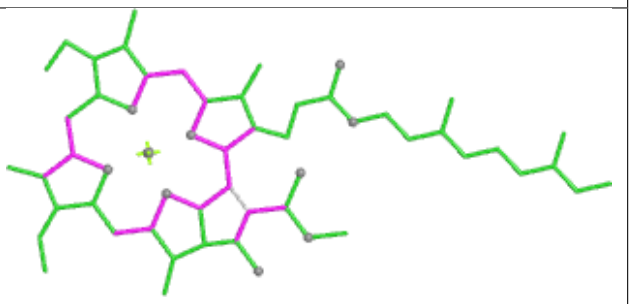
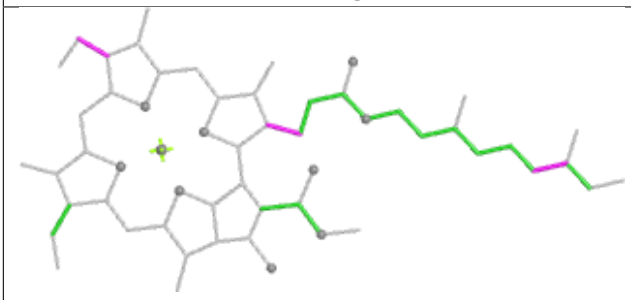
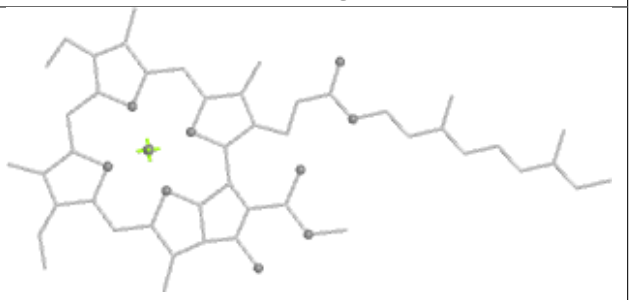
Bond angles

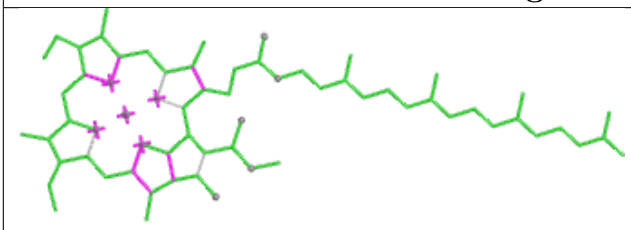
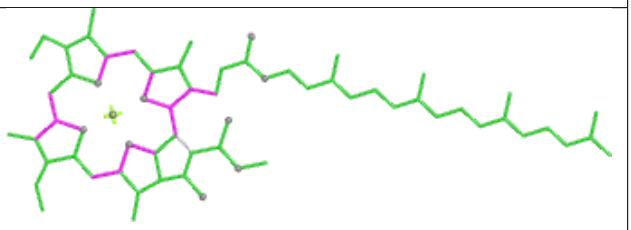
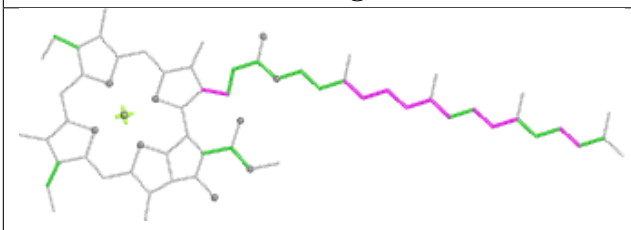
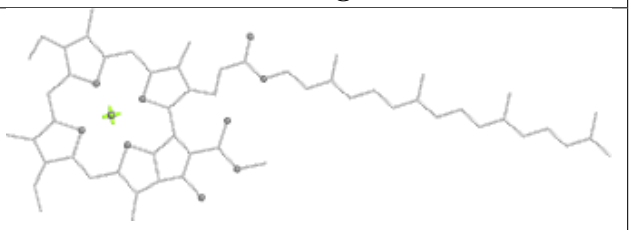


Torsions

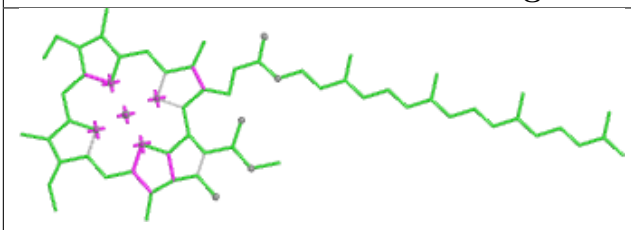


Rings

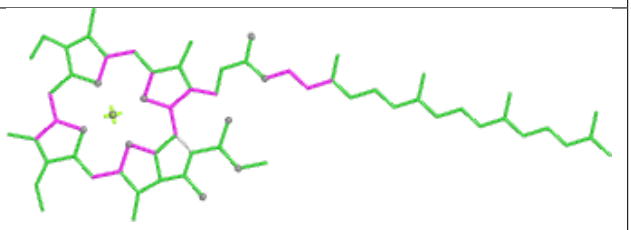
Ligand CLA A 608	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA F 601	
	
Bond lengths	Bond angles
	
Torsions	Rings

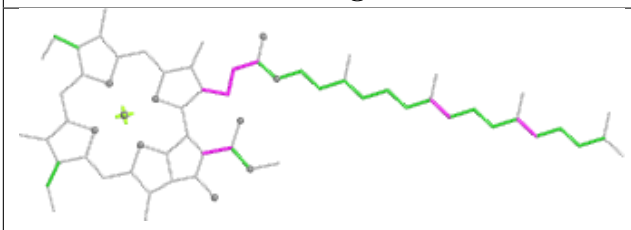
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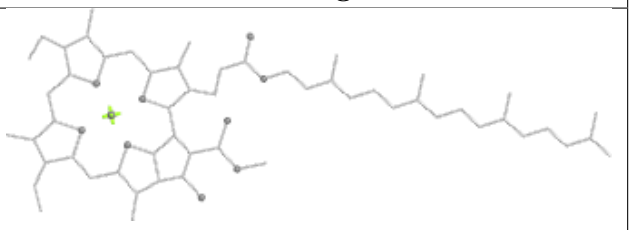
Bond lengths



Bond angles

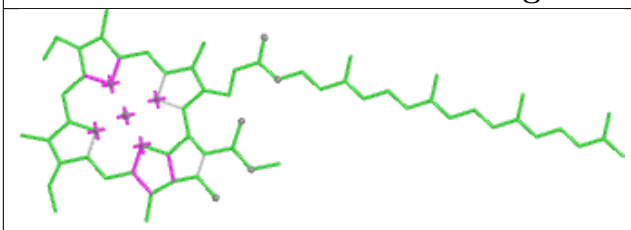


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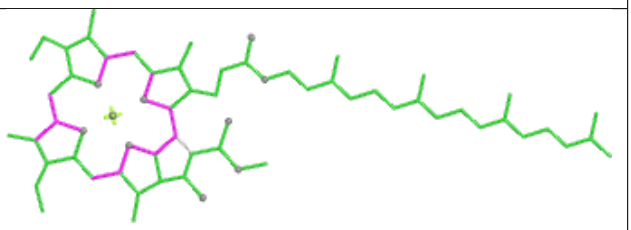


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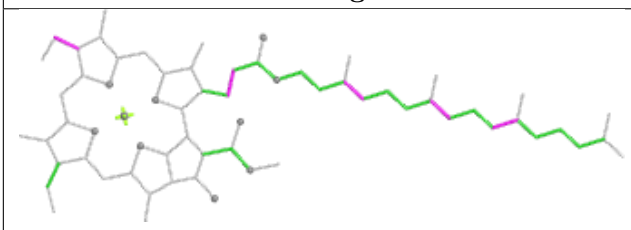
Ligand CLA K 604



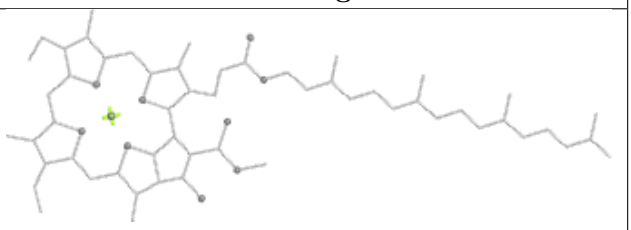
Bond lengths



Bond angles

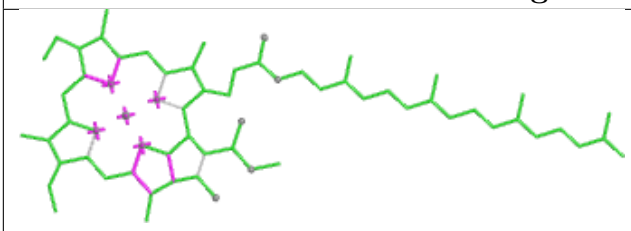


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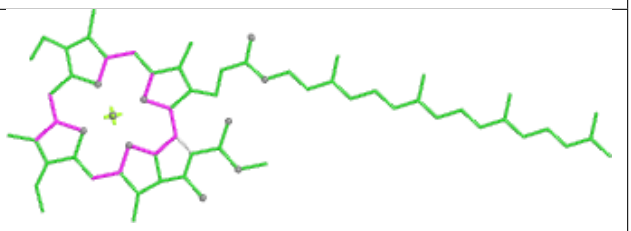


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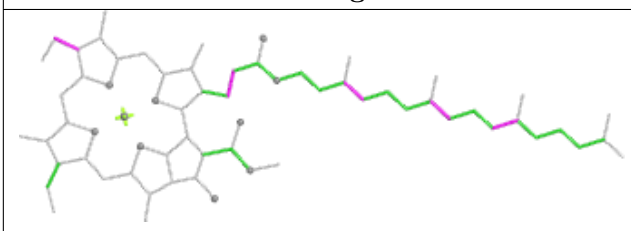
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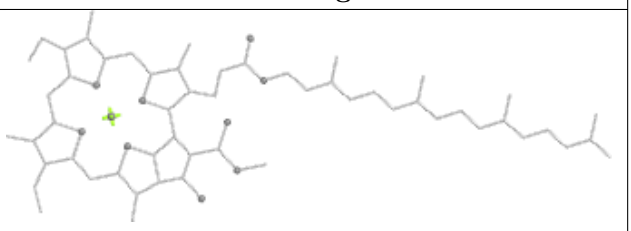
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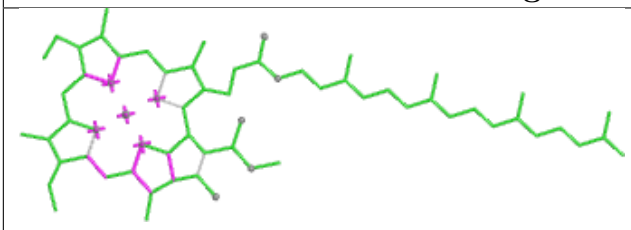
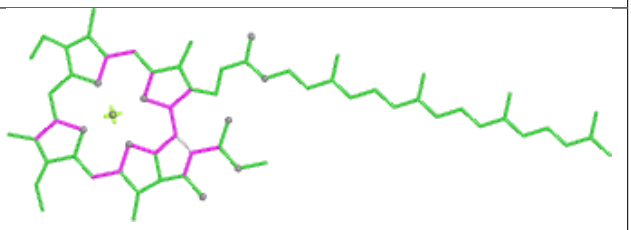
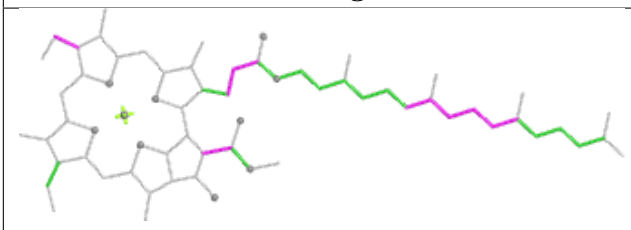
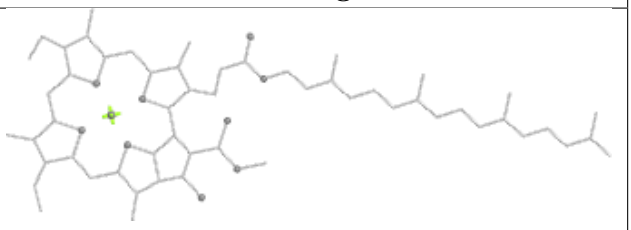
Bond angles

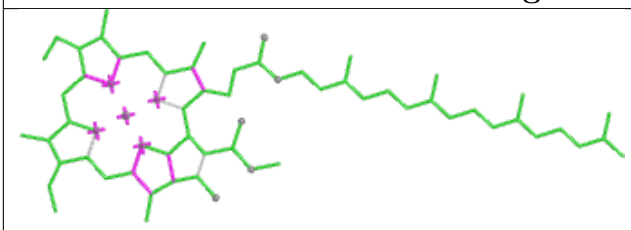
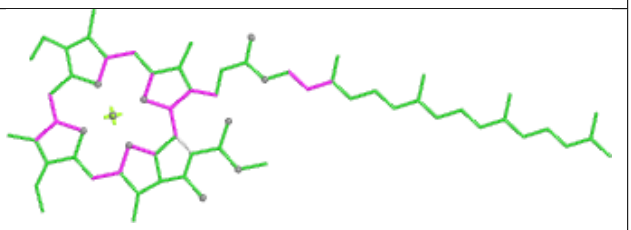
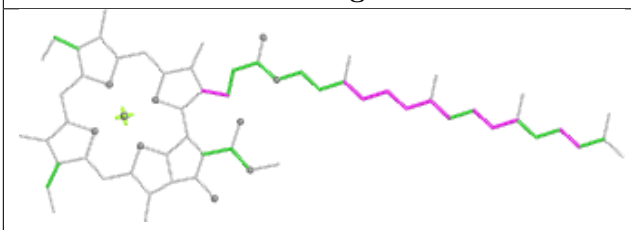
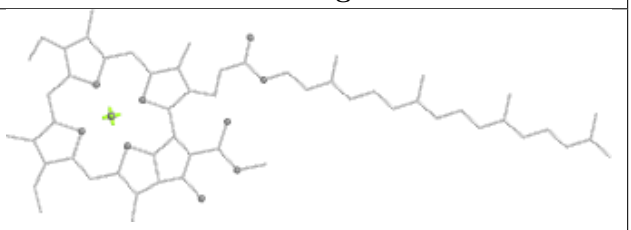


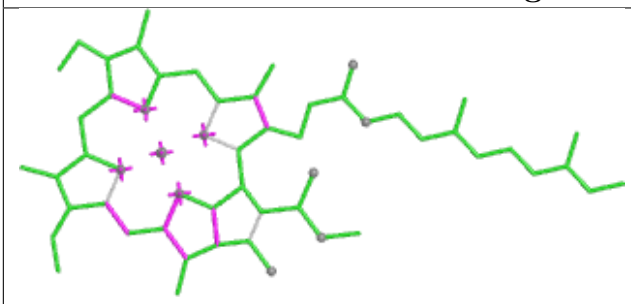
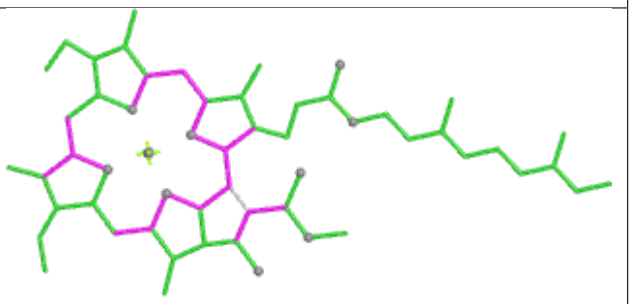
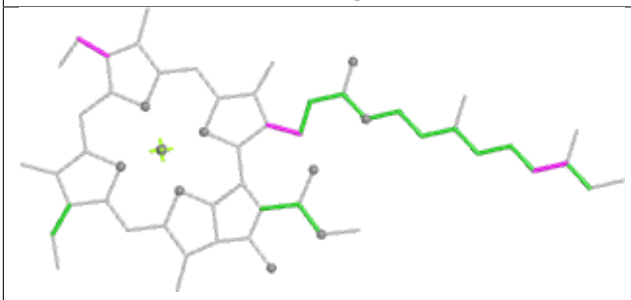
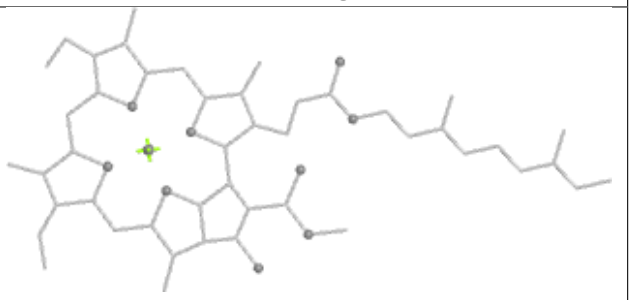
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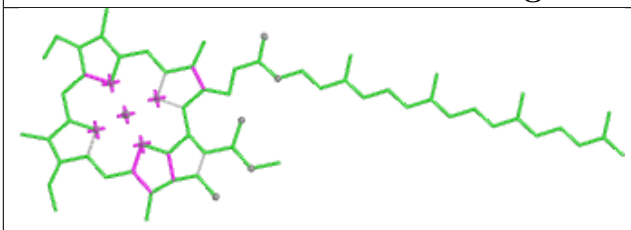
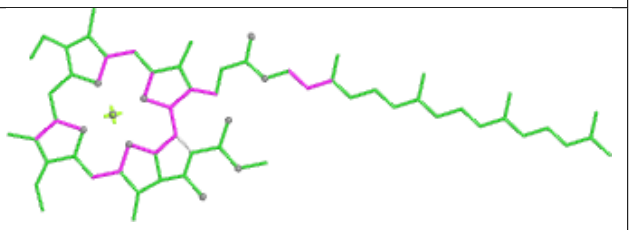
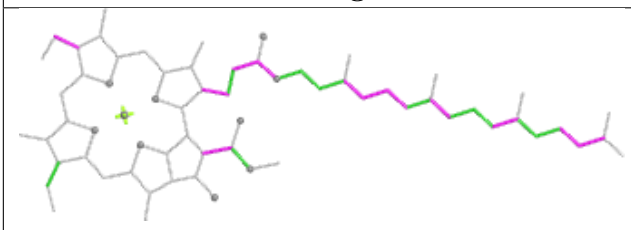
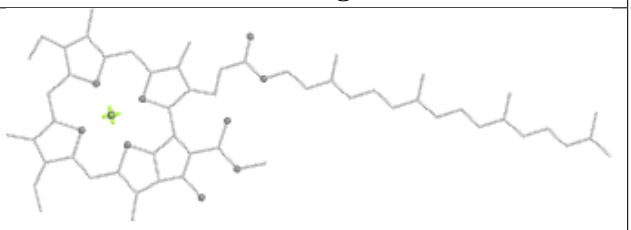


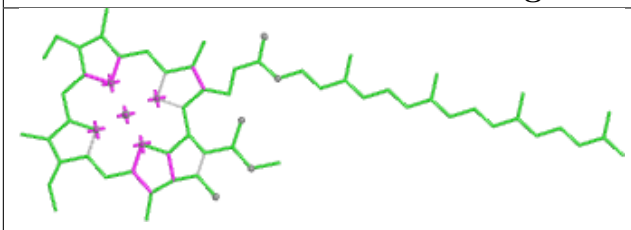
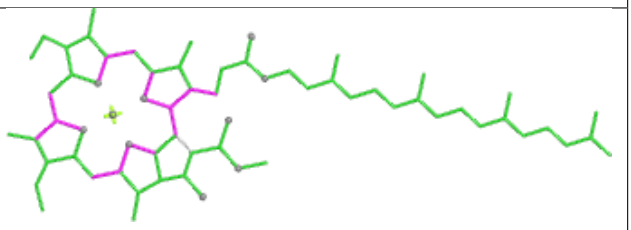
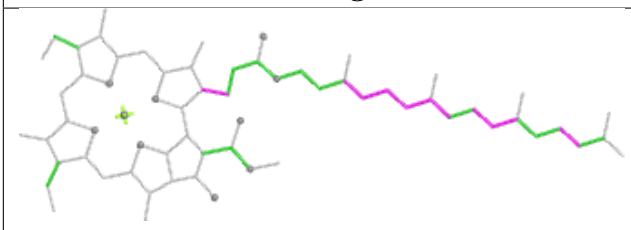
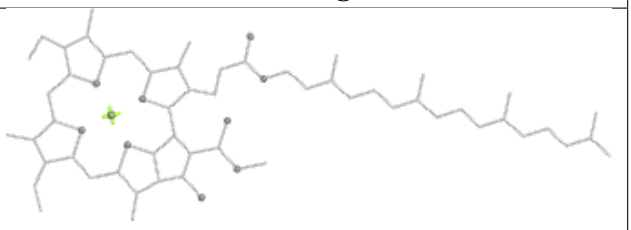
Rings

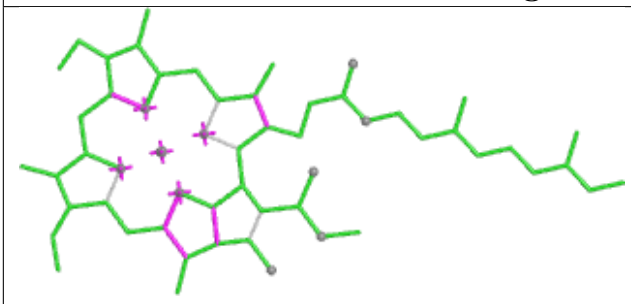
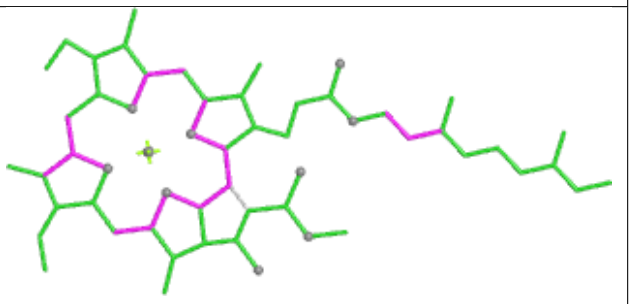
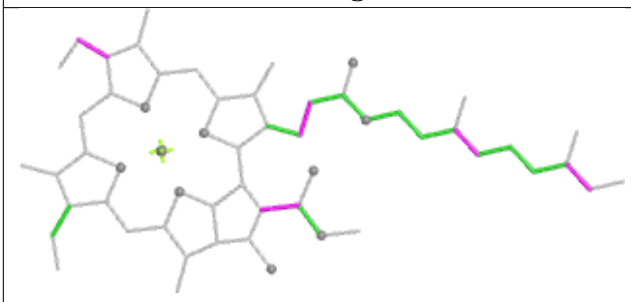
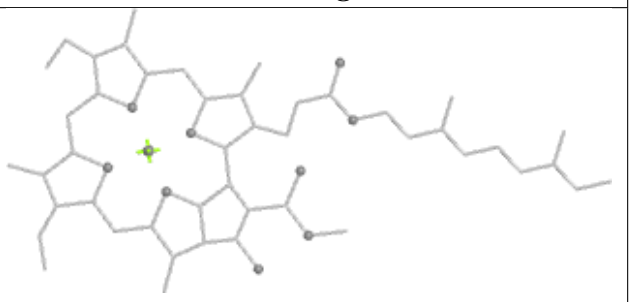
Ligand CLA D 610	
	
Bond lengths	**Bond angles**
	
Torsions	**Rings**

Ligand CLA K 601	
	
Bond lengths	**Bond angles**
	
Torsions	**Rings**

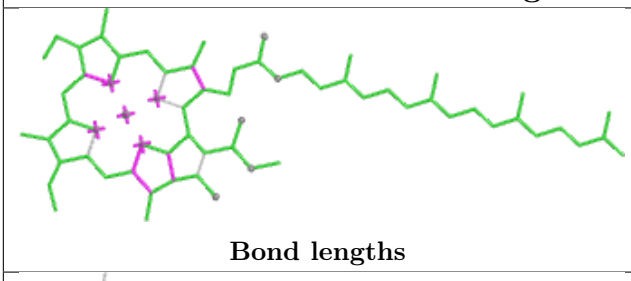
Ligand CLA B 608	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA C 605	
	
Bond lengths	Bond angles
	
Torsions	Rings

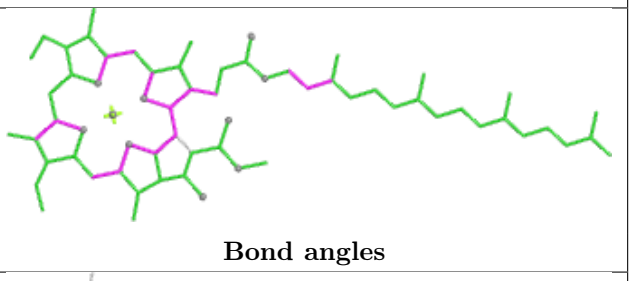
Ligand CLA B 601	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand CLA K 611	
 Bond lengths	 Bond angles
 Torsions	 Rings

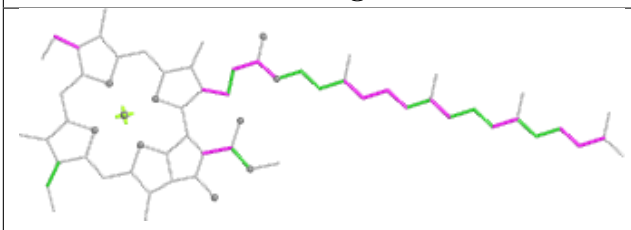
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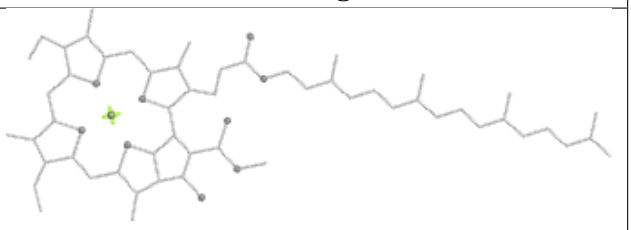
Bond lengths



Bond angles

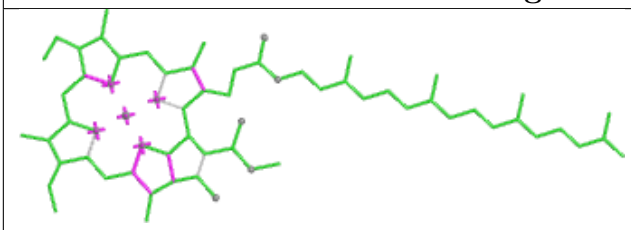


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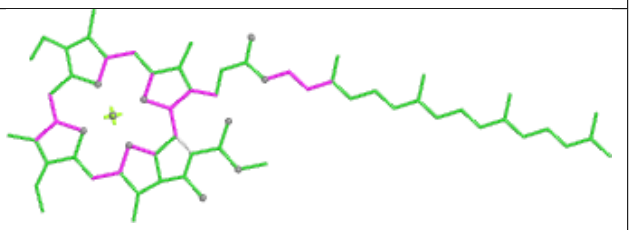


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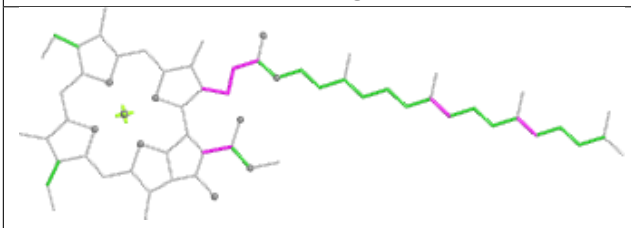
Ligand CLA B 602



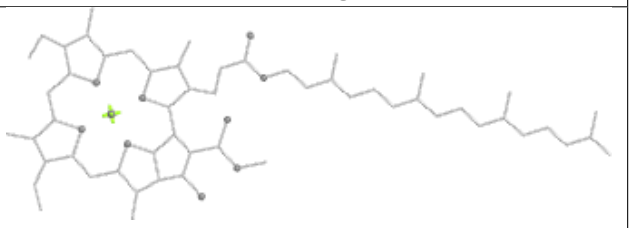
Bond lengths



Bond angles

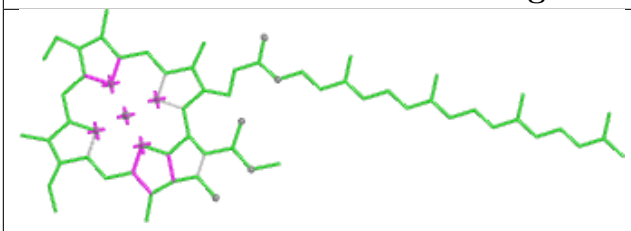


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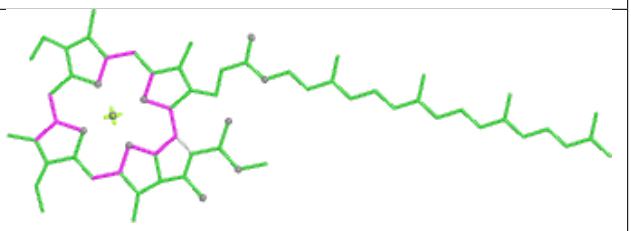


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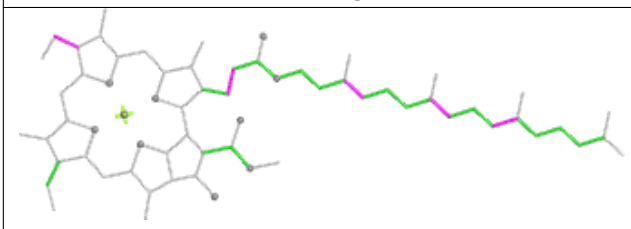
Ligand CLA J 604



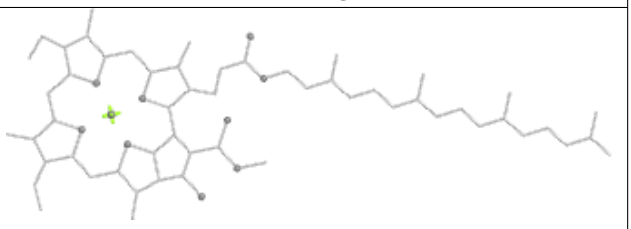
Bond lengths



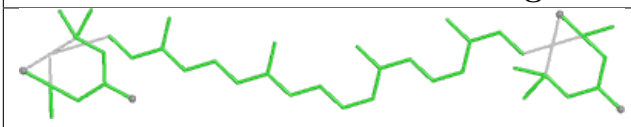
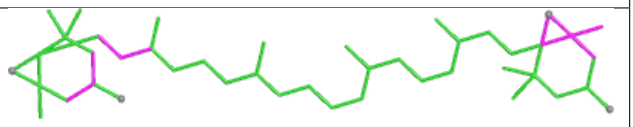
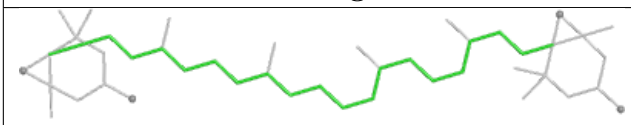
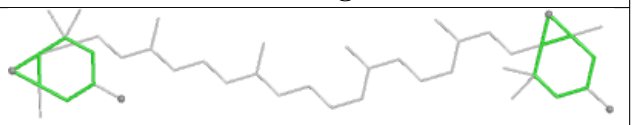
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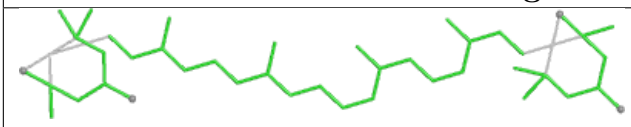
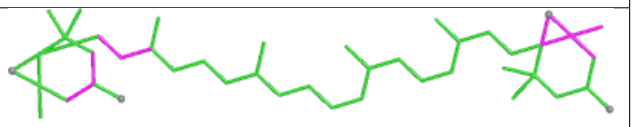
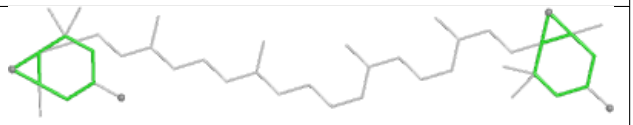


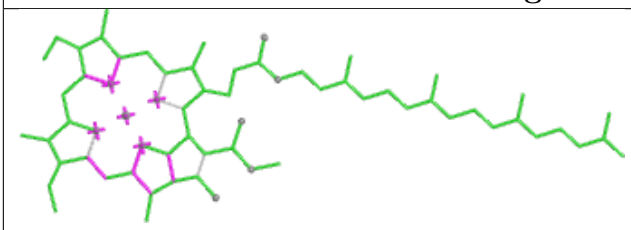
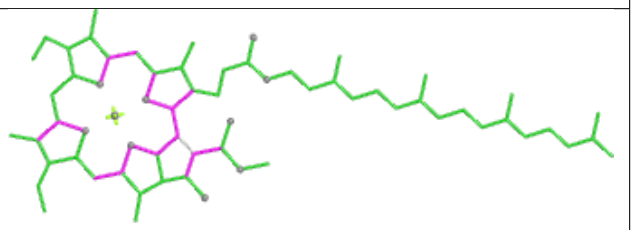
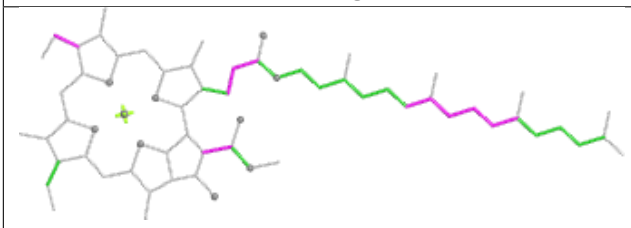
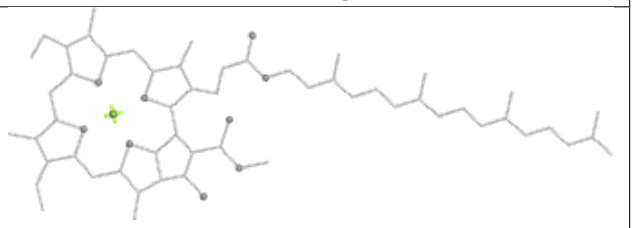
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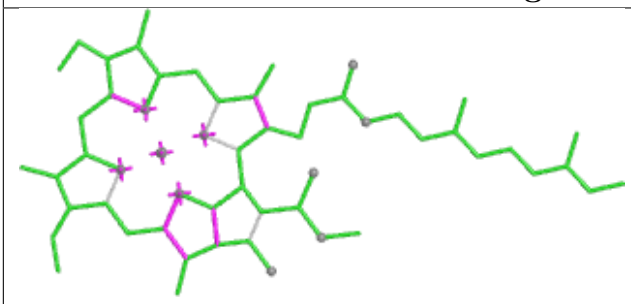
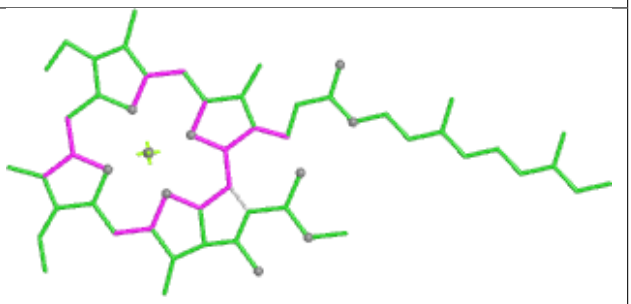
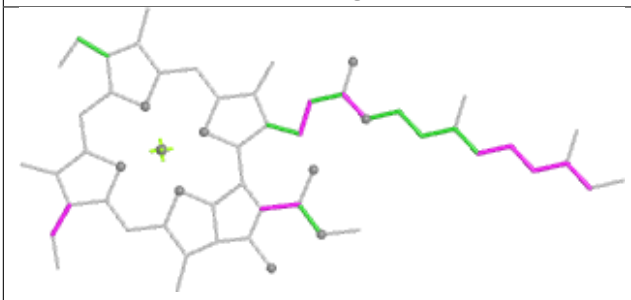
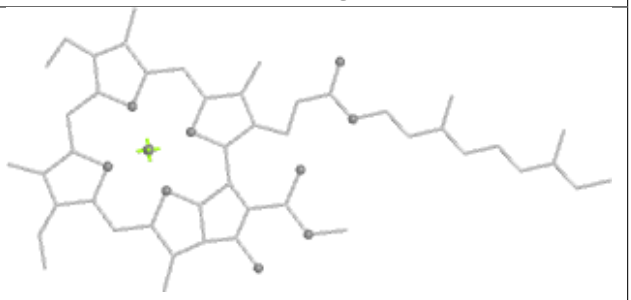


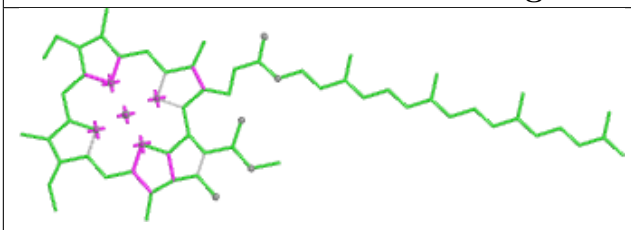
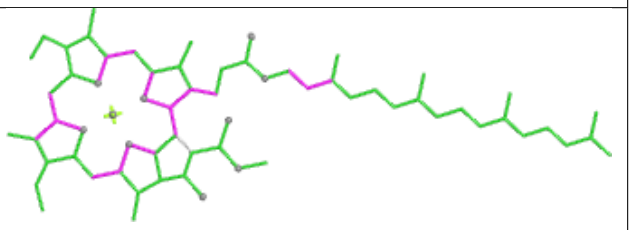
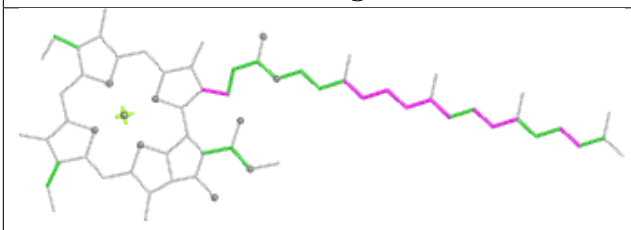
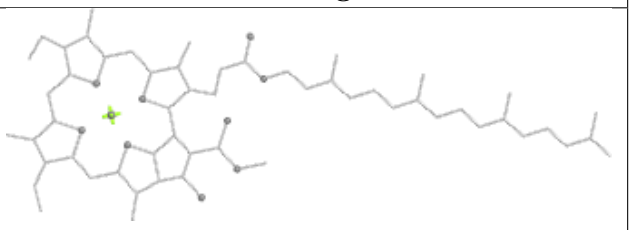
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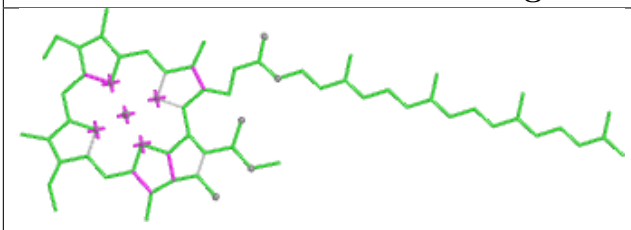
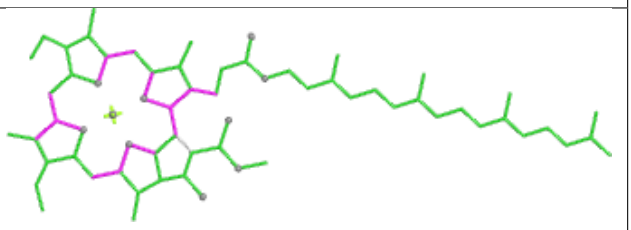
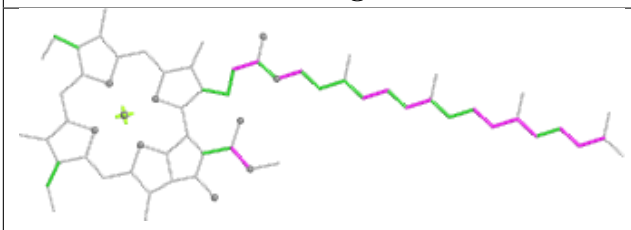
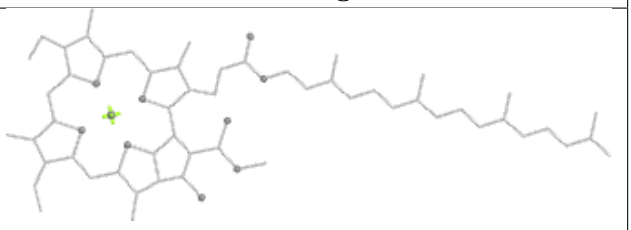
Ligand XAT E 612	
 Bond lengths	 Bond angles
 Torsions	 Rings

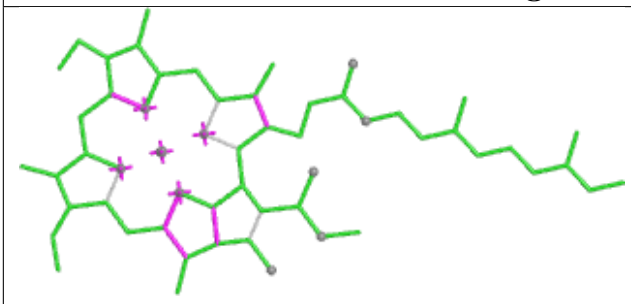
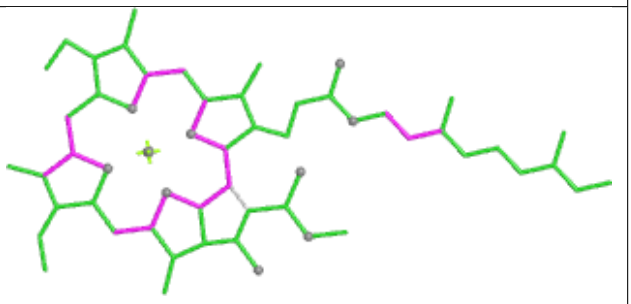
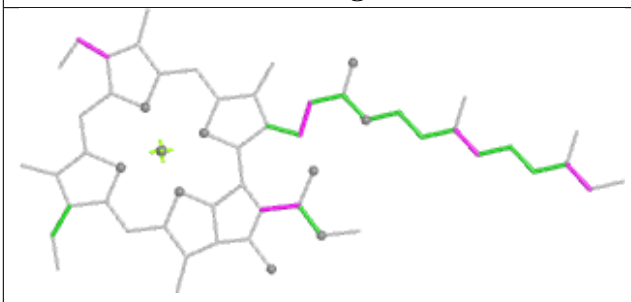
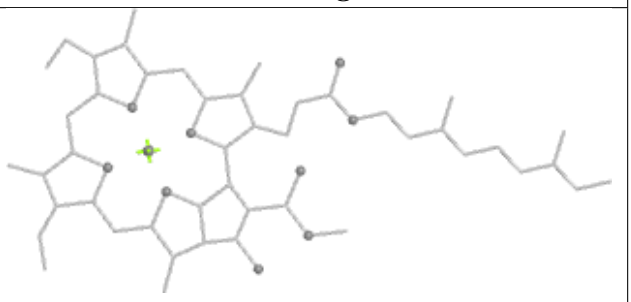
Ligand XAT C 612	
 Bond lengths	 Bond angles
 Torsions	 Rings

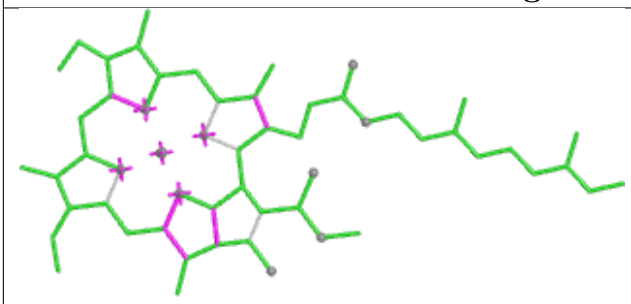
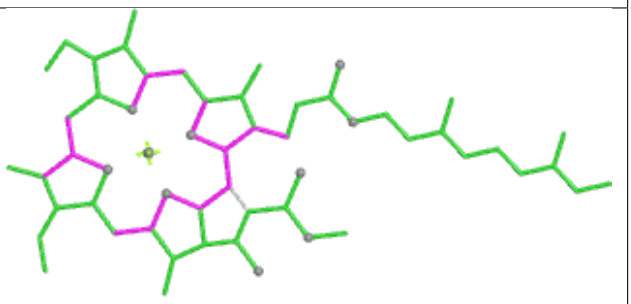
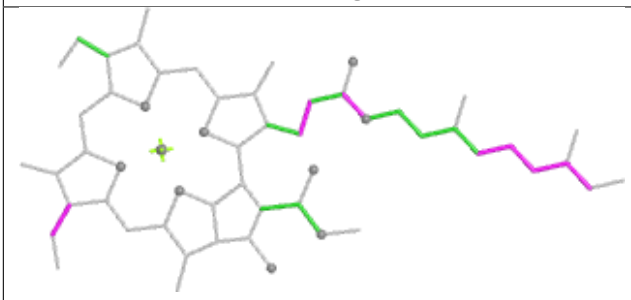
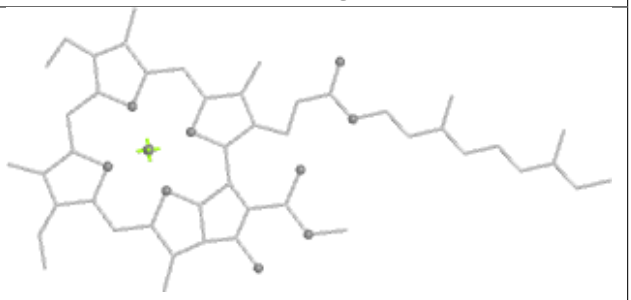
Ligand CLA B 610	
 Bond lengths	 Bond angles
 Torsions	 Rings

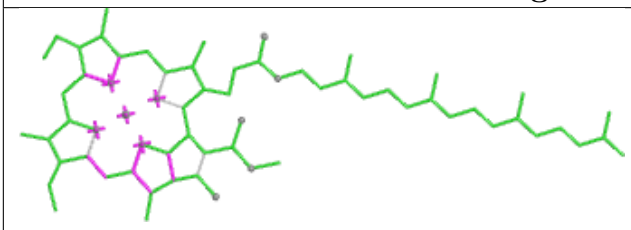
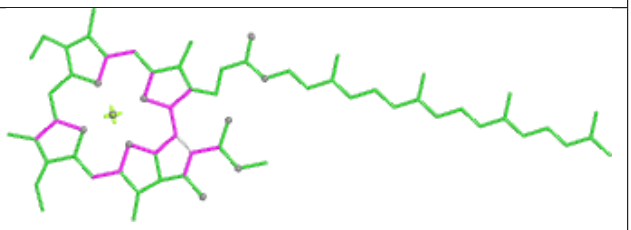
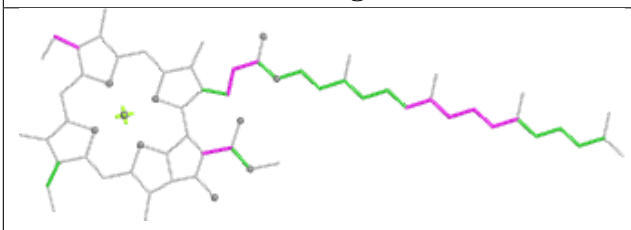
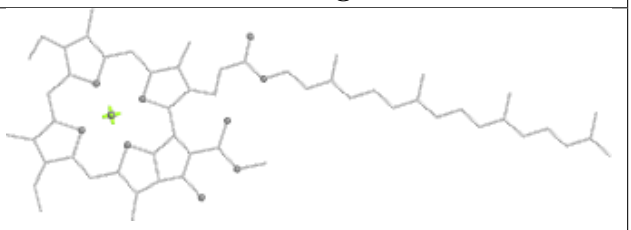
Ligand CLA C 603	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA E 601	
	
Bond lengths	Bond angles
	
Torsions	Rings

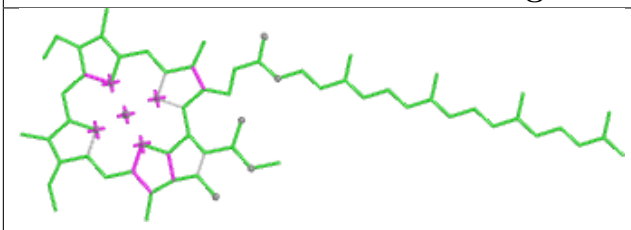
Ligand CLA C 607	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA E 611	
	
Bond lengths	Bond angles
	
Torsions	Rings

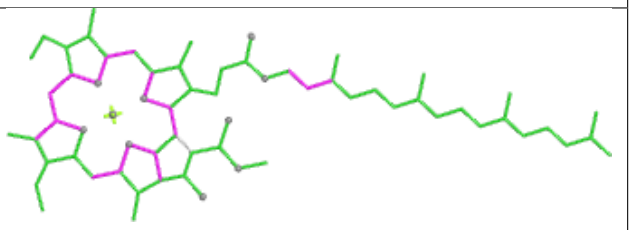
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Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA I 610	
	
Bond lengths	Bond angles
	
Torsions	Rings

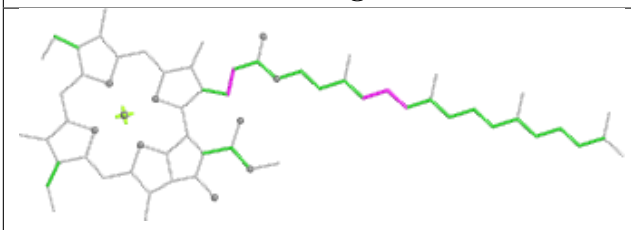
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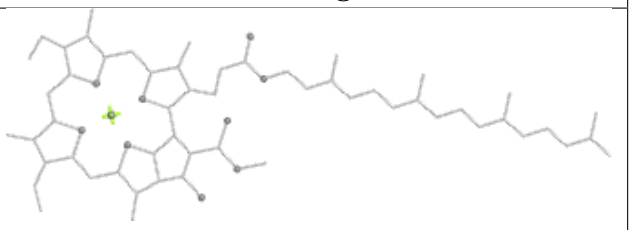
Bond lengths



Bond angles

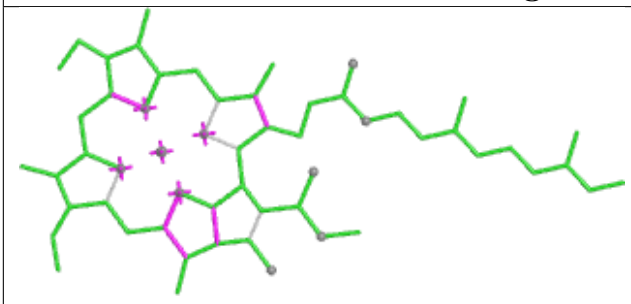


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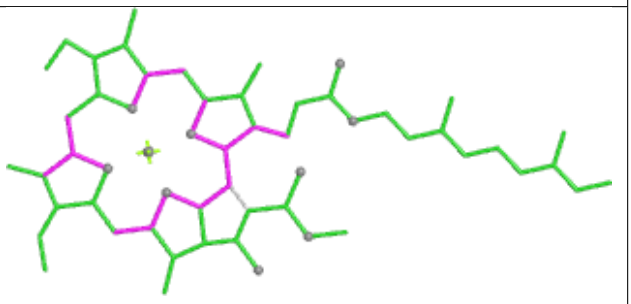


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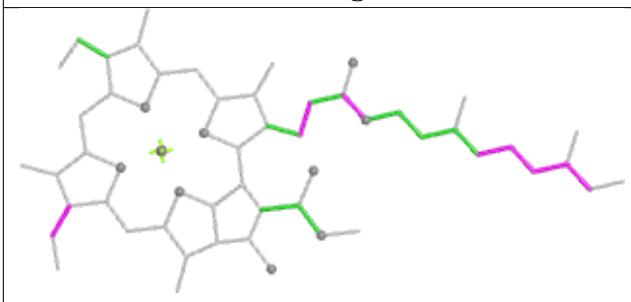
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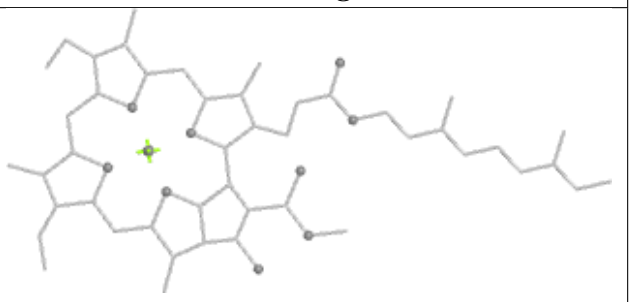
Bond lengths



Bond angles

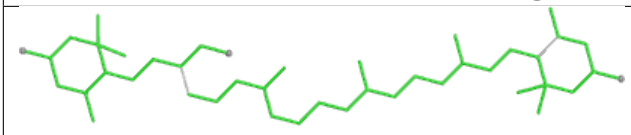


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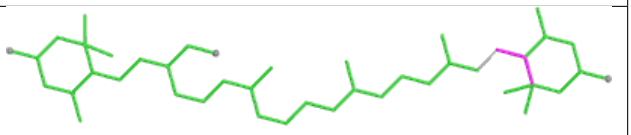


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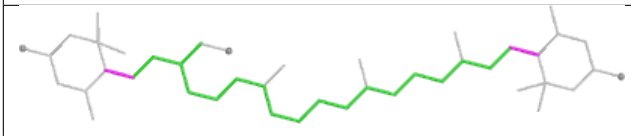
Ligand 32N H 613



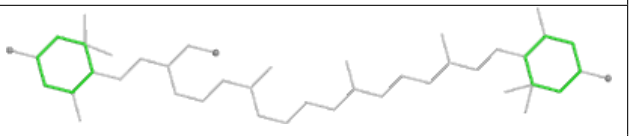
Bond lengths



Bond angles

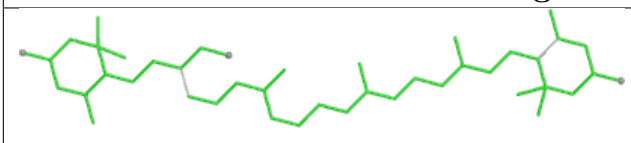


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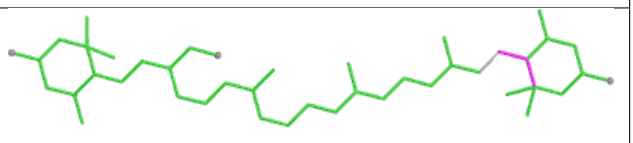


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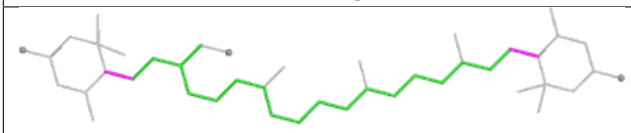
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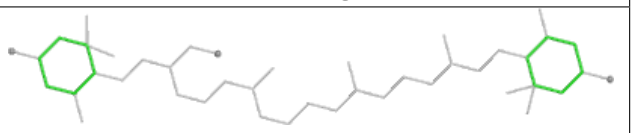
Bond lengths



Bond angles

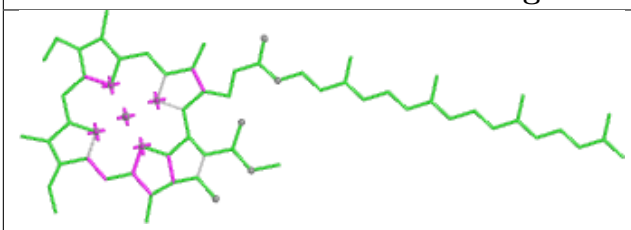


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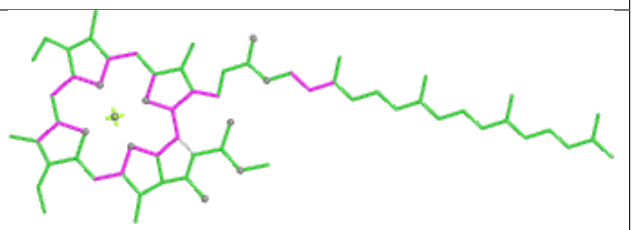


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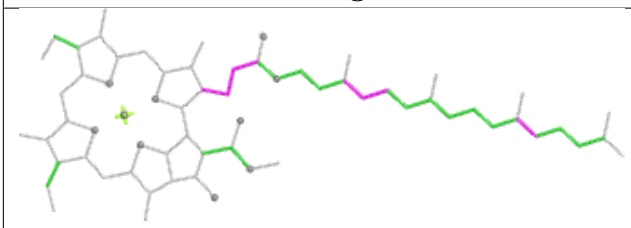
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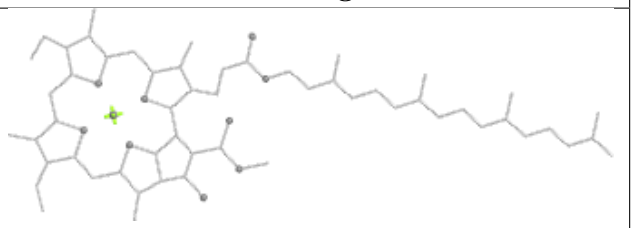
Bond lengths



Bond angles

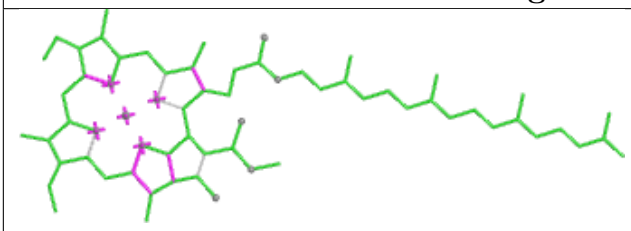


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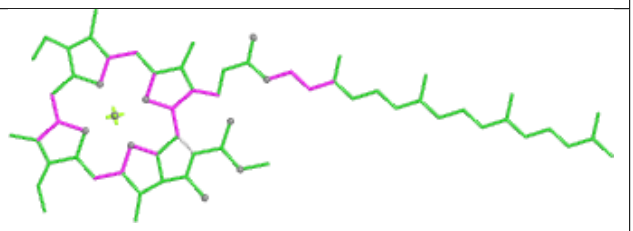


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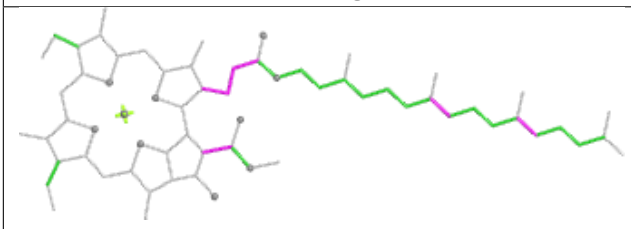
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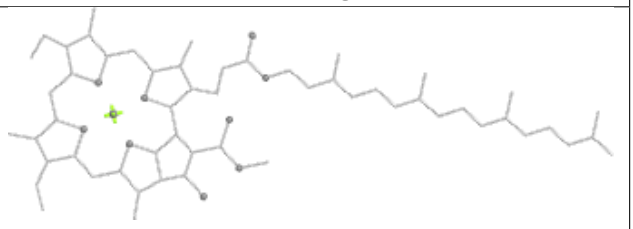
Bond lengths



Bond angles

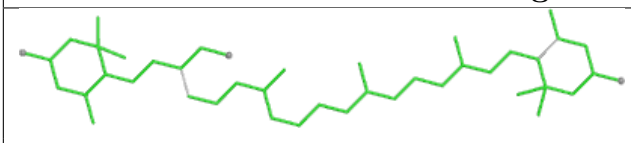


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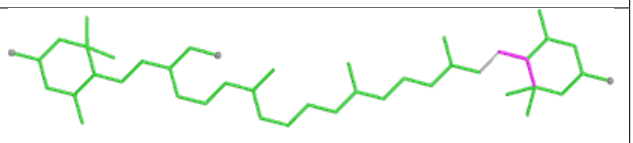


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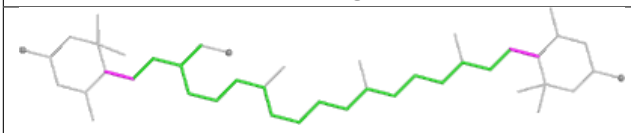
Ligand 32N F 613



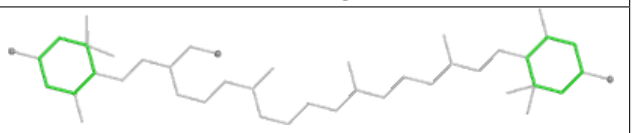
Bond lengths



Bond angles

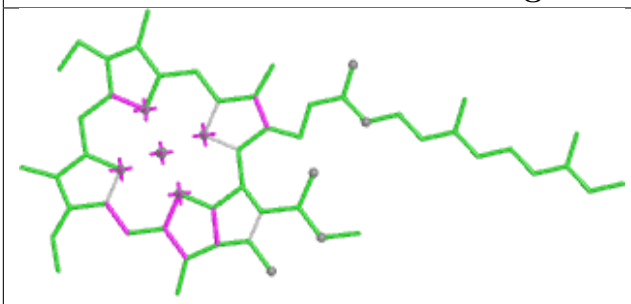


Torsions

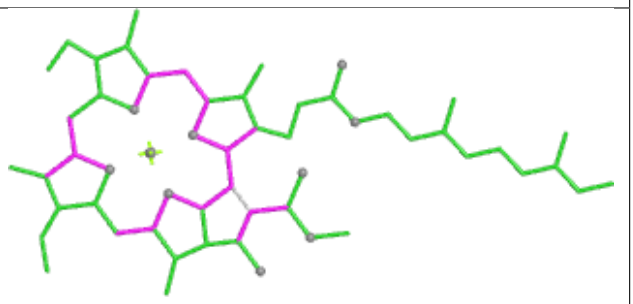


Rings

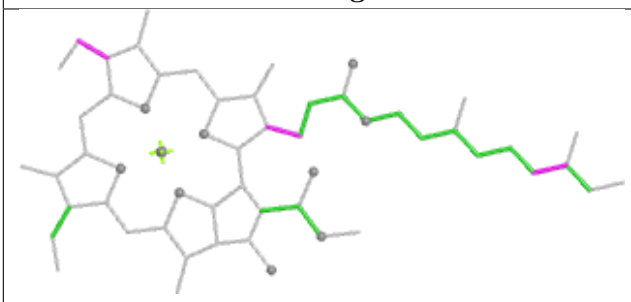
Ligand CLA C 608



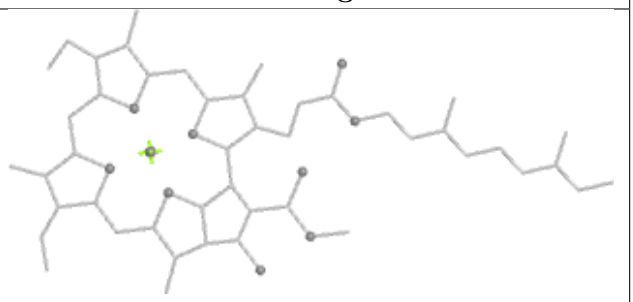
Bond lengths



Bond angles

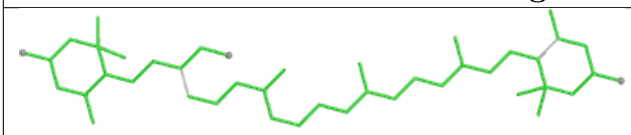


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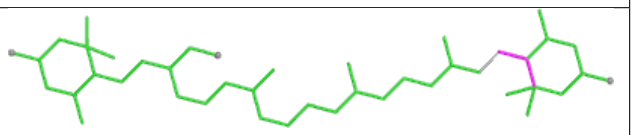


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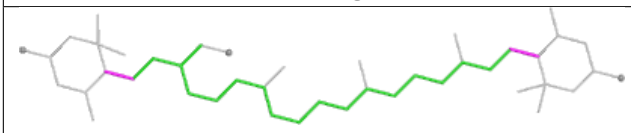
Ligand 32N A 613



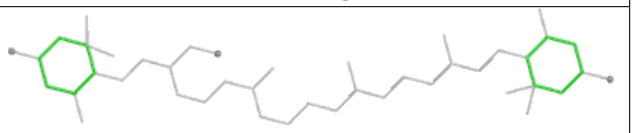
Bond lengths



Bond angles

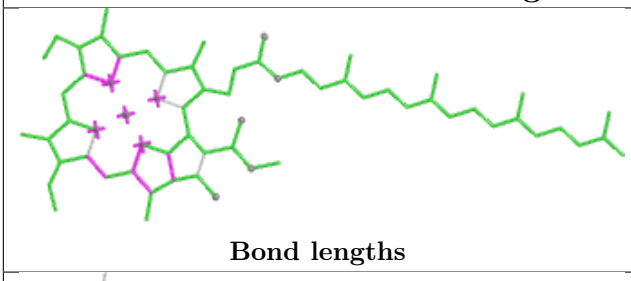


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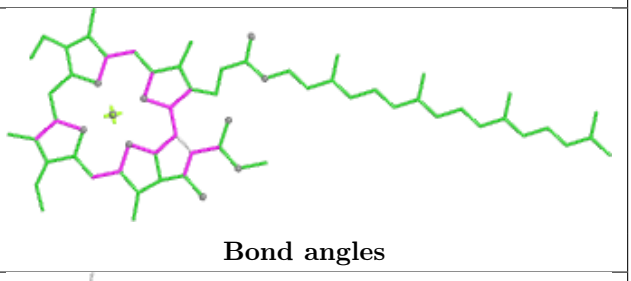


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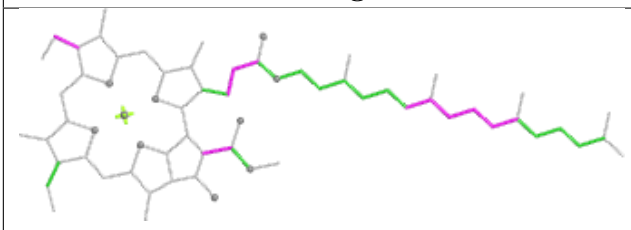
Ligand CLA G 610



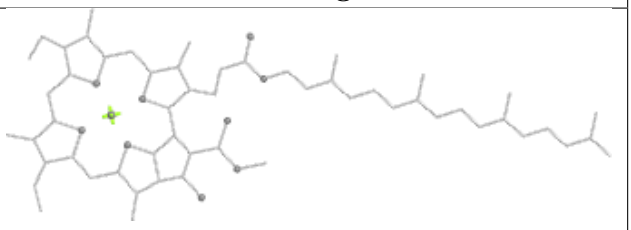
Bond lengths



Bond angles

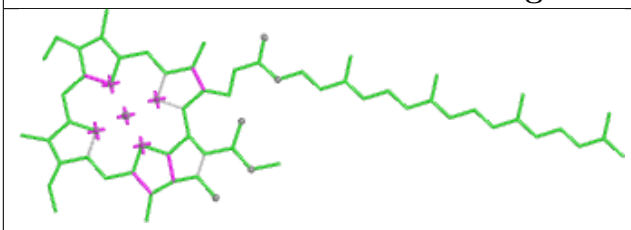


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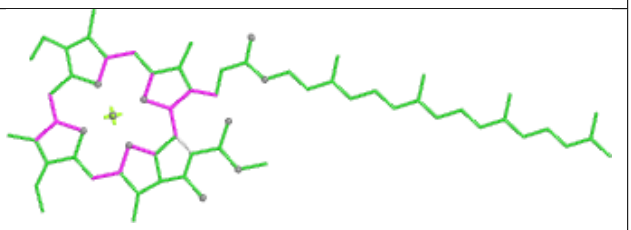


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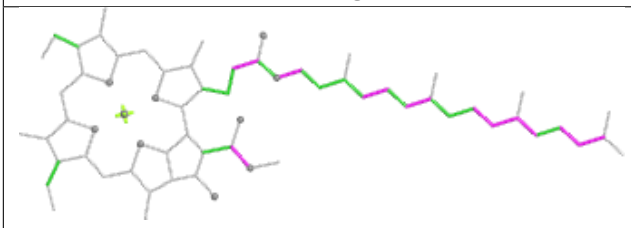
Ligand CLA E 607



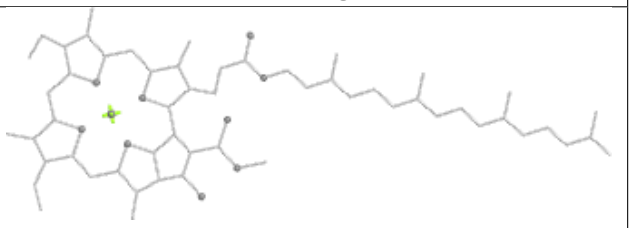
Bond lengths



Bond angles

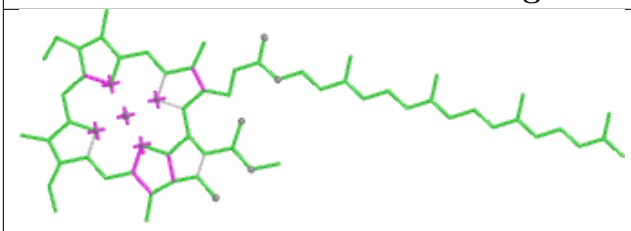


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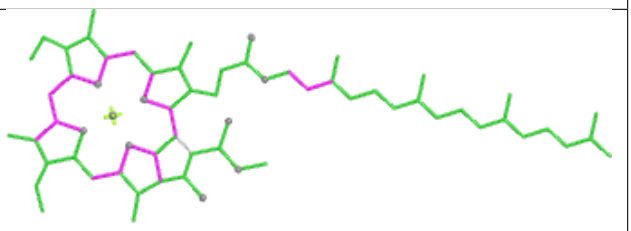


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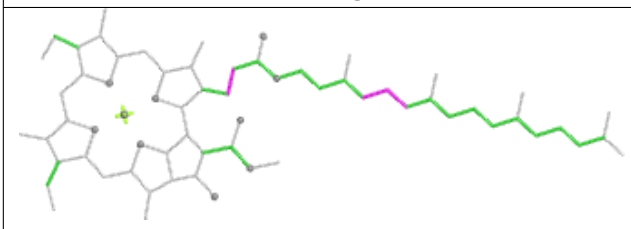
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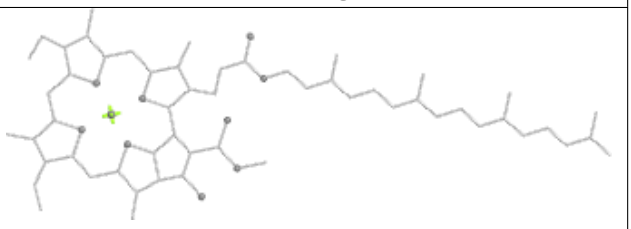
Bond lengths



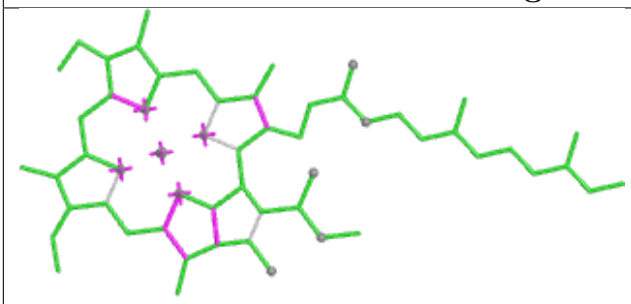
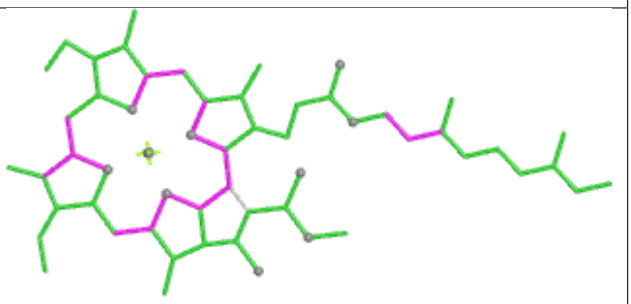
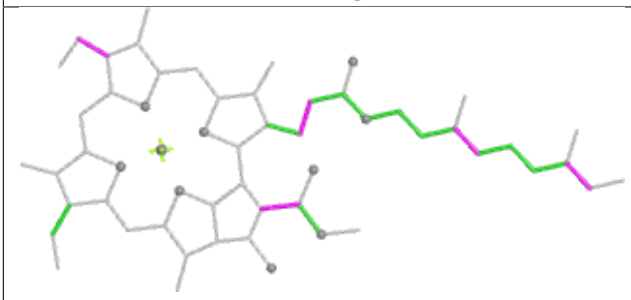
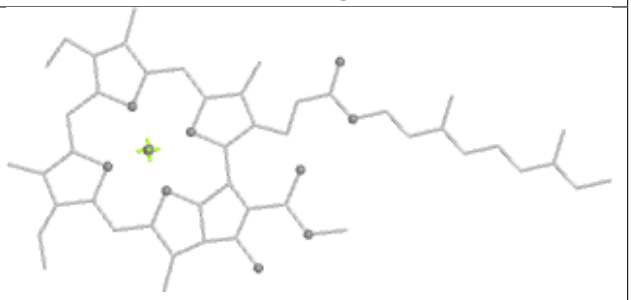
Bond angles

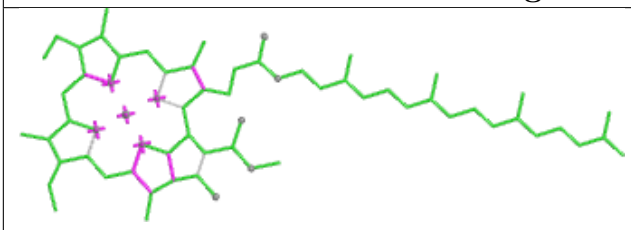
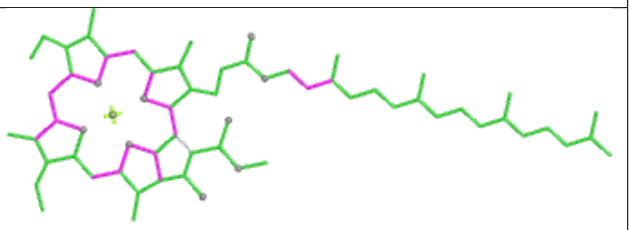
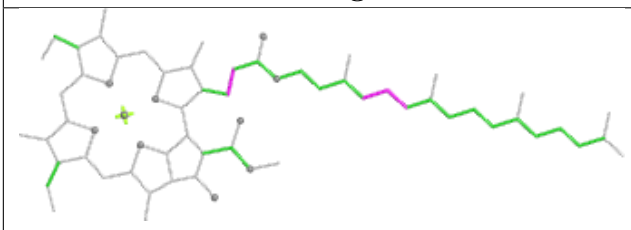
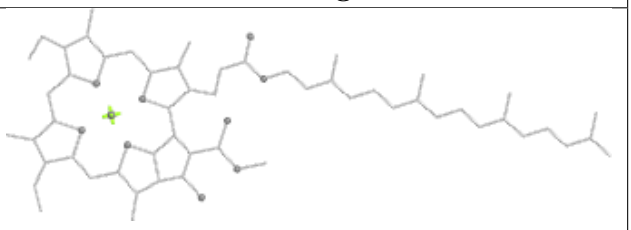


Torsions

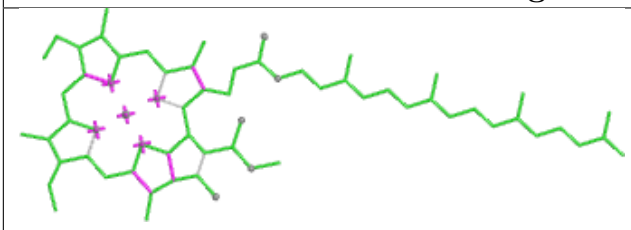


Rings

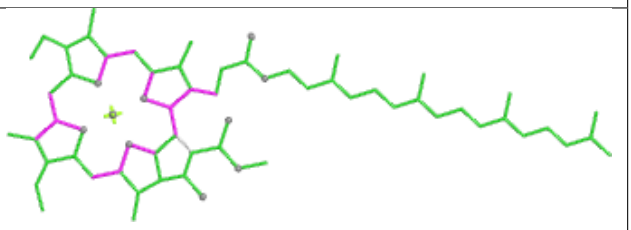
Ligand CLA C 611	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA B 609	
	
Bond lengths	Bond angles
	
Torsions	Rings

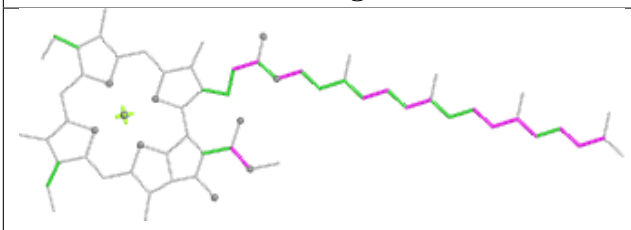
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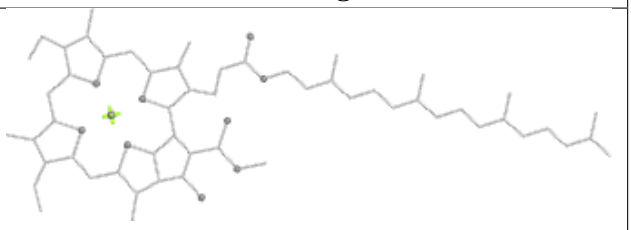
Bond lengths



Bond angles

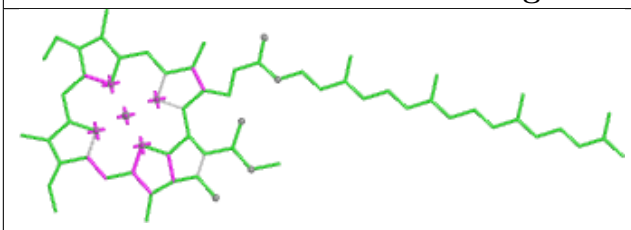


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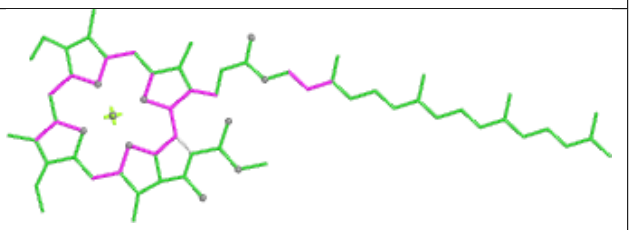


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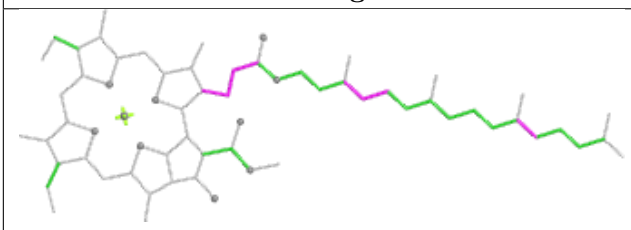
Ligand CLA K 606



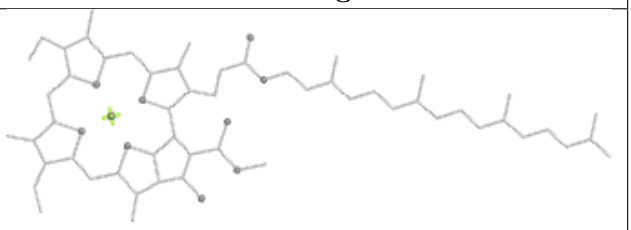
Bond lengths



Bond angles

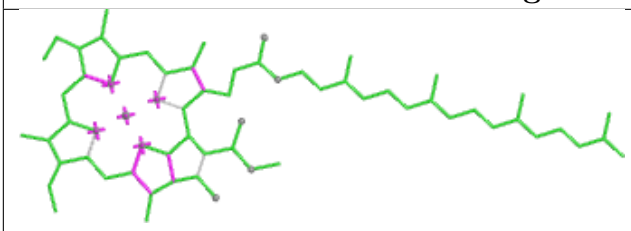


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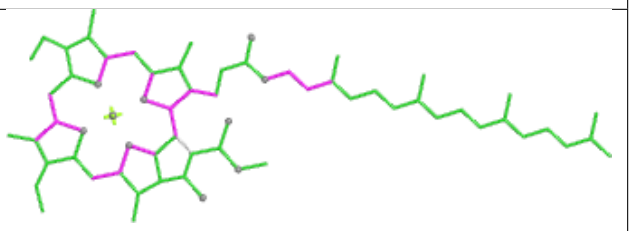


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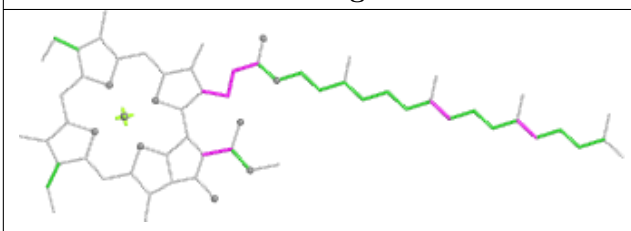
Ligand CLA H 602



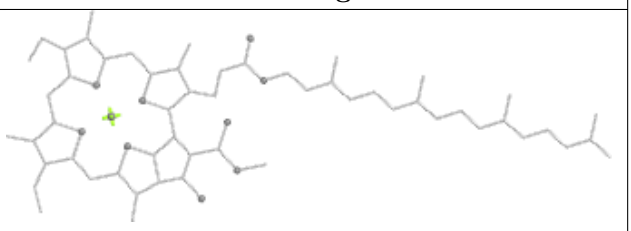
Bond lengths



Bond angles

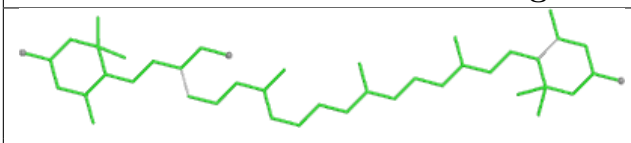


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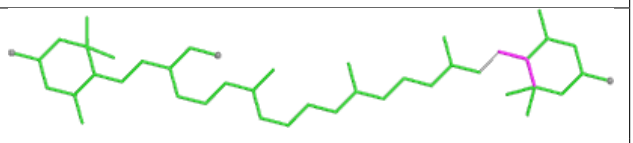


Rings

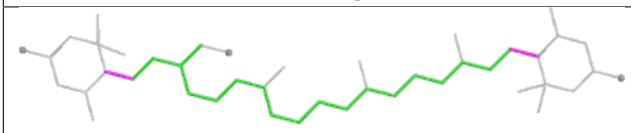
Ligand 32N B 613



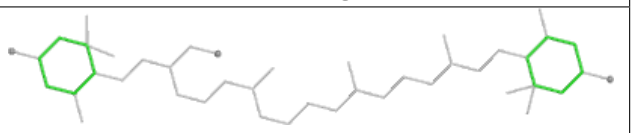
Bond lengths



Bond angles

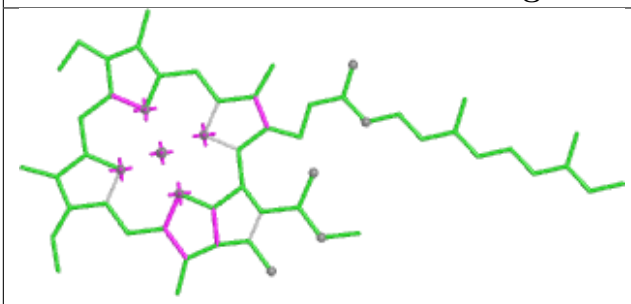


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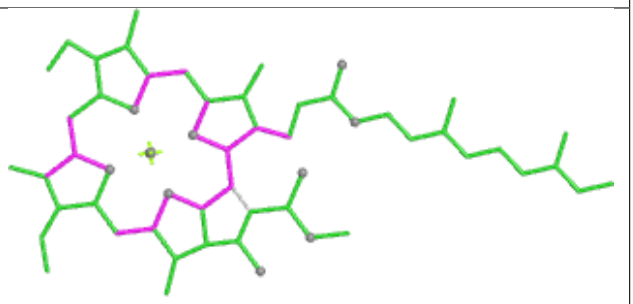


Rings

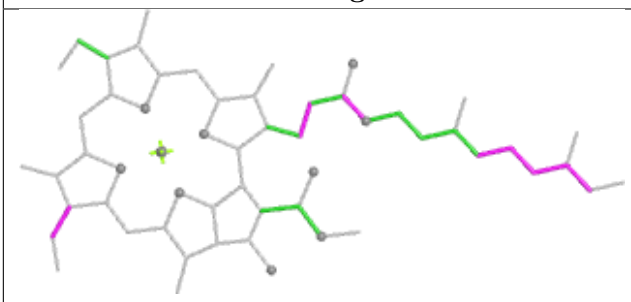
Ligand CLA G 603



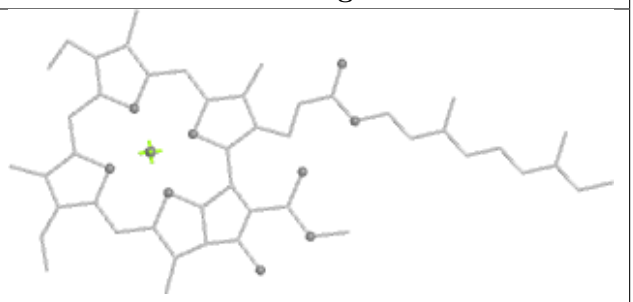
Bond lengths



Bond angles

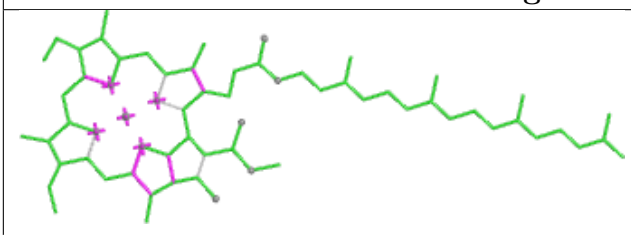


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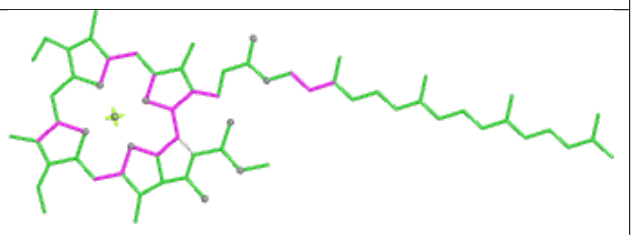


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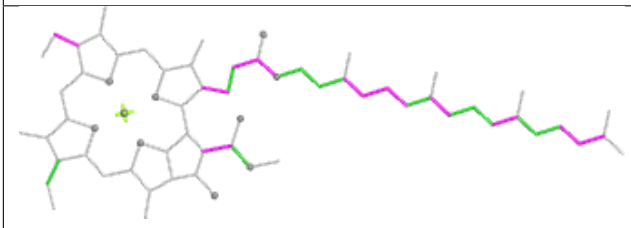
Ligand CLA H 605



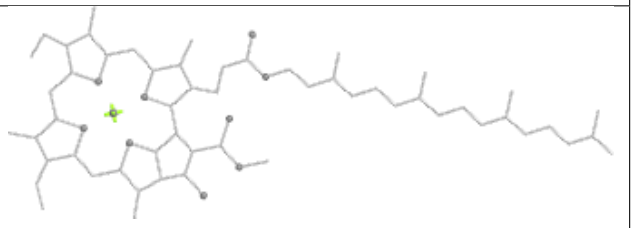
Bond lengths



Bond angles

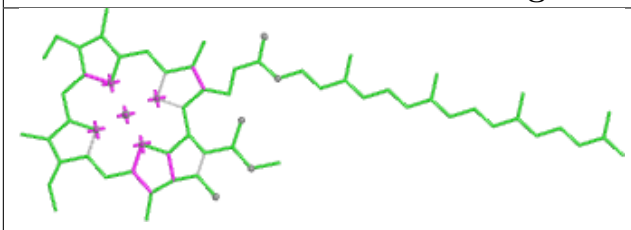


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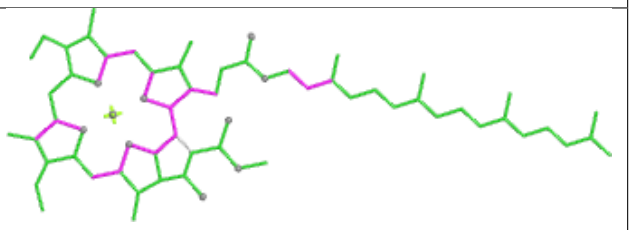


Rings

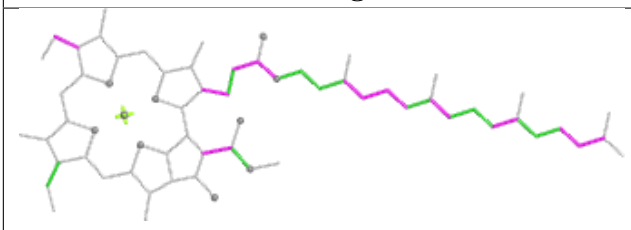
Ligand CLA J 605



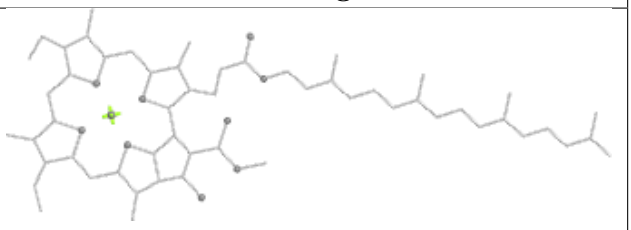
Bond lengths



Bond angles

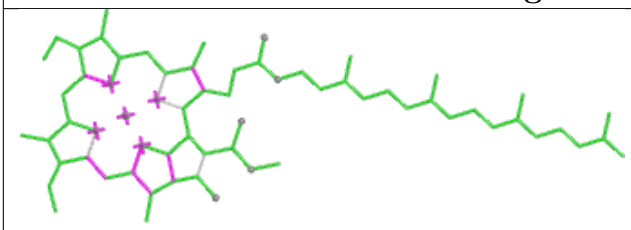


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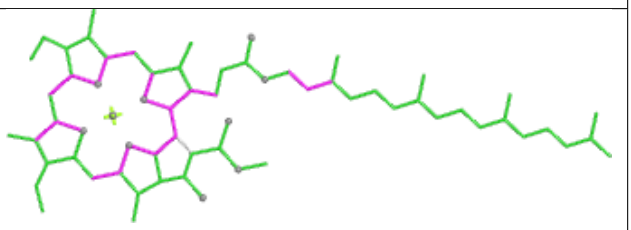


Rings

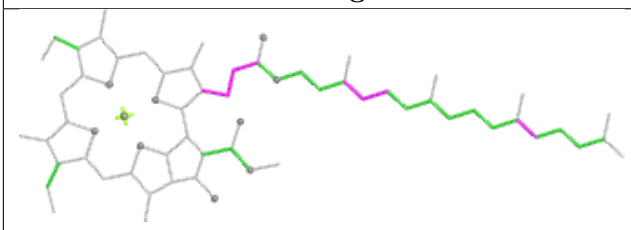
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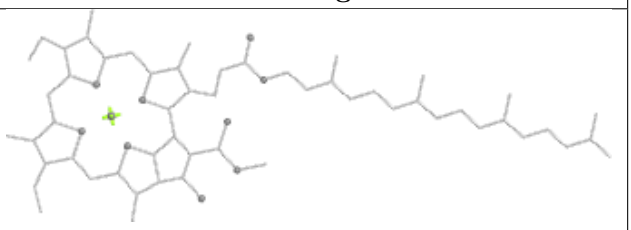
Bond lengths



Bond angles

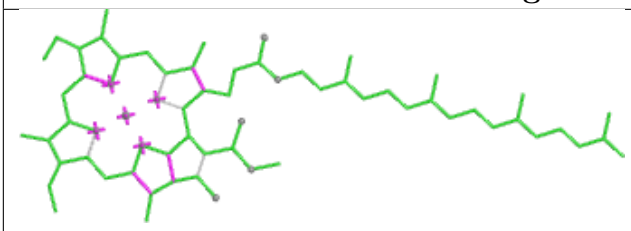


Torsions

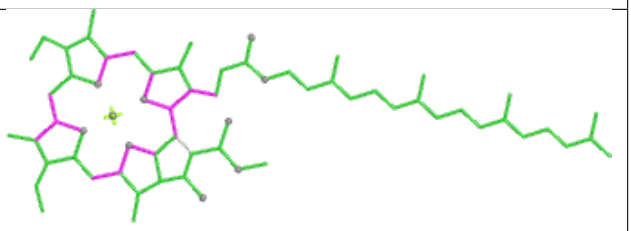


Rings

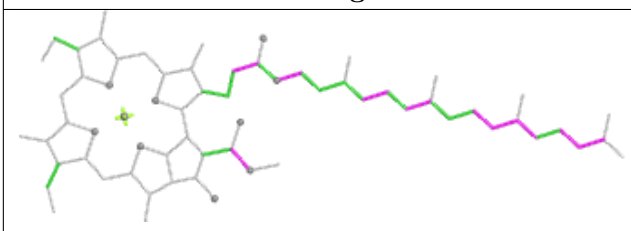
Ligand CLA K 607



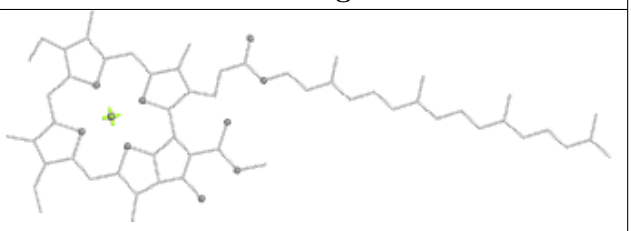
Bond lengths



Bond angles

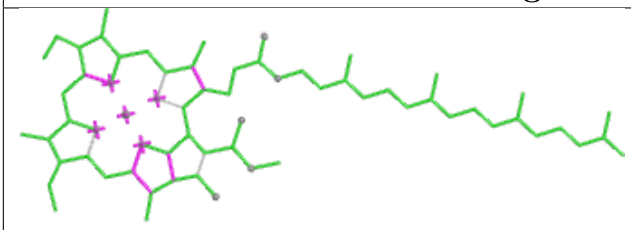


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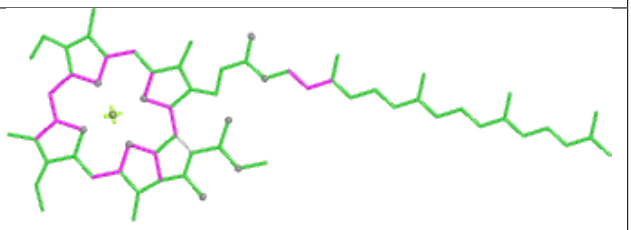


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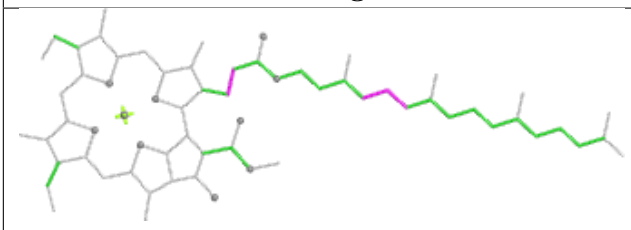
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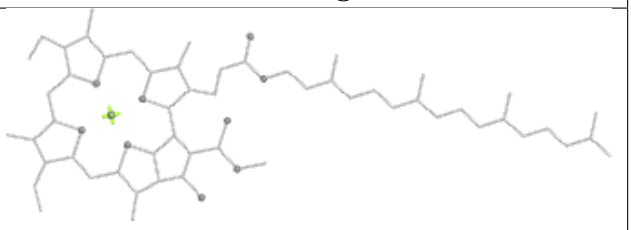
Bond lengths



Bond angles

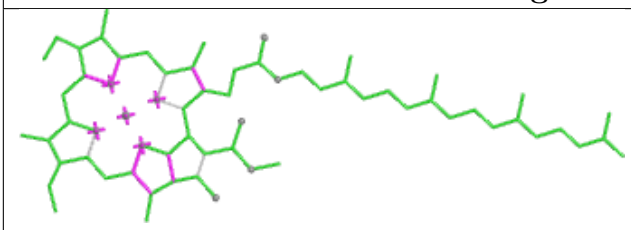


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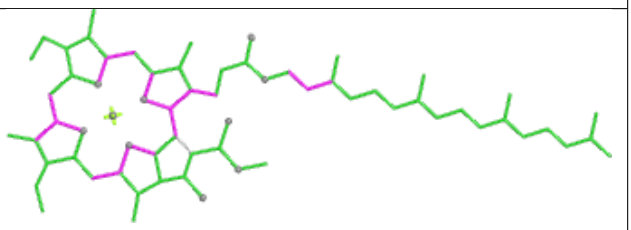


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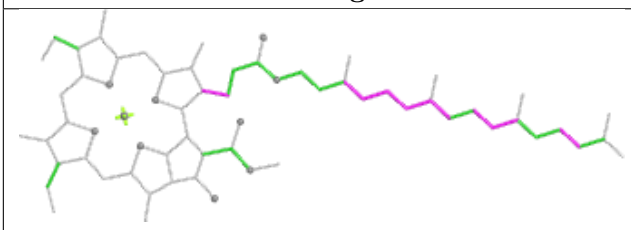
Ligand CLA I 601



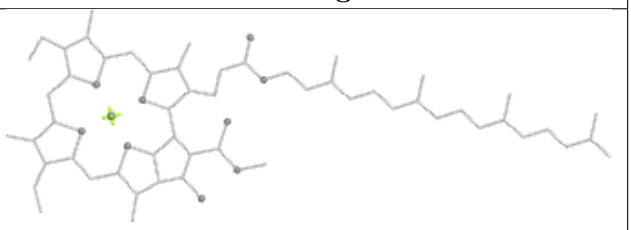
Bond lengths



Bond angles

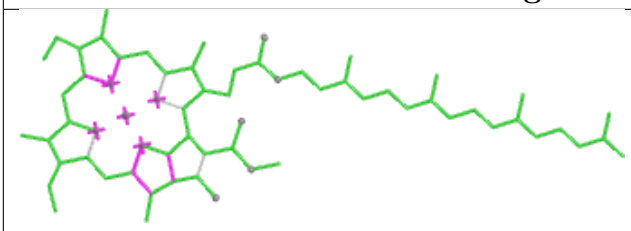


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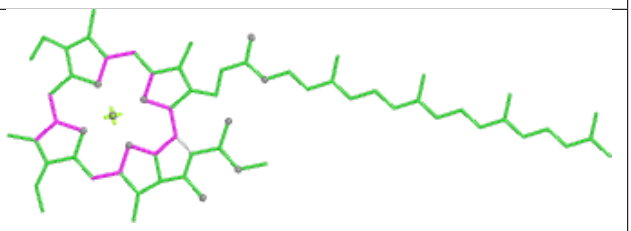


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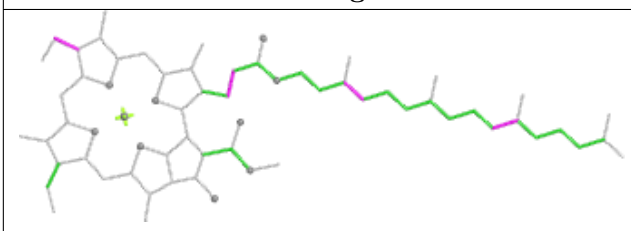
Ligand CLA A 604



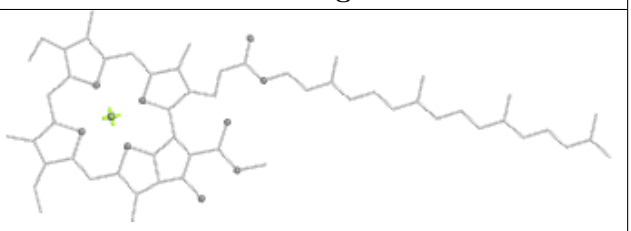
Bond lengths



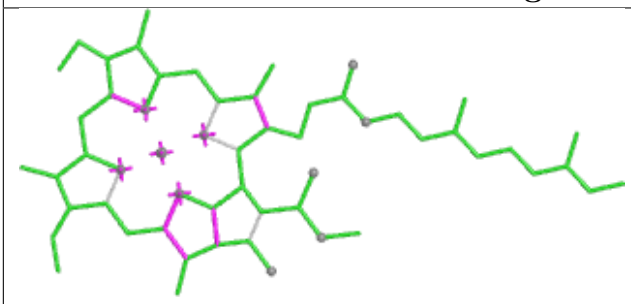
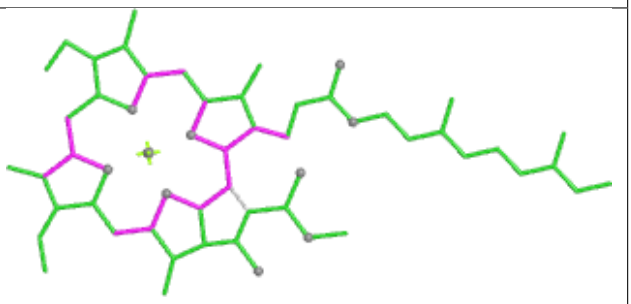
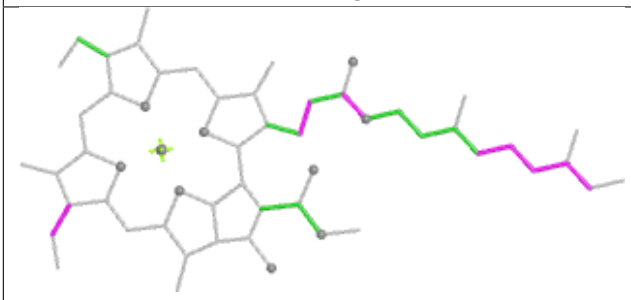
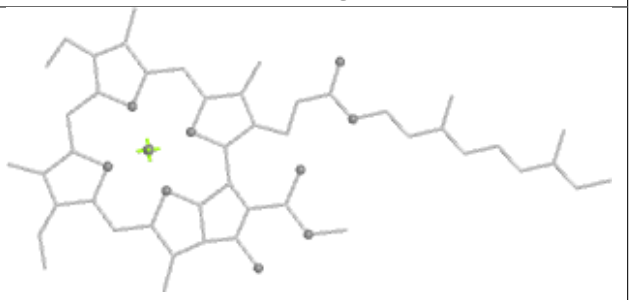
Bond angles

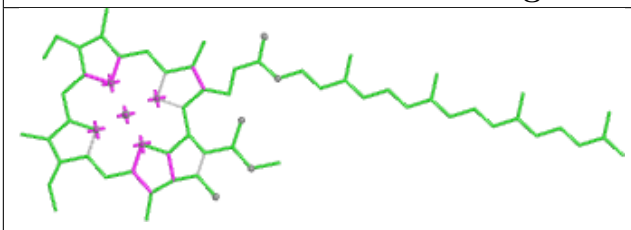
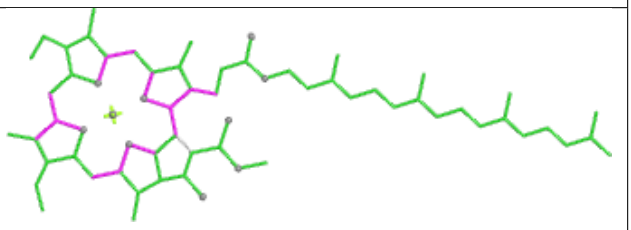
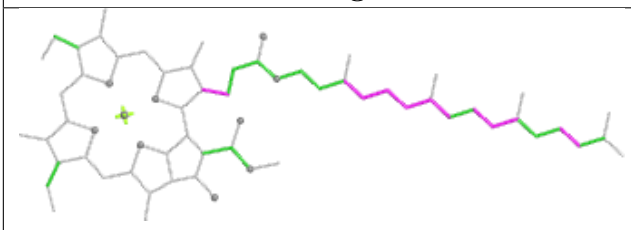
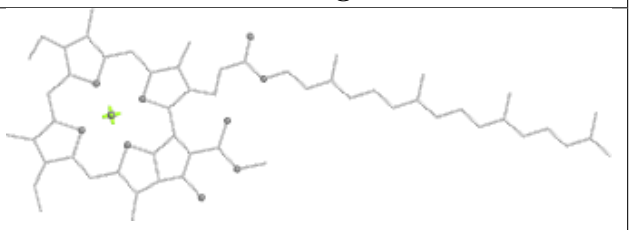


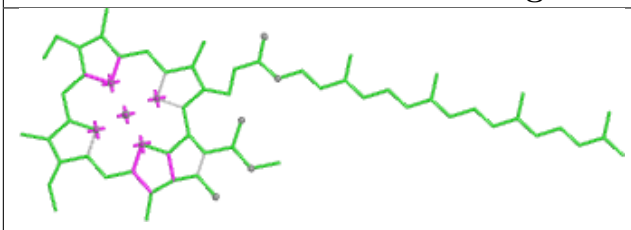
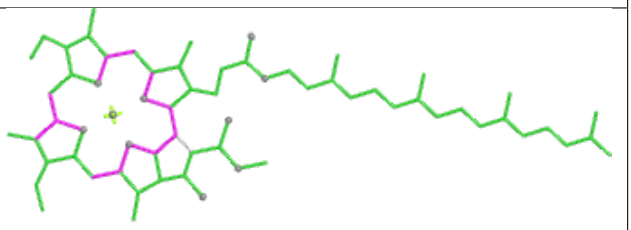
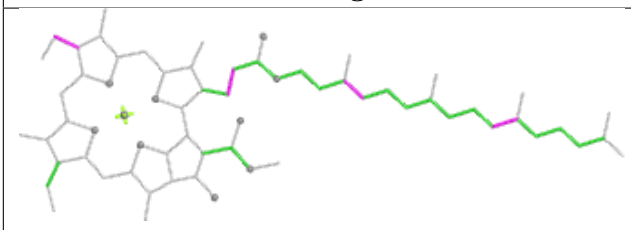
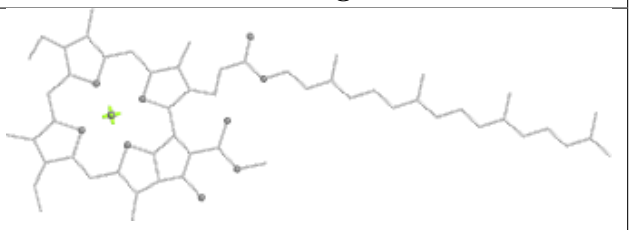
Torsions

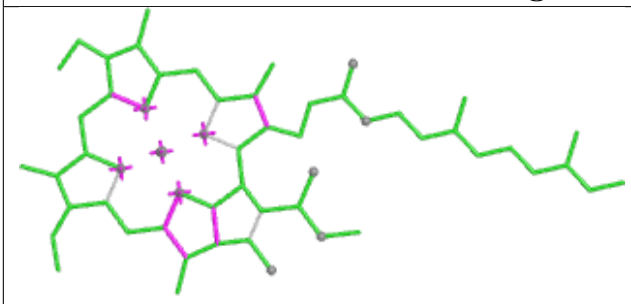
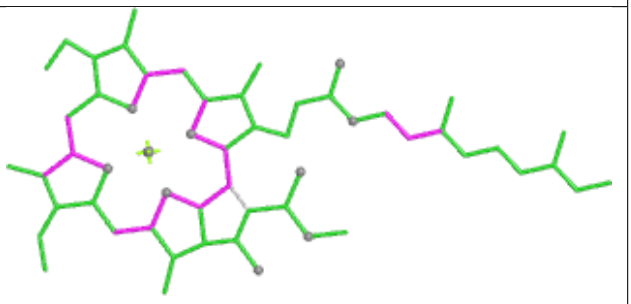
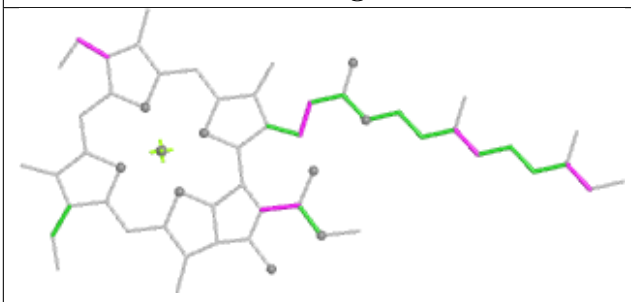
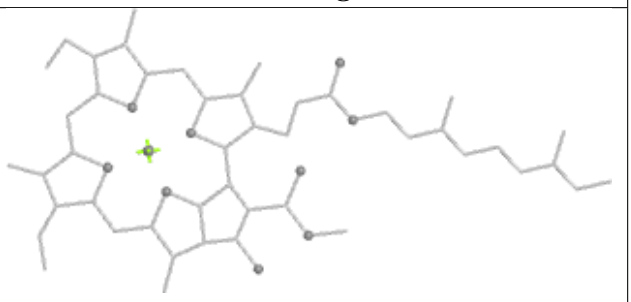


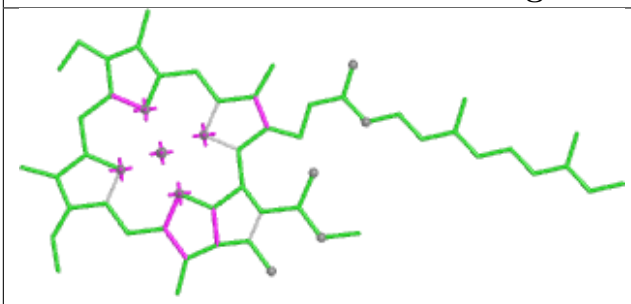
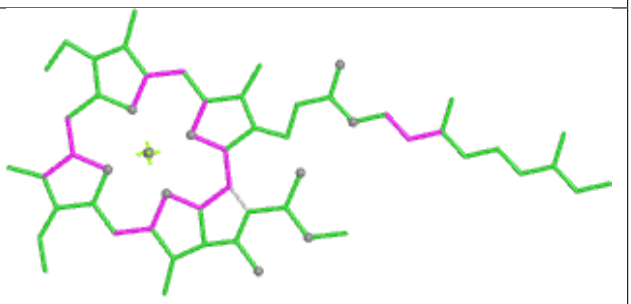
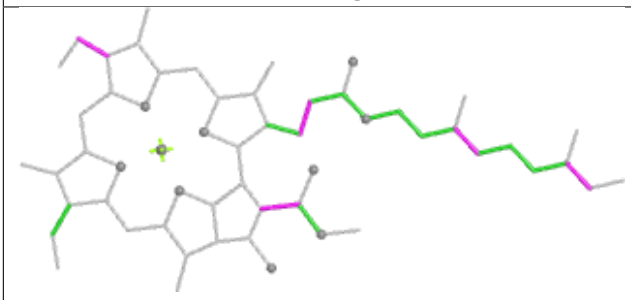
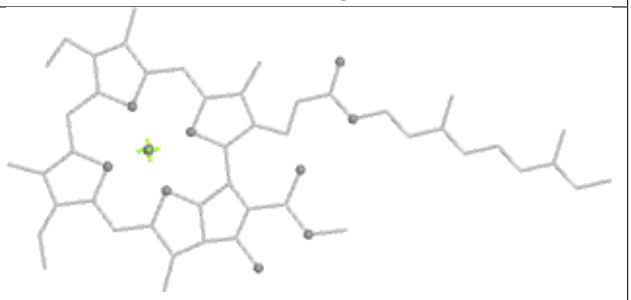
Rings

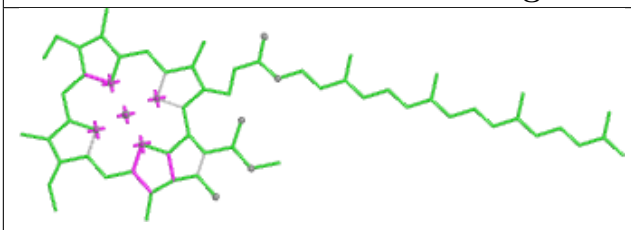
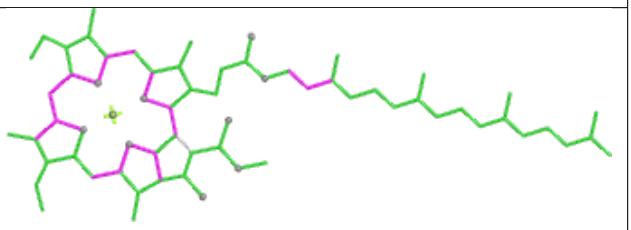
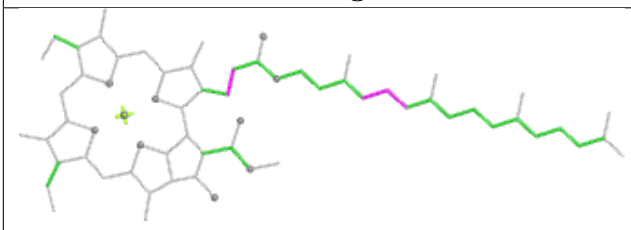
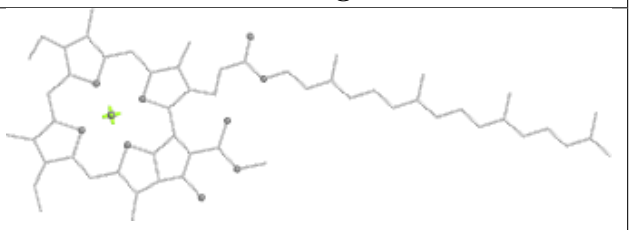
Ligand CLA B 603	
	
Bond lengths	Bond angles
	
Torsions	Rings

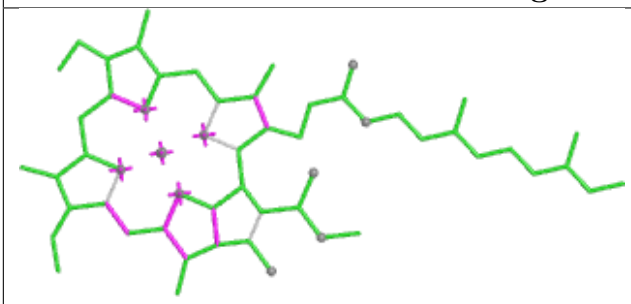
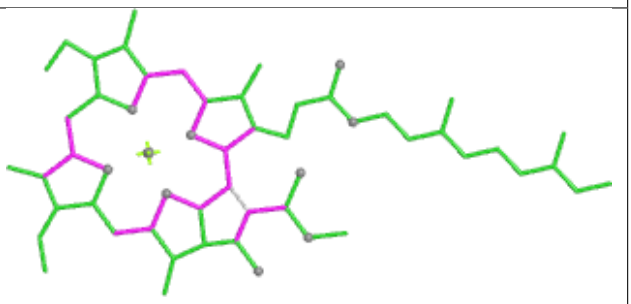
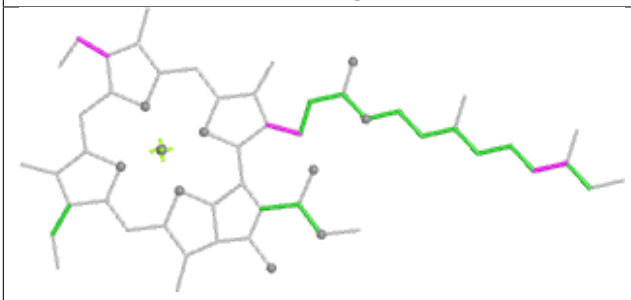
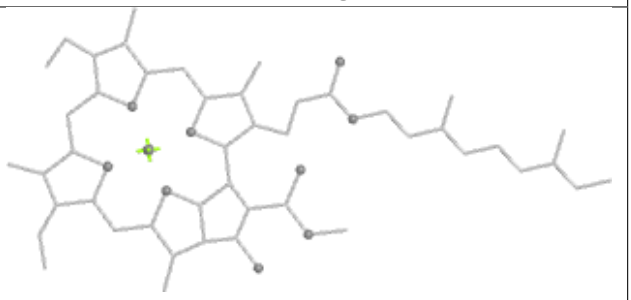
Ligand CLA H 601	
	
Bond lengths	Bond angles
	
Torsions	Rings

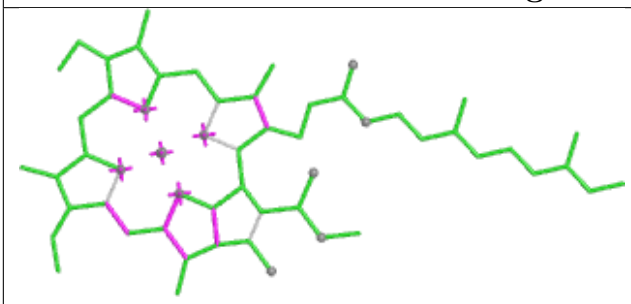
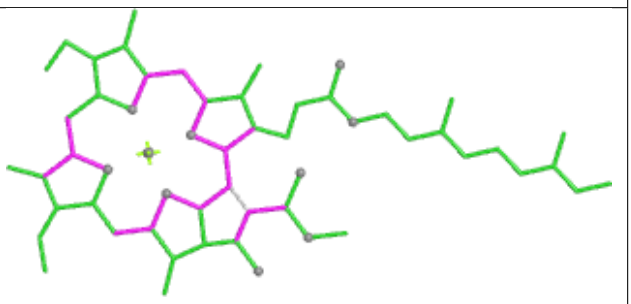
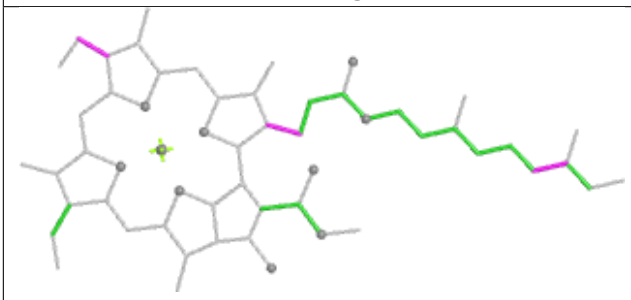
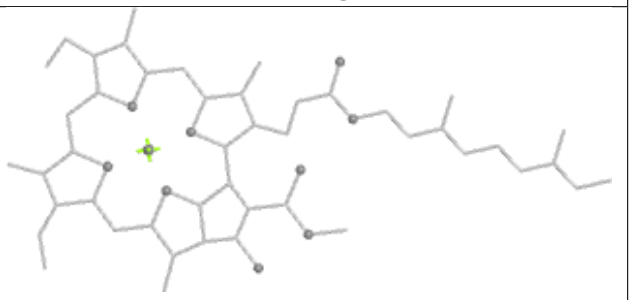
Ligand CLA F 604	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand CLA J 611	
 Bond lengths	 Bond angles
 Torsions	 Rings

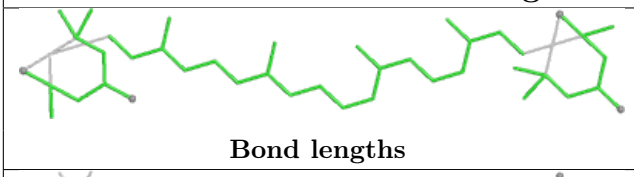
Ligand CLA F 611	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA K 609	
	
Bond lengths	Bond angles
	
Torsions	Rings

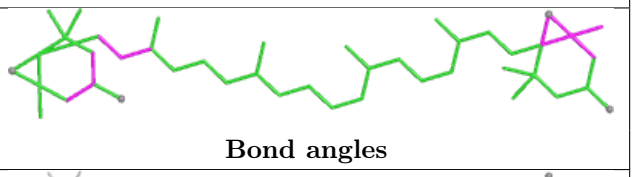
Ligand CLA I 608	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA F 608	
	
Bond lengths	Bond angles
	
Torsions	Rings

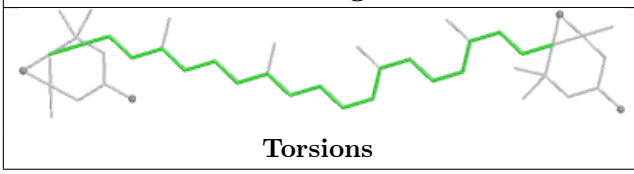
Ligand XAT G 612



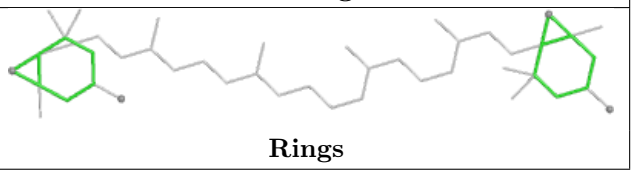
Bond lengths



Bond angles

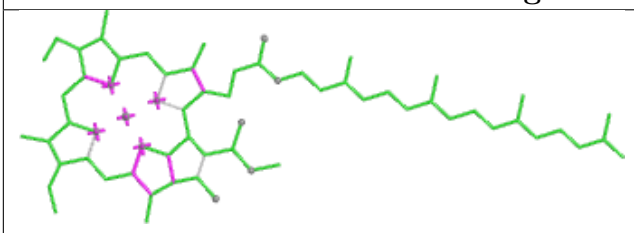


Torsions

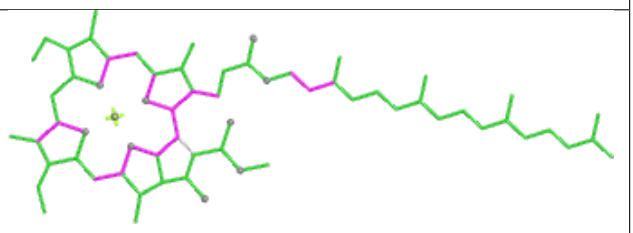


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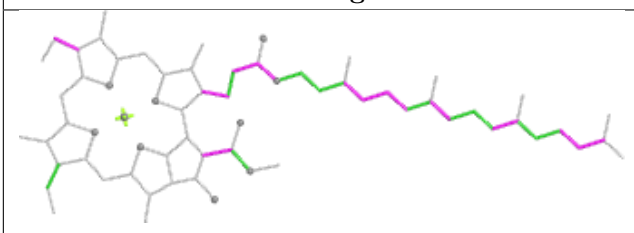
Ligand CLA A 605



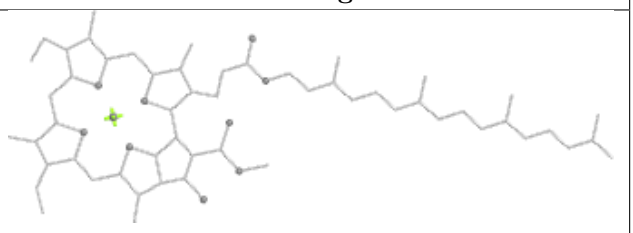
Bond lengths



Bond angles

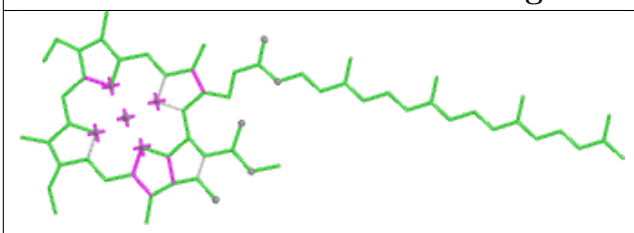


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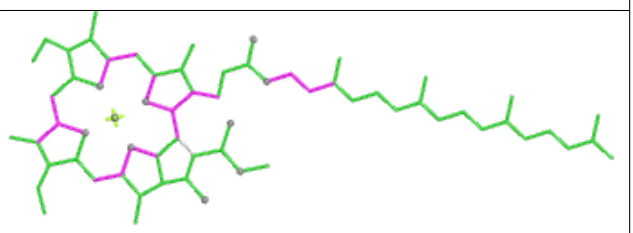


Rings

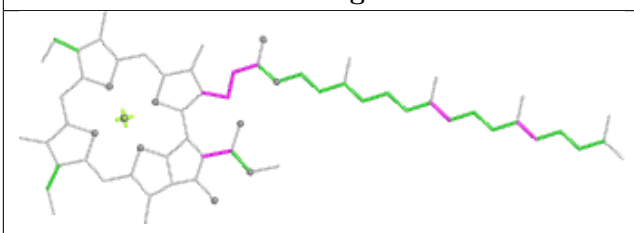
Ligand CLA E 602



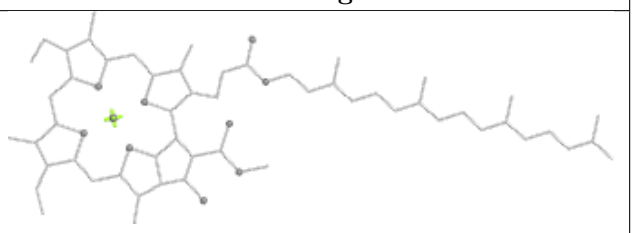
Bond lengths



Bond angles

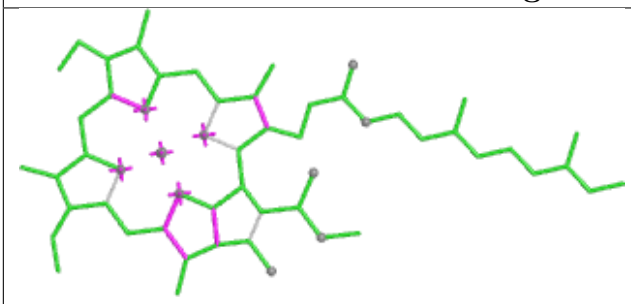


Torsions

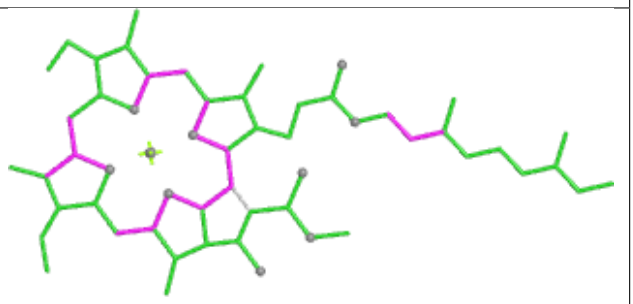


Rings

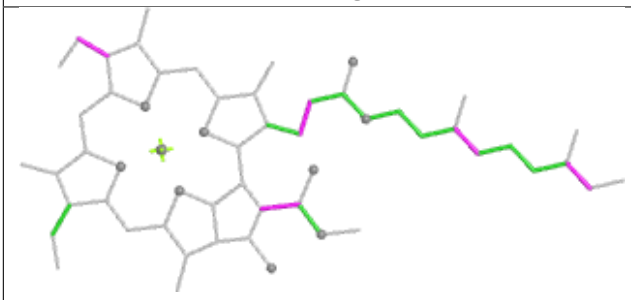
Ligand CLA H 611



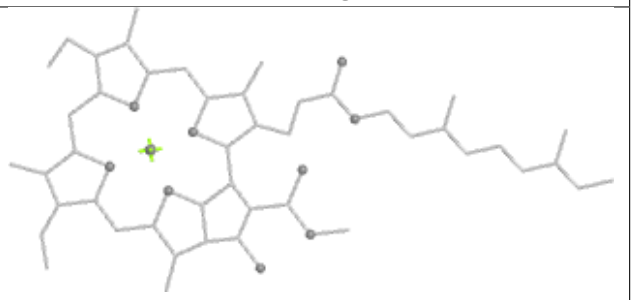
Bond lengths



Bond angles

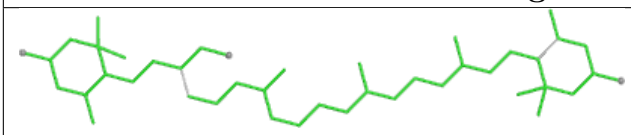


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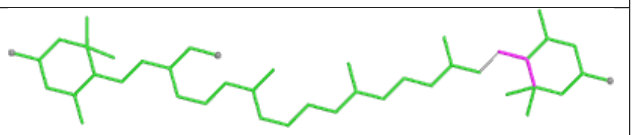


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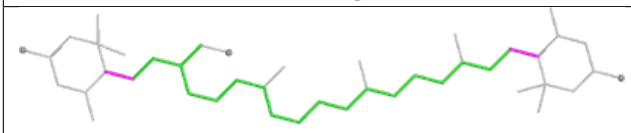
Ligand 32N G 613



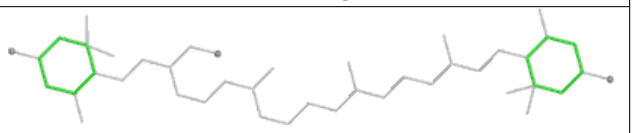
Bond lengths



Bond angles

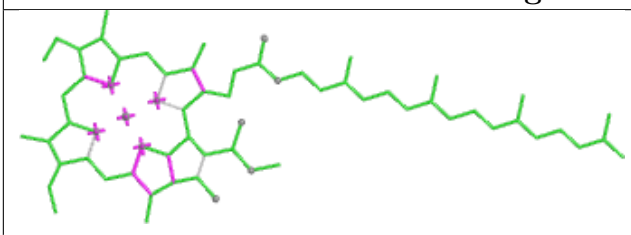


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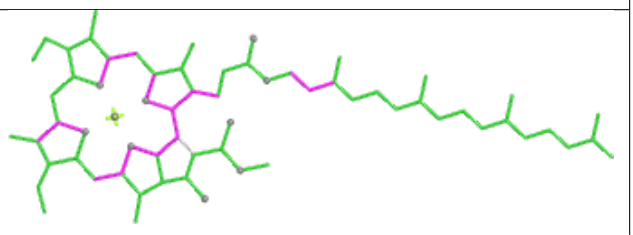


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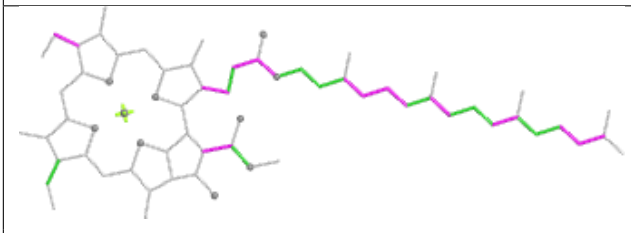
Ligand CLA B 605



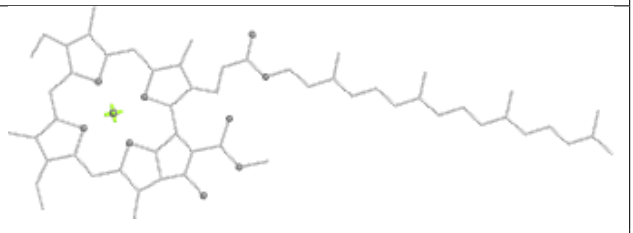
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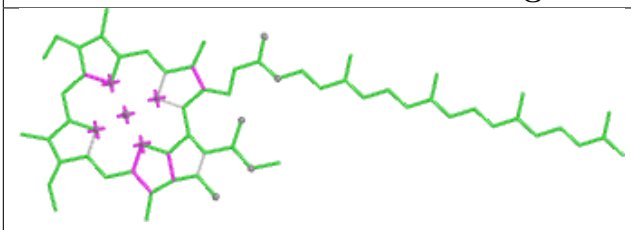
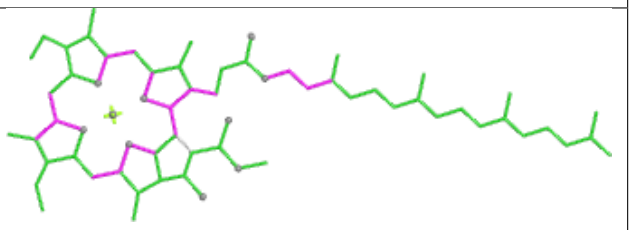
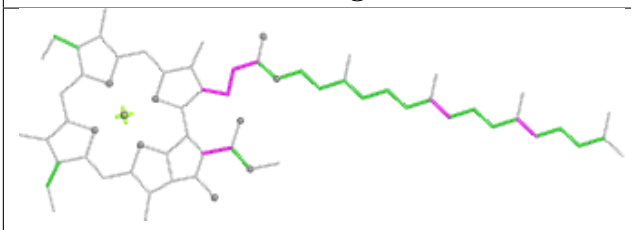
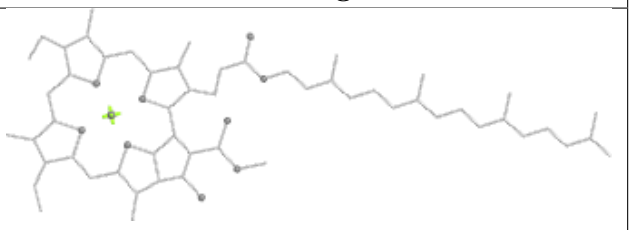
Bond angles

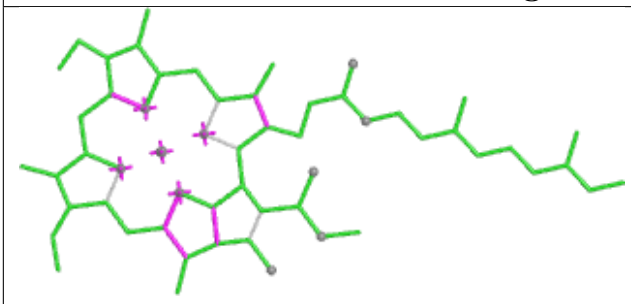
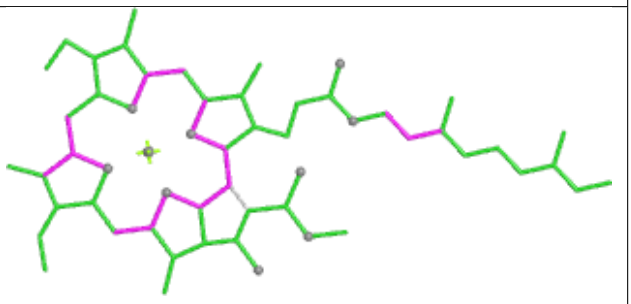
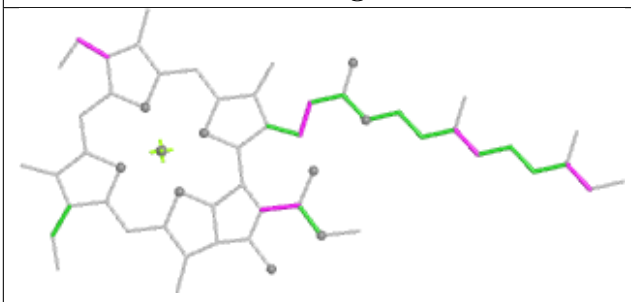
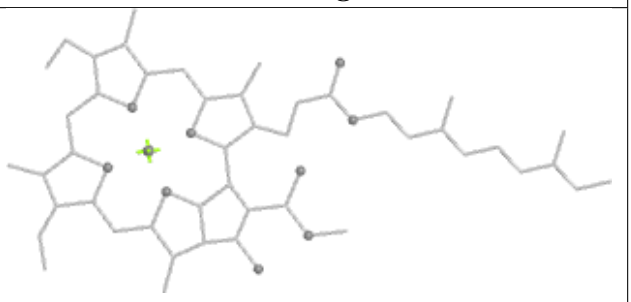


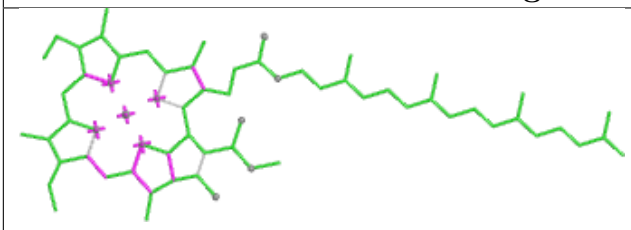
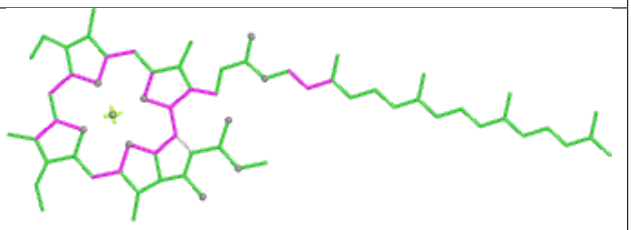
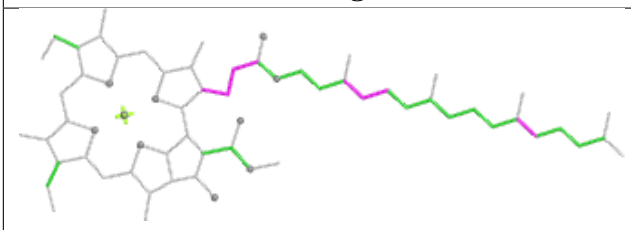
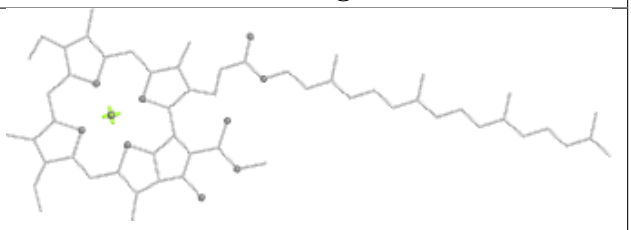
Torsions

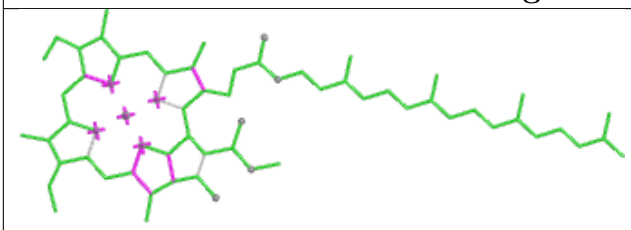
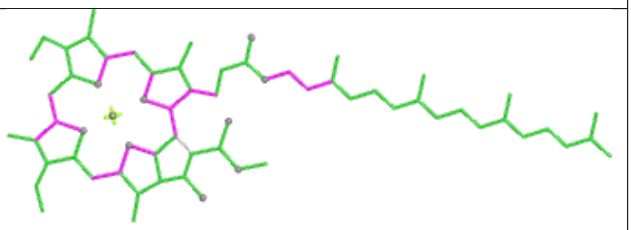
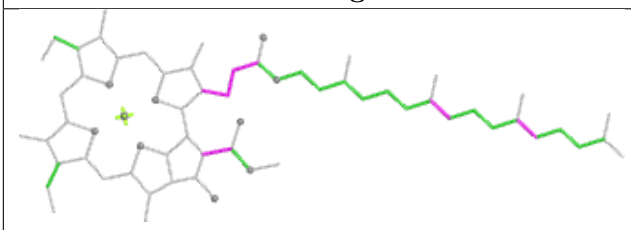
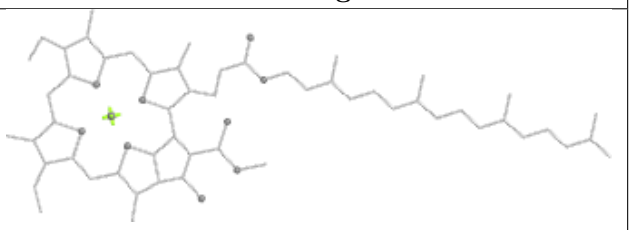


Rings

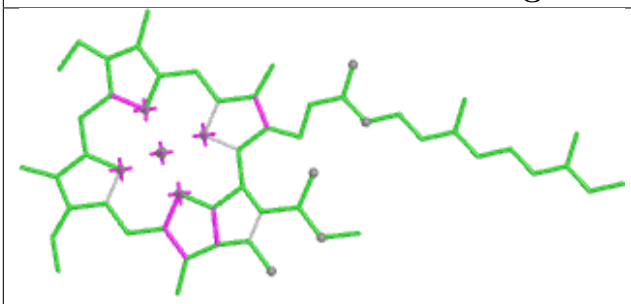
Ligand CLA C 602	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand CLA G 611	
 Bond lengths	 Bond angles
 Torsions	 Rings

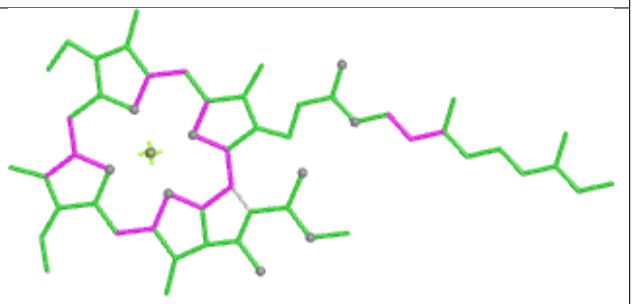
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Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA G 602	
	
Bond lengths	Bond angles
	
Torsions	Rings

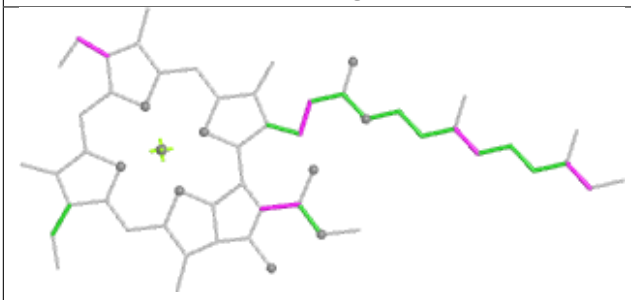
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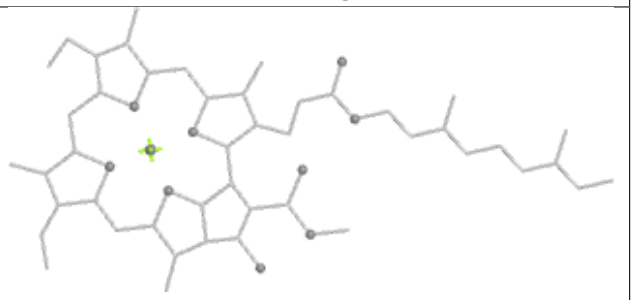
Bond lengths



Bond angles

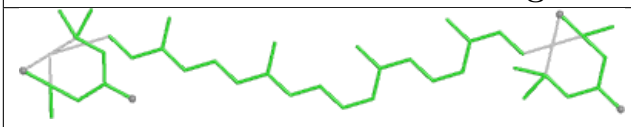


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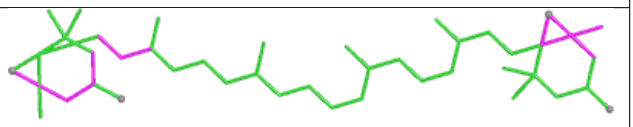


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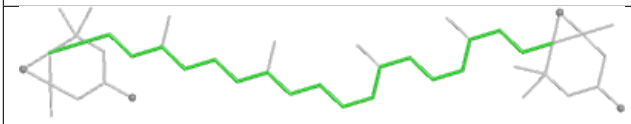
Ligand XAT J 612



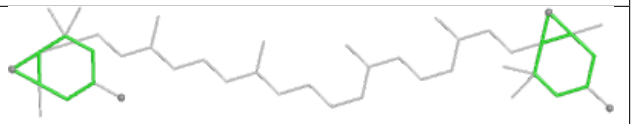
Bond lengths



Bond angles

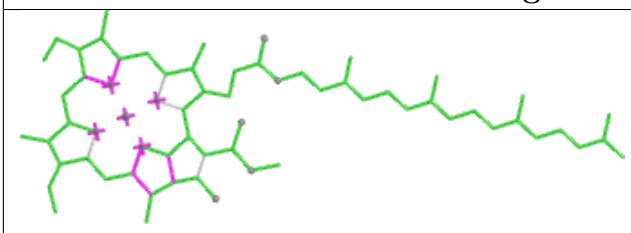


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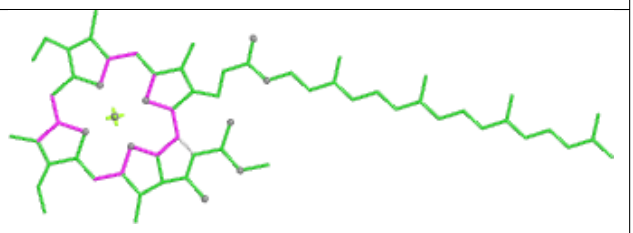


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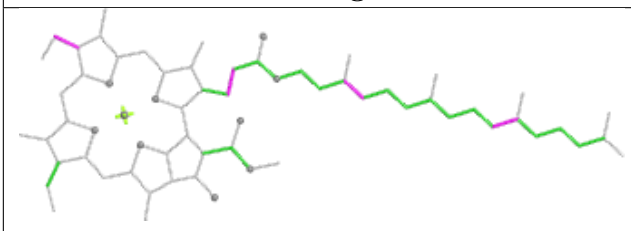
Ligand CLA G 604



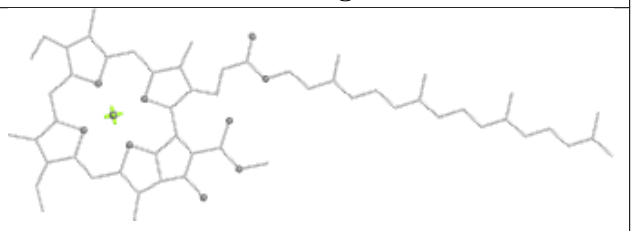
Bond lengths



Bond angles

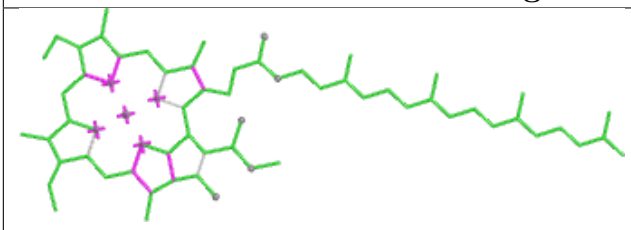


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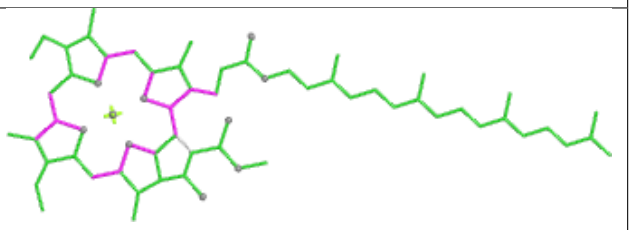


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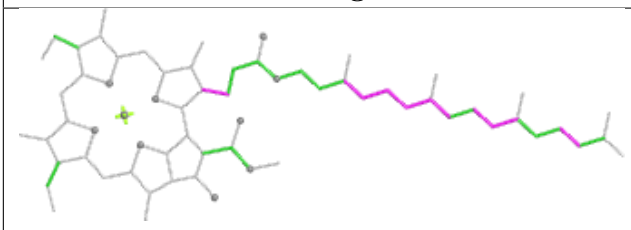
Ligand CLA D 601



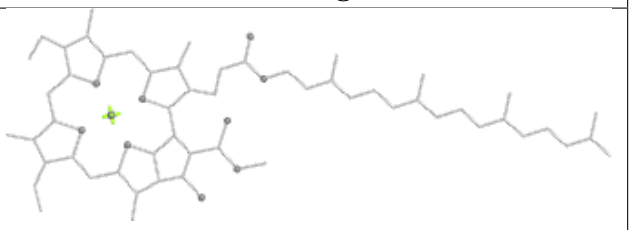
Bond lengths



Bond angles

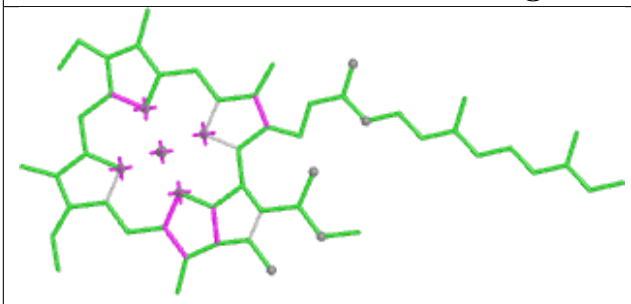


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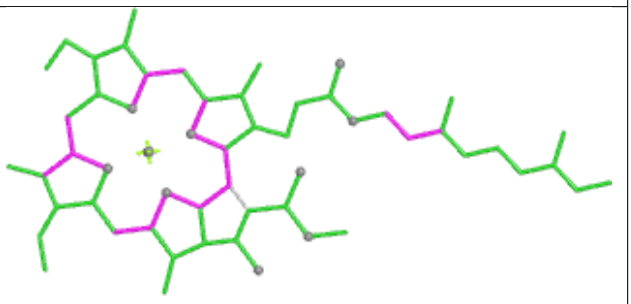


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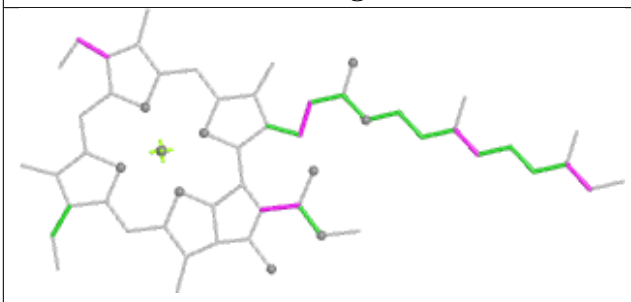
Ligand CLA B 611



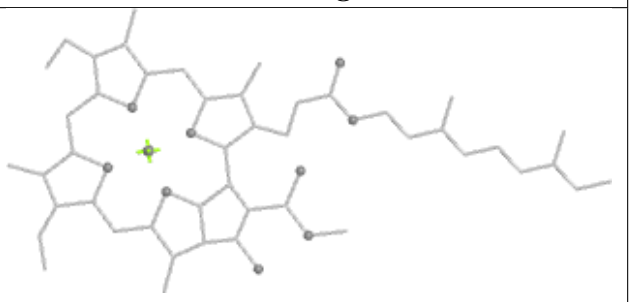
Bond lengths



Bond angles

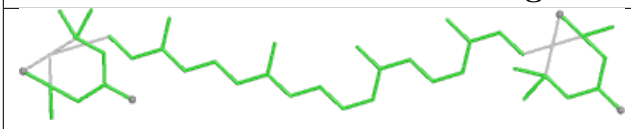


Torsions



Rings

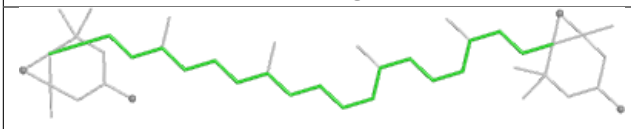
Ligand XAT F 612



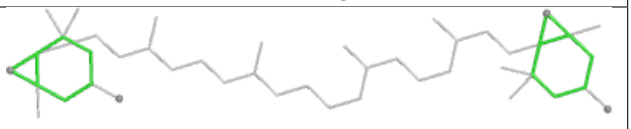
Bond lengths



Bond angles

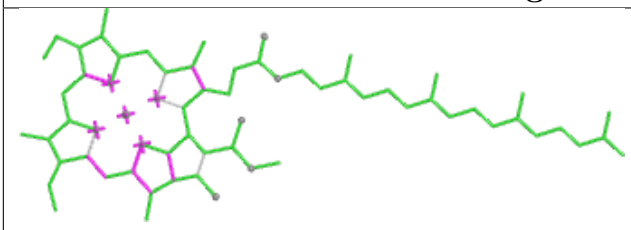


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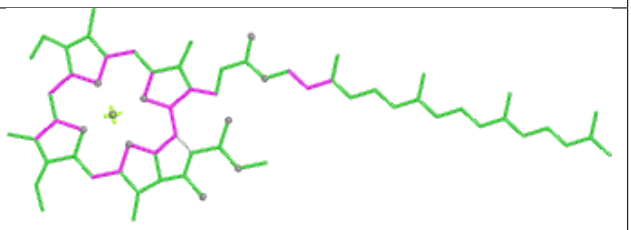


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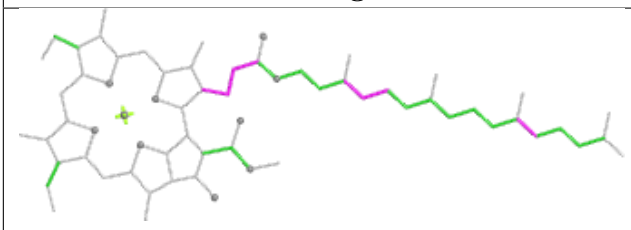
Ligand CLA F 606



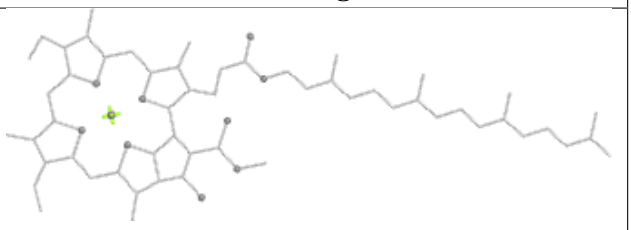
Bond lengths



Bond angles

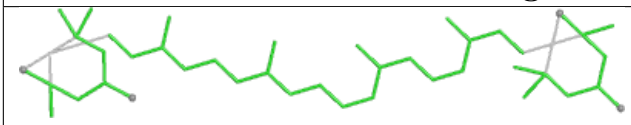


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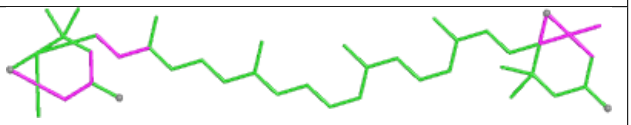


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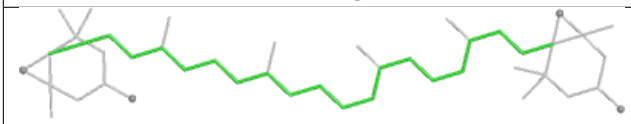
Ligand XAT K 612



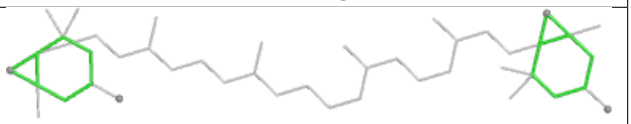
Bond lengths



Bond angles

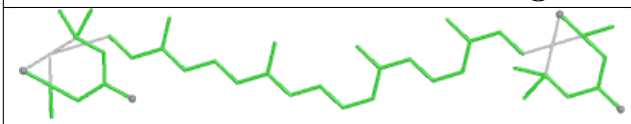


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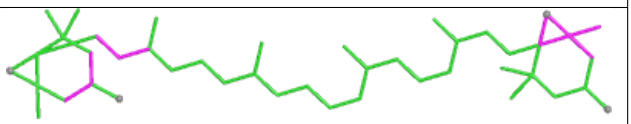


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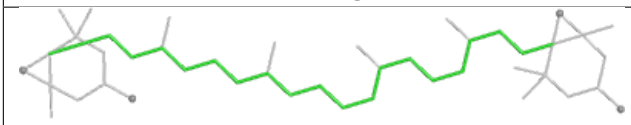
Ligand XAT I 612



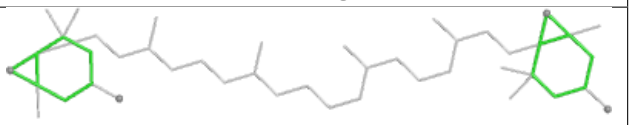
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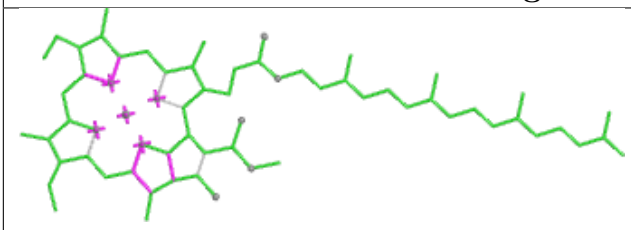


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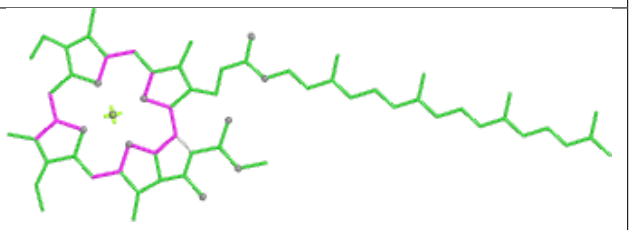


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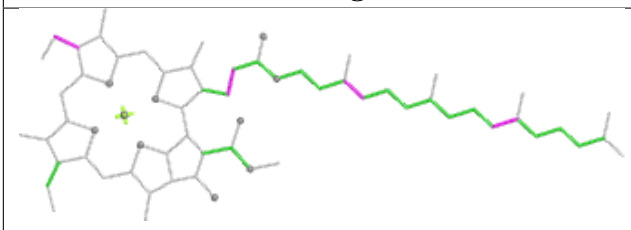
Ligand CLA D 604



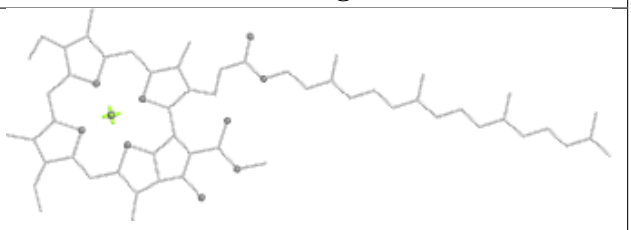
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Bond angles

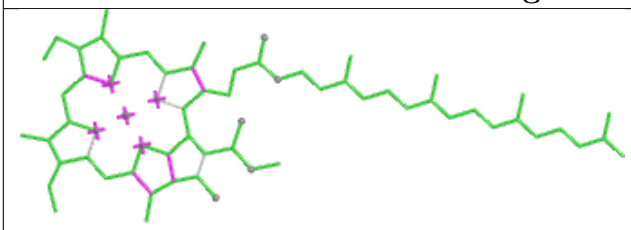


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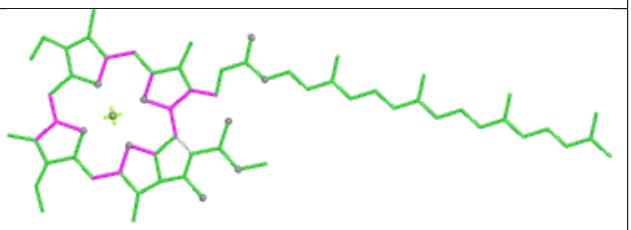


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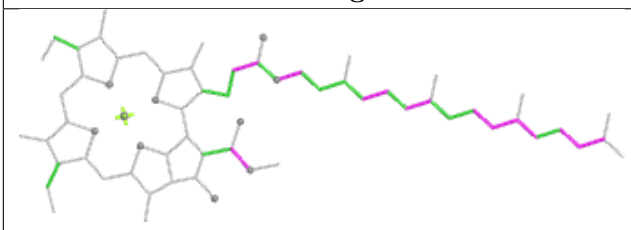
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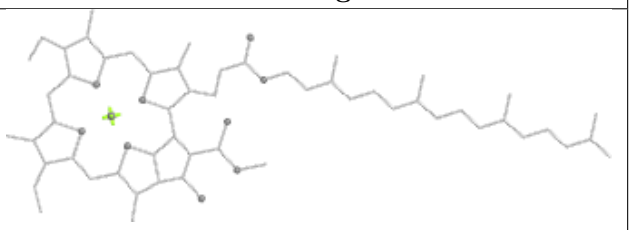
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Bond angles

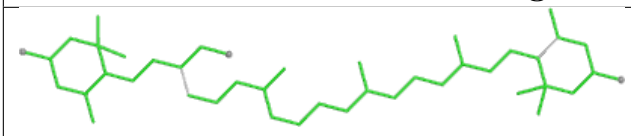


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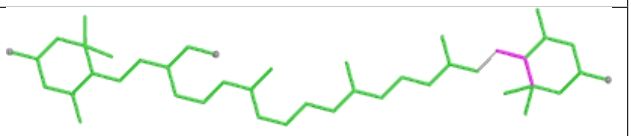


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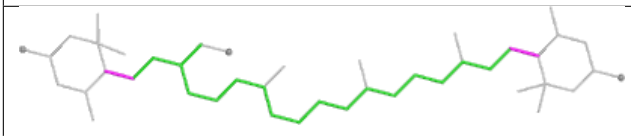
Ligand 32N K 613



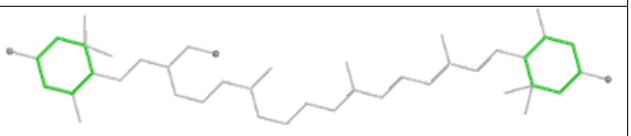
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Bond angles

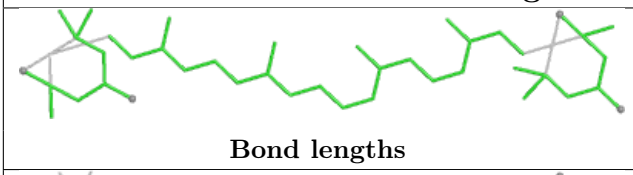


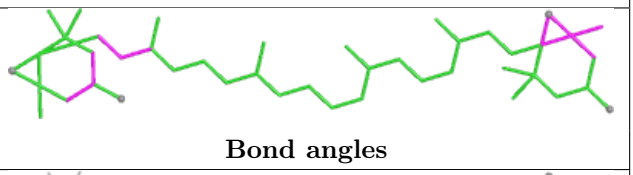
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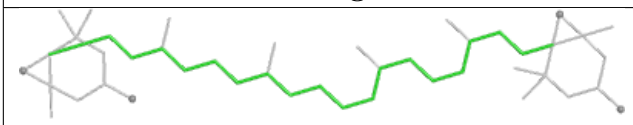


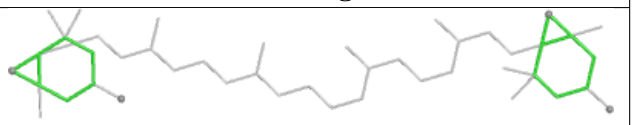
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Ligand XAT A 612

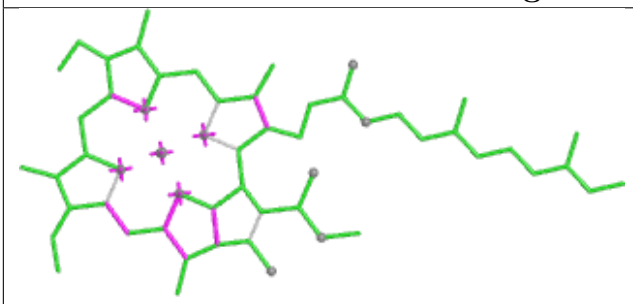

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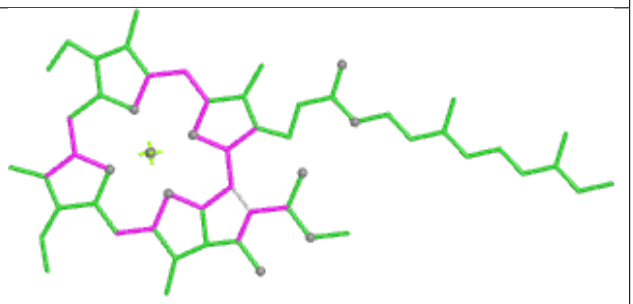

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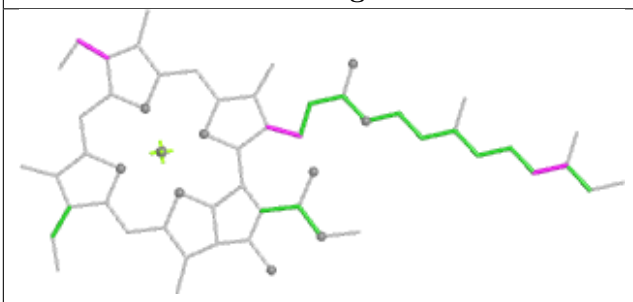

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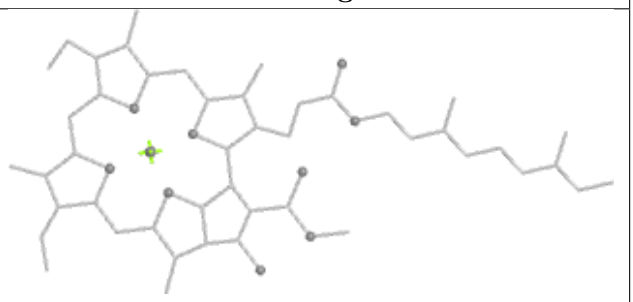

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Ligand CLA H 608

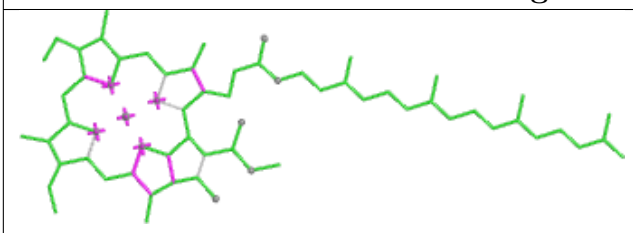

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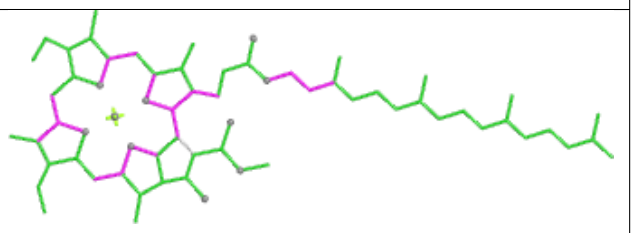

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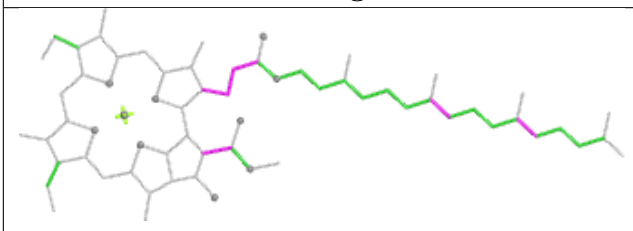

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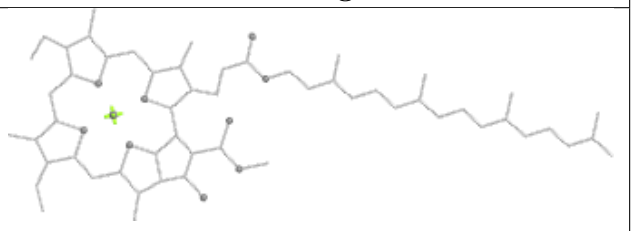

Rings

Ligand CLA I 602

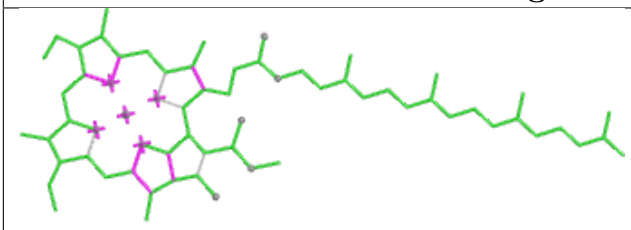

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Bond angles

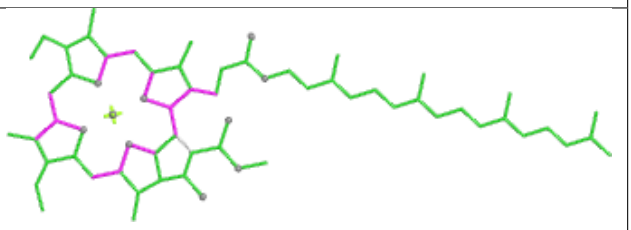

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Rings

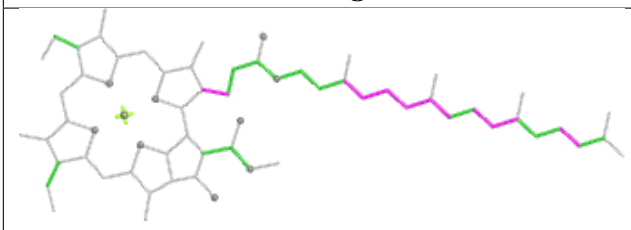
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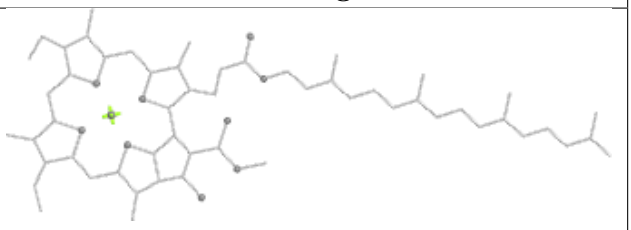
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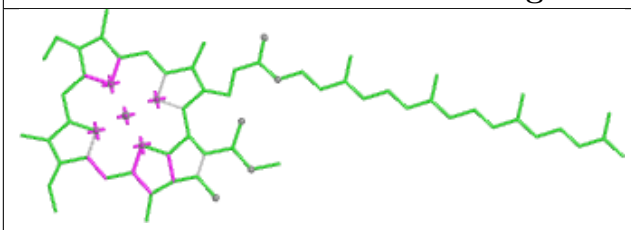


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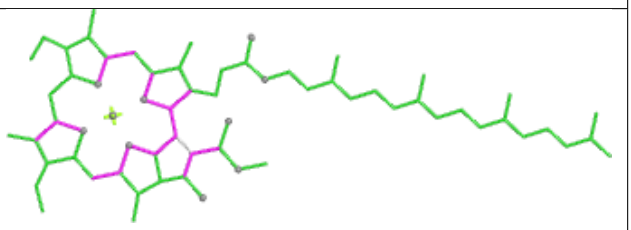


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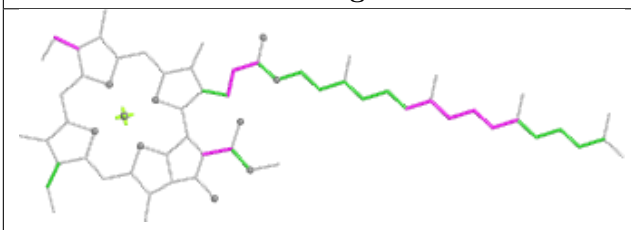
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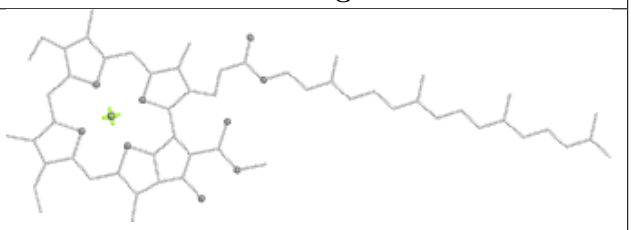
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Bond angles

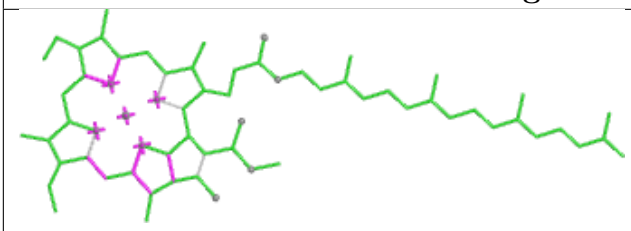


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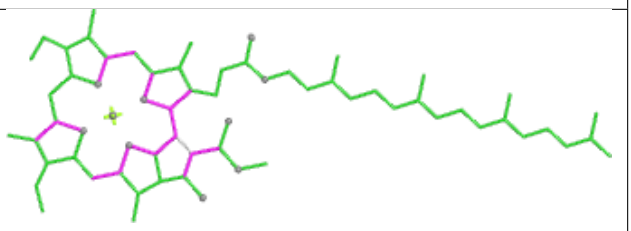


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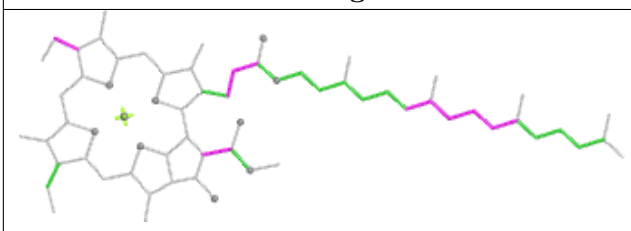
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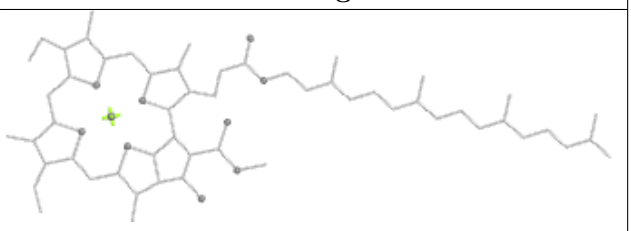
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Bond angles

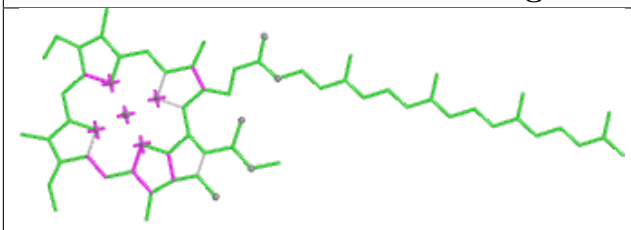


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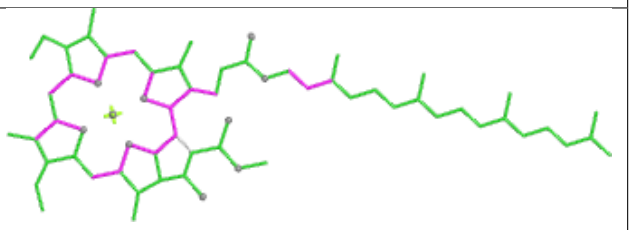


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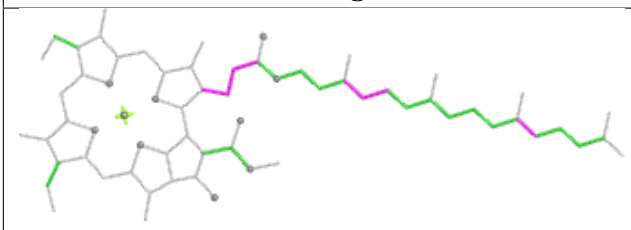
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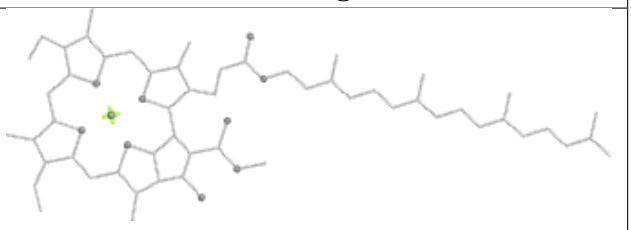
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Bond angles

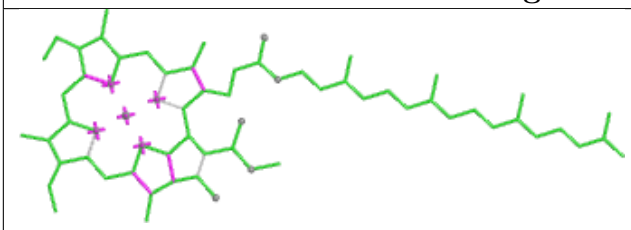


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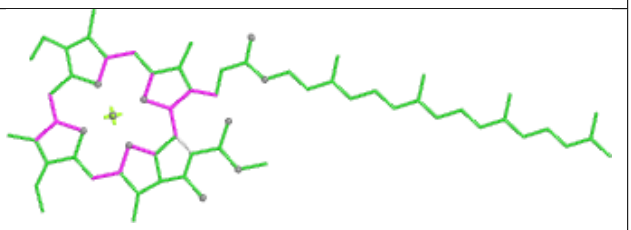


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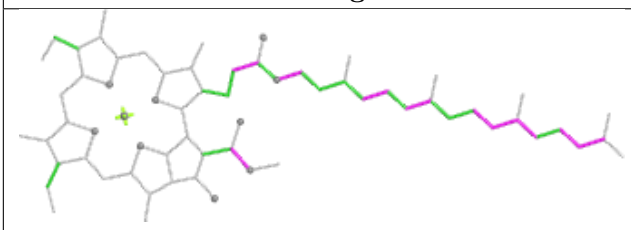
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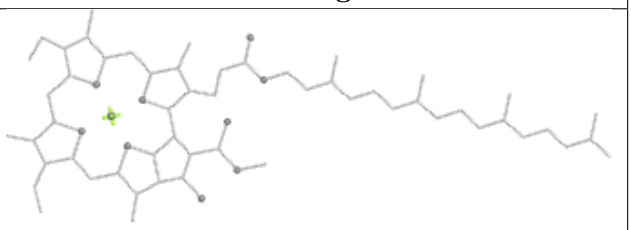
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Bond angles

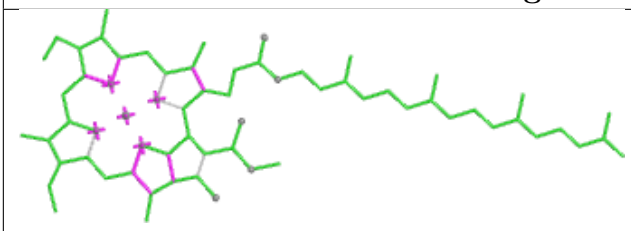


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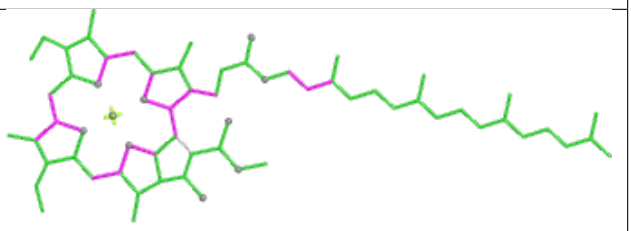


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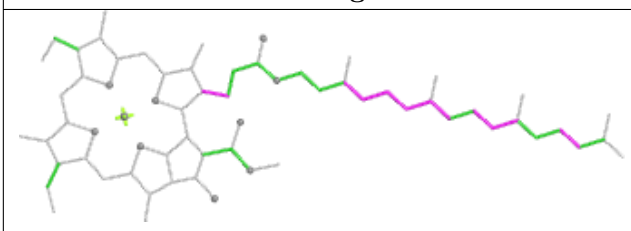
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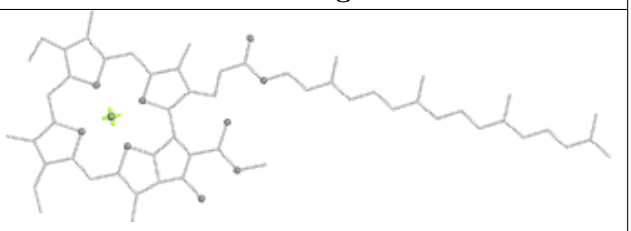
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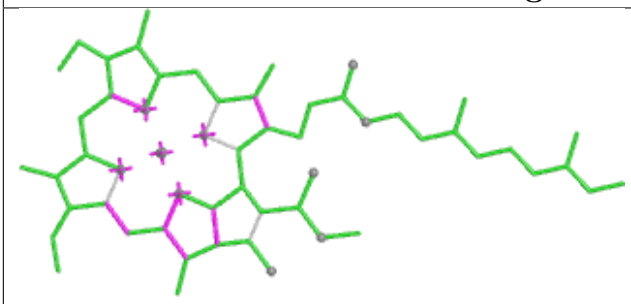


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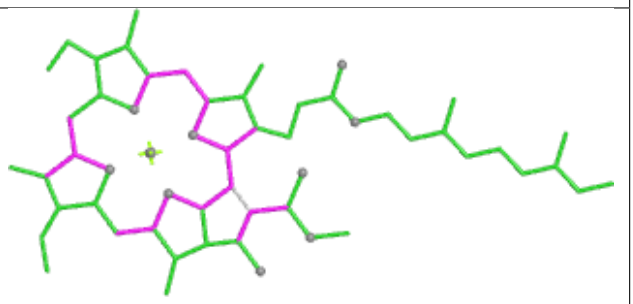


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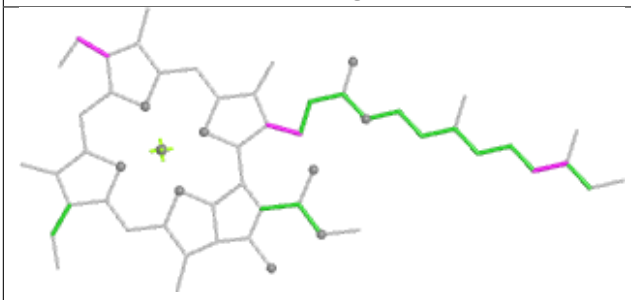
Ligand CLA E 608



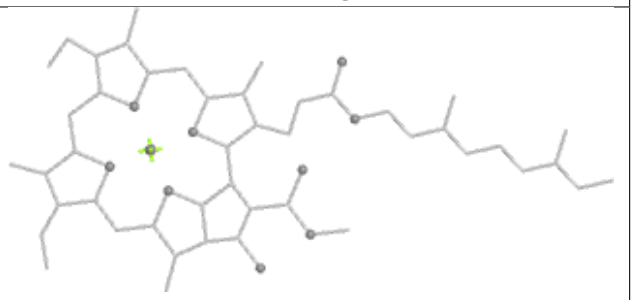
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Bond angles

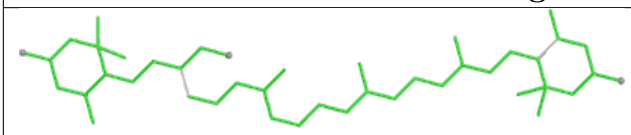


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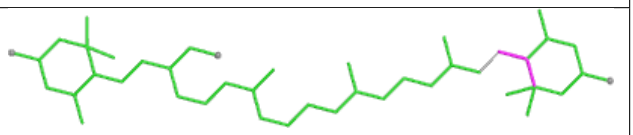


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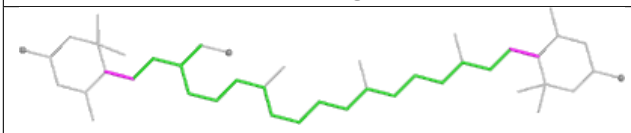
Ligand 32N D 613



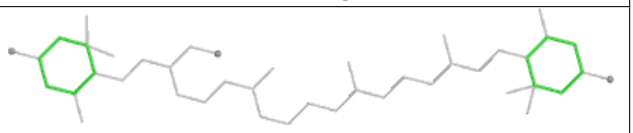
Bond lengths



Bond angles

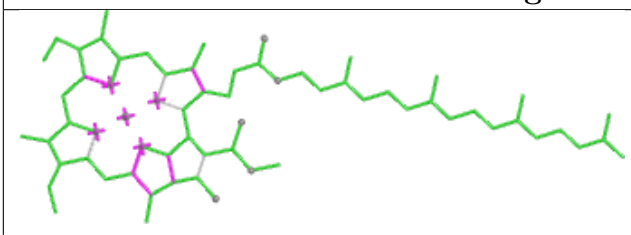


Torsions

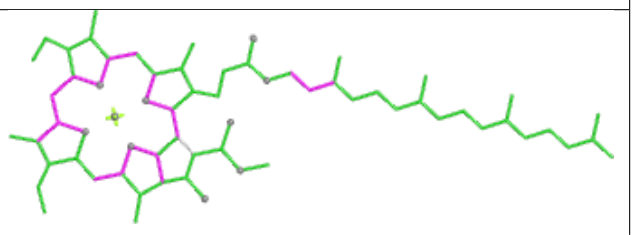


Rings

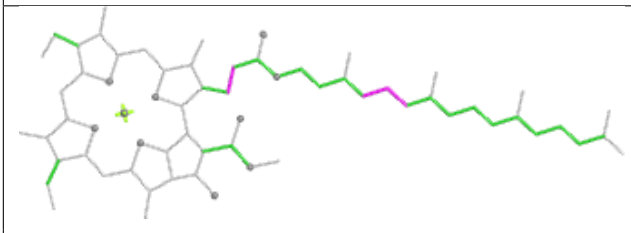
Ligand CLA C 609



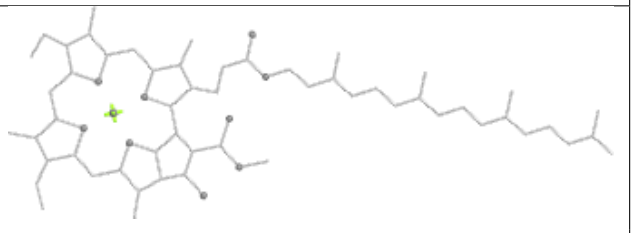
Bond lengths



Bond angles

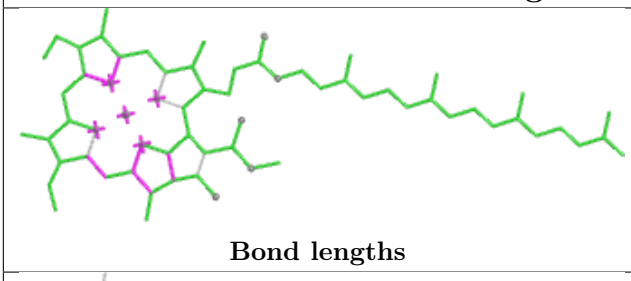


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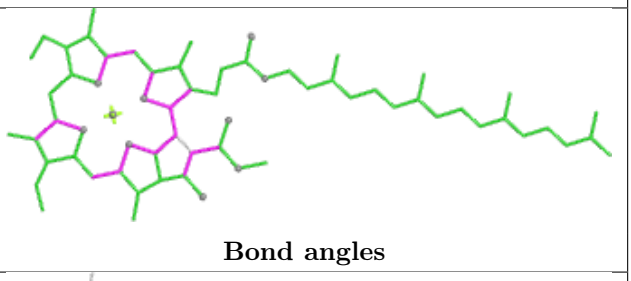


Rings

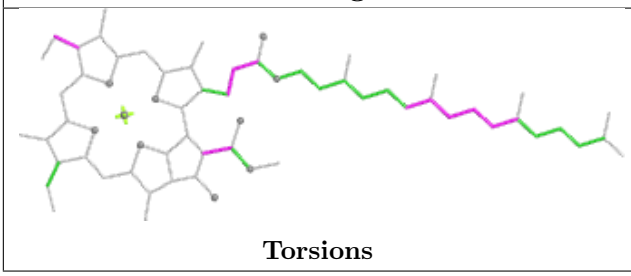
Ligand CLA E 610



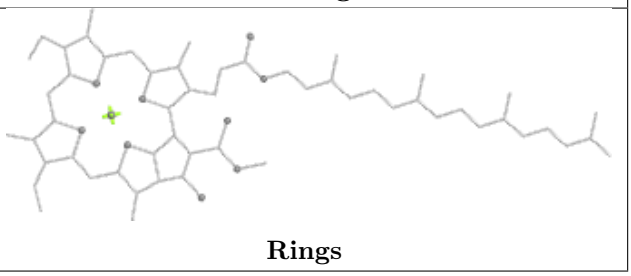
Bond lengths



Bond angles

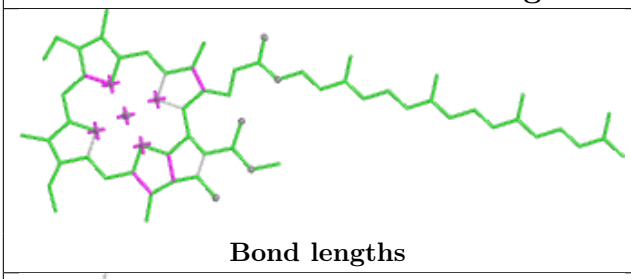


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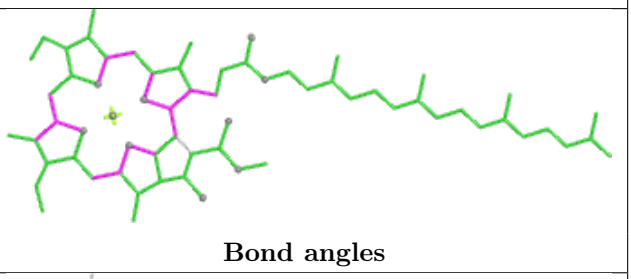


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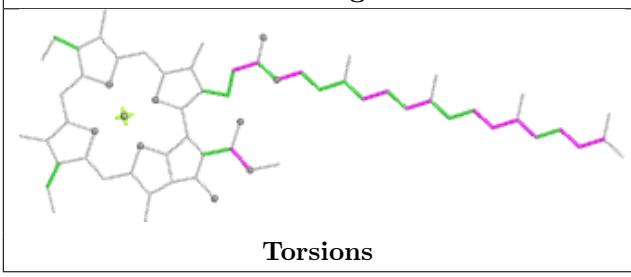
Ligand CLA A 607



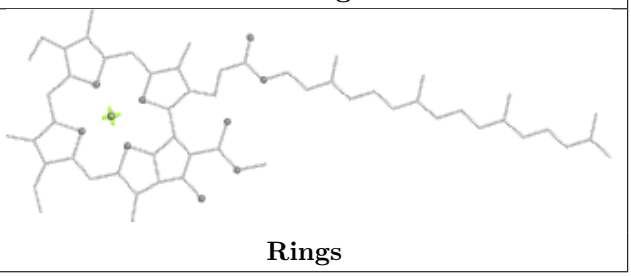
Bond lengths



Bond angles

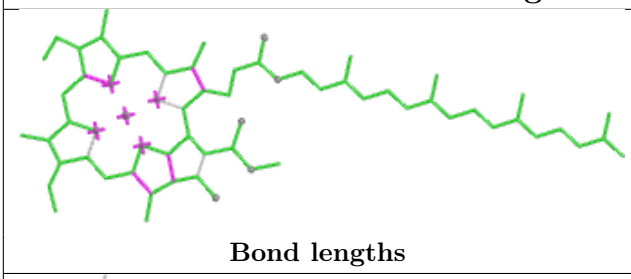


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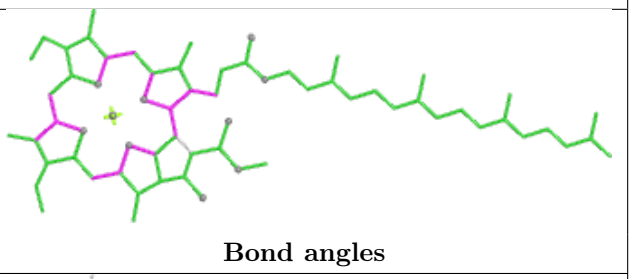


Rings

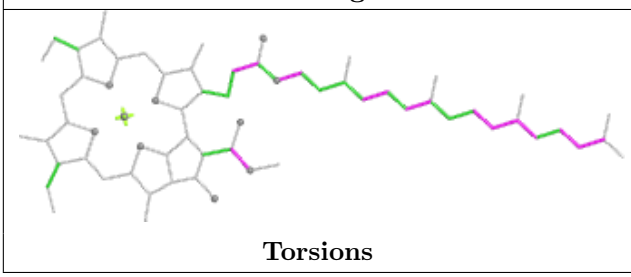
Ligand CLA B 607



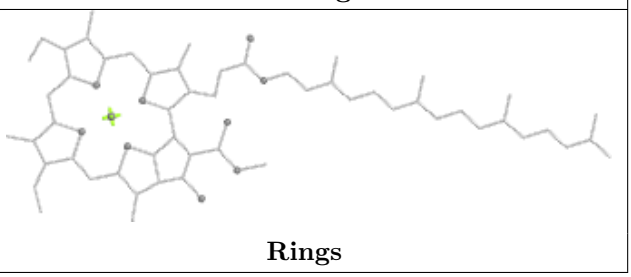
Bond lengths



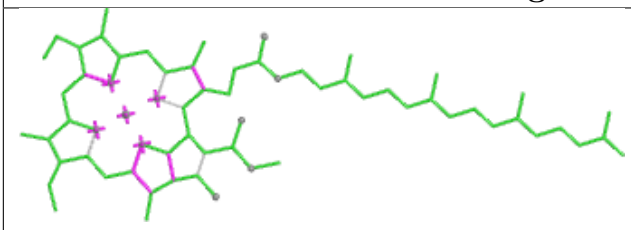
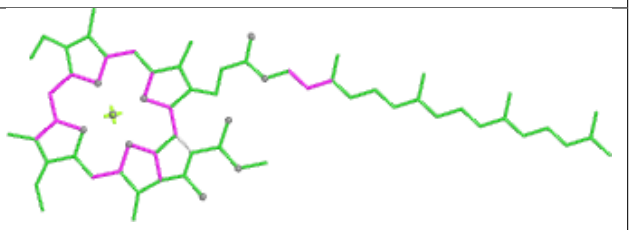
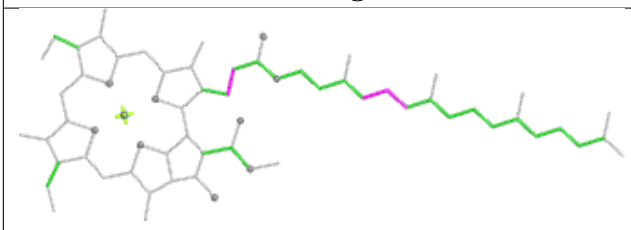
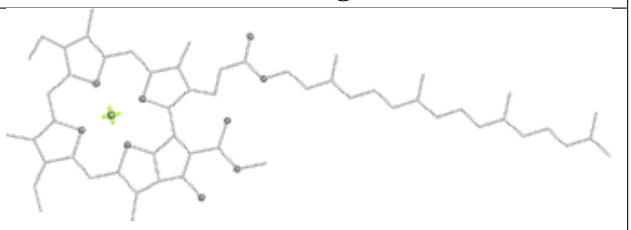
Bond angles

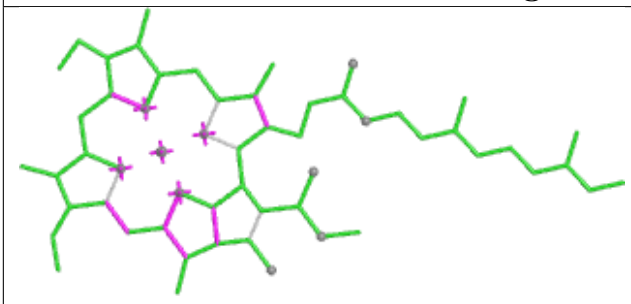
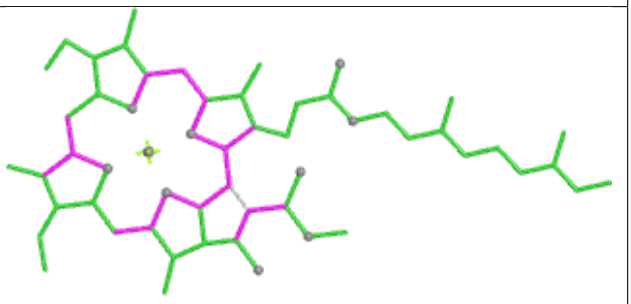
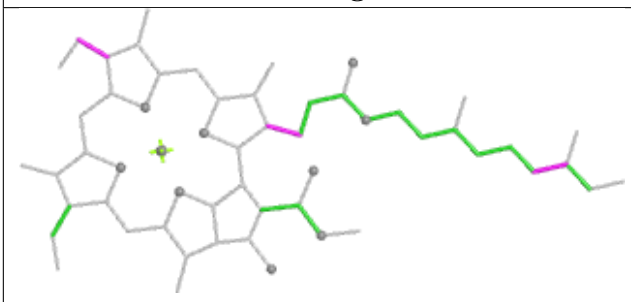
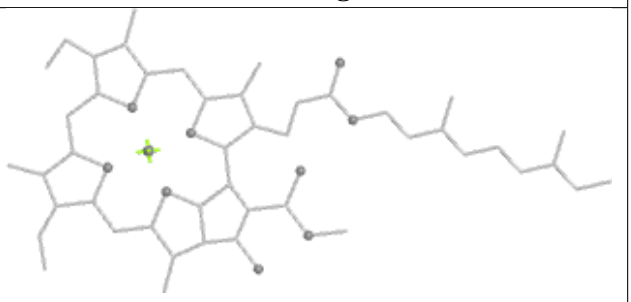


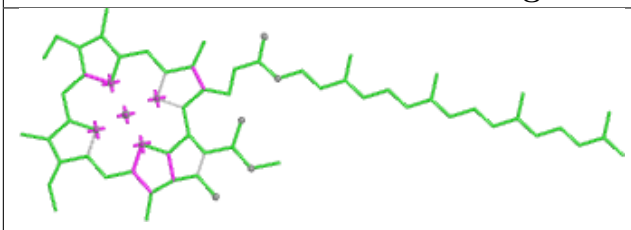
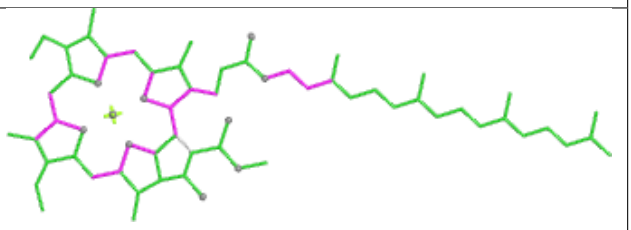
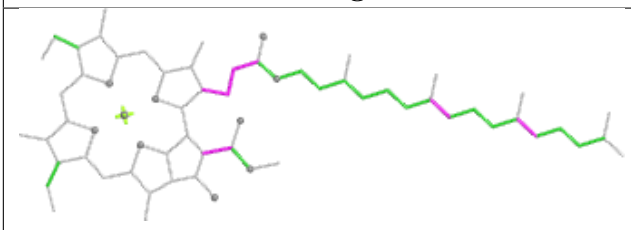
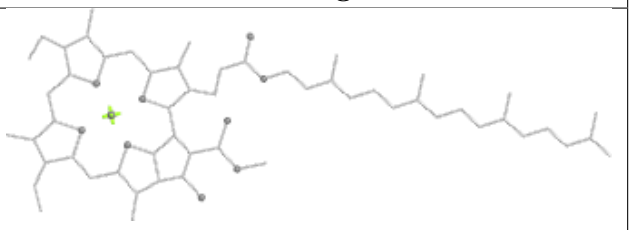
Torsions

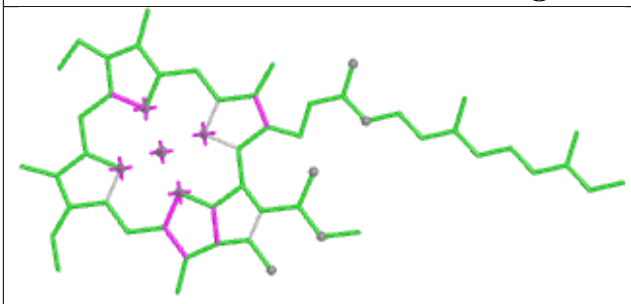
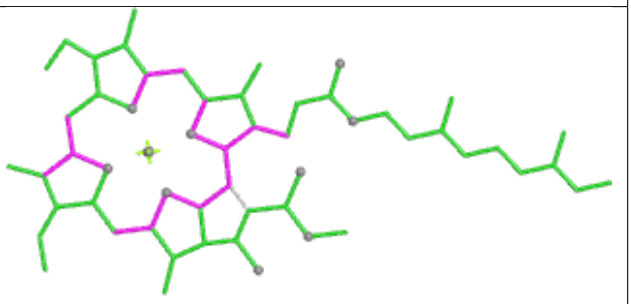
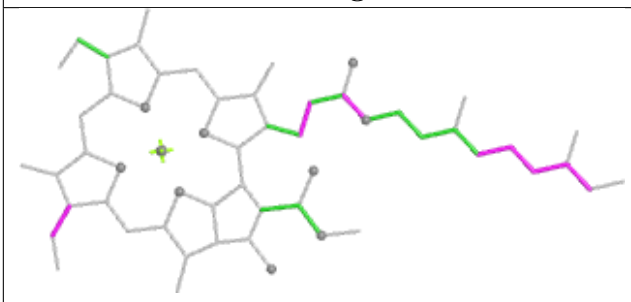
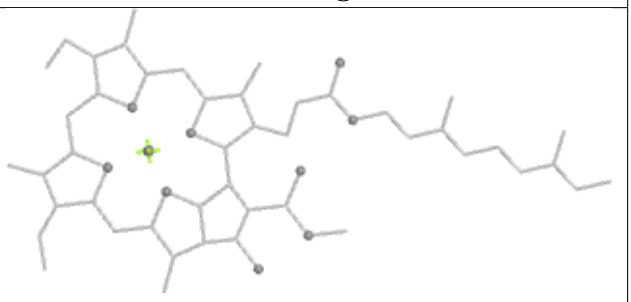


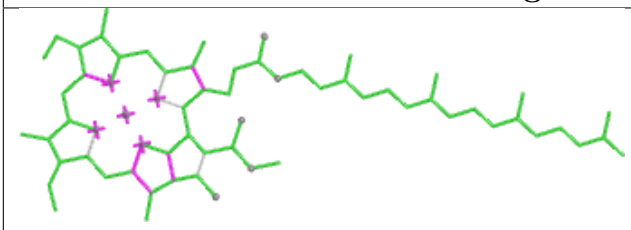
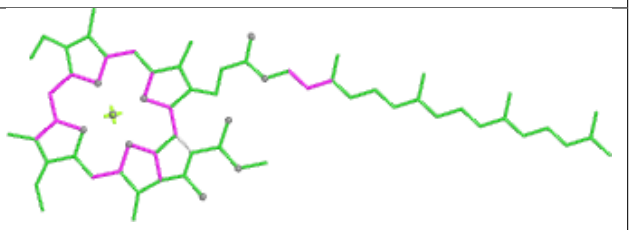
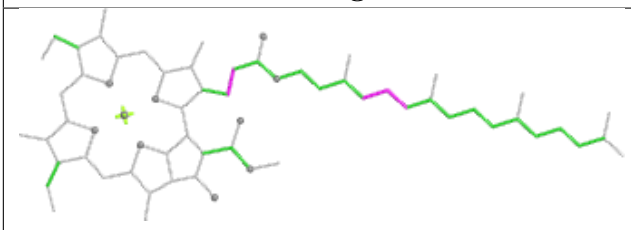
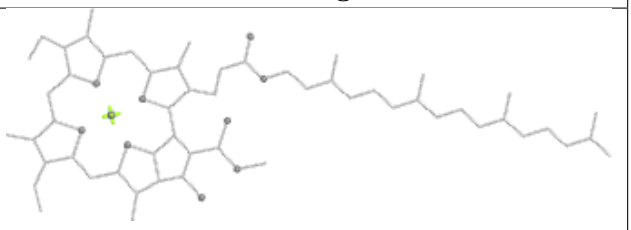
Rings

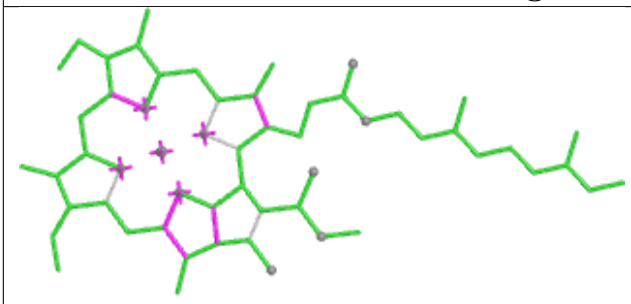
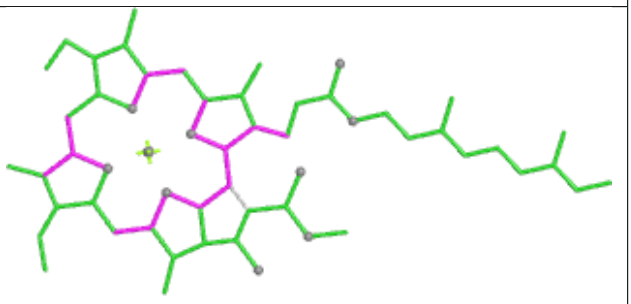
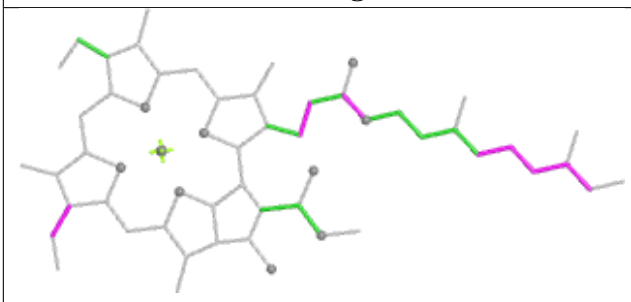
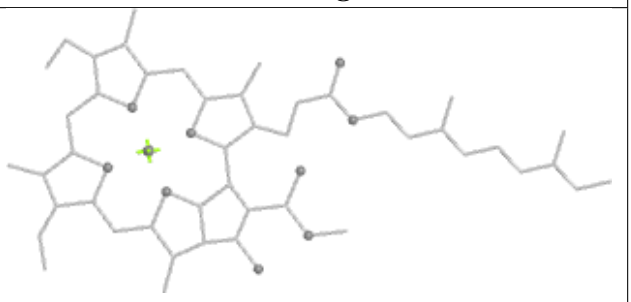
Ligand CLA J 609	
 Bond lengths	 Bond angles
 Torsions	 Rings

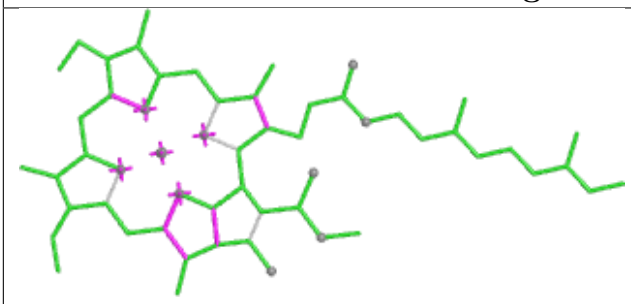
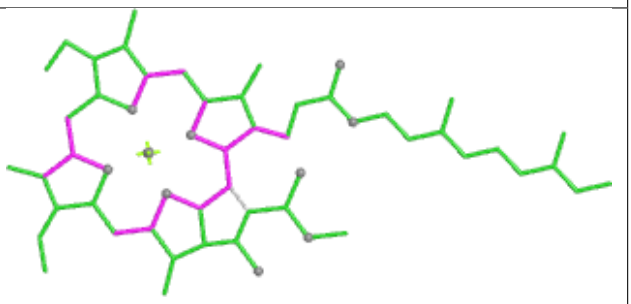
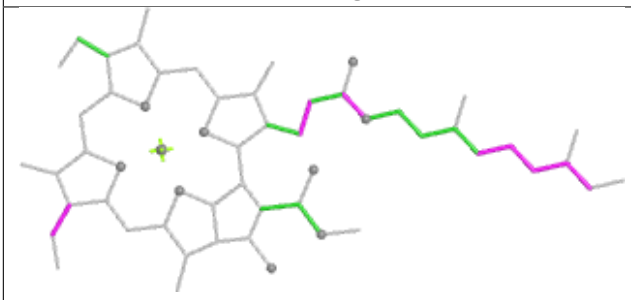
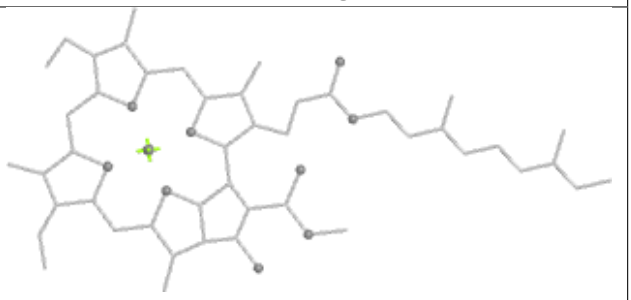
Ligand CLA K 608	
 Bond lengths	 Bond angles
 Torsions	 Rings

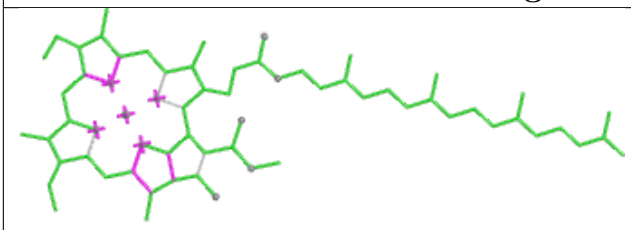
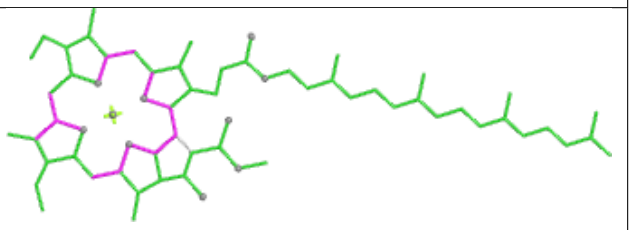
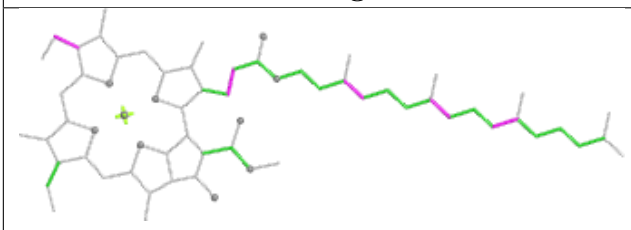
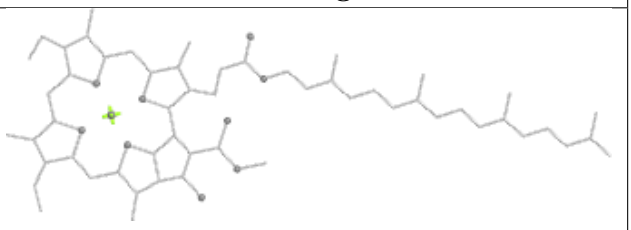
Ligand CLA K 602	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand CLA I 603	
 Bond lengths	 Bond angles
 Torsions	 Rings

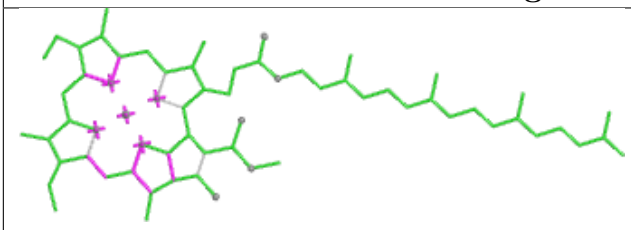
Ligand CLA A 609	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand CLA F 603	
 Bond lengths	 Bond angles
 Torsions	 Rings

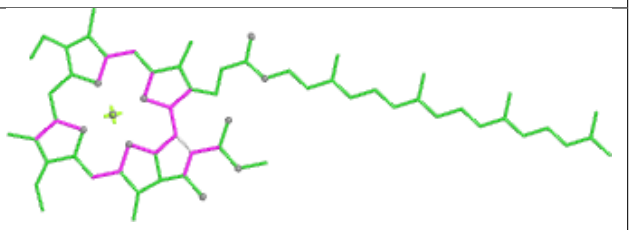
Ligand CLA K 603	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA B 604	
	
Bond lengths	Bond angles
	
Torsions	Rings

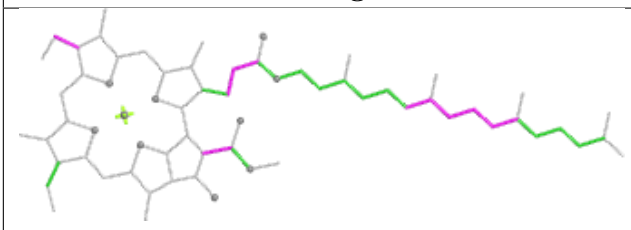
Ligand CLA C 610



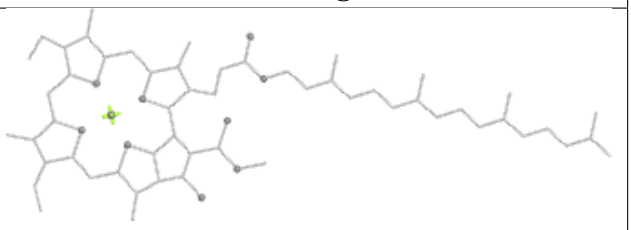
Bond lengths



Bond angles

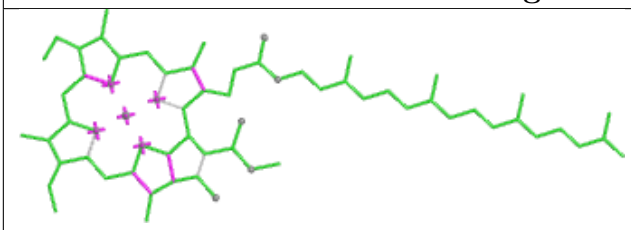


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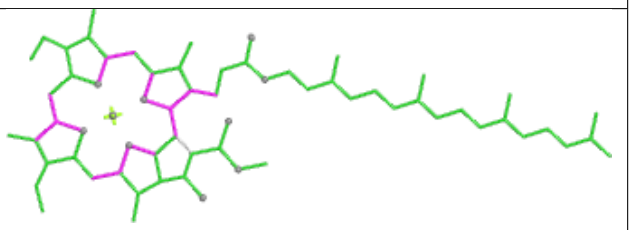


Rings

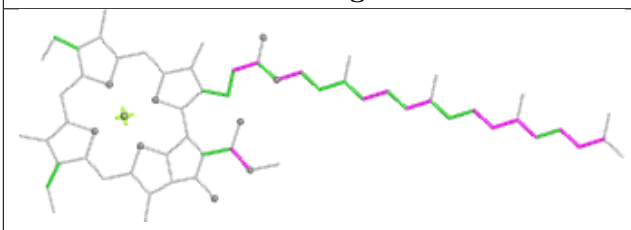
Ligand CLA D 607



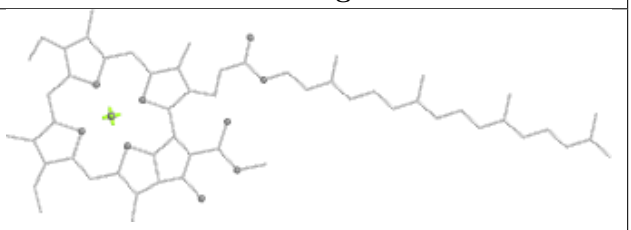
Bond lengths



Bond angles

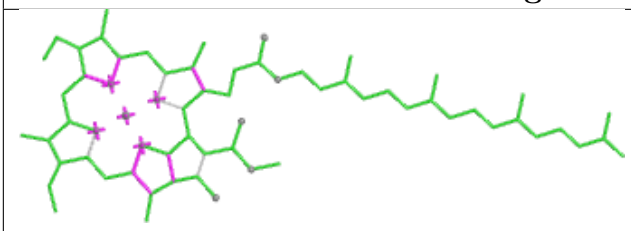


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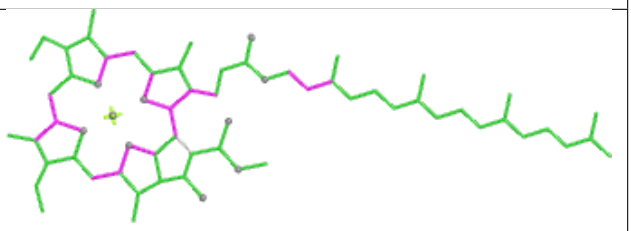


Rings

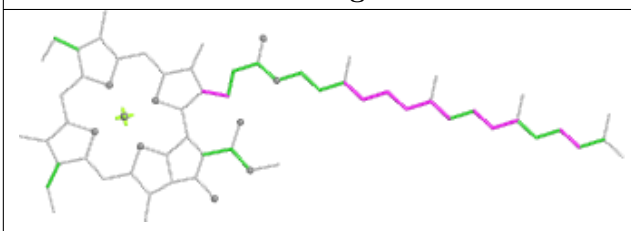
Ligand CLA C 601



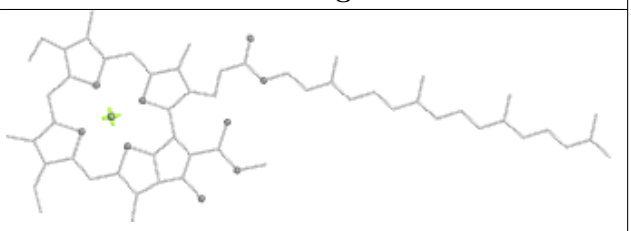
Bond lengths



Bond angles

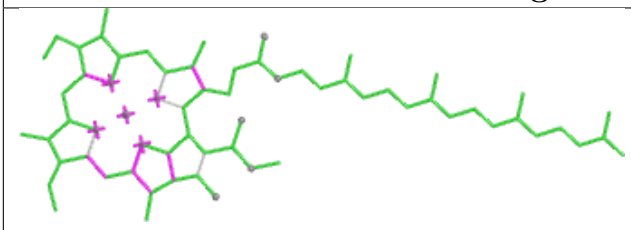


Torsions

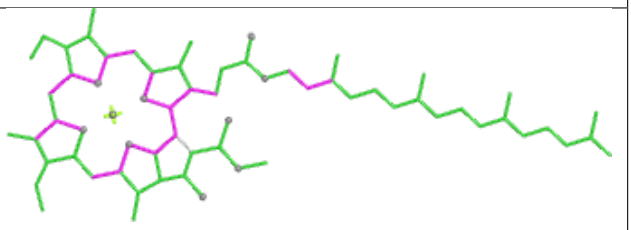


Rings

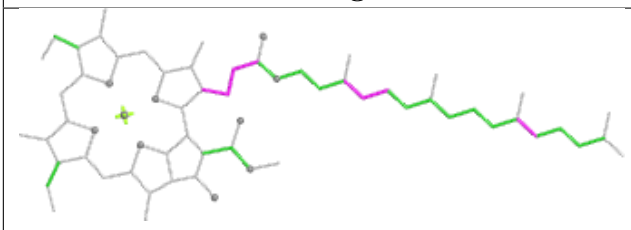
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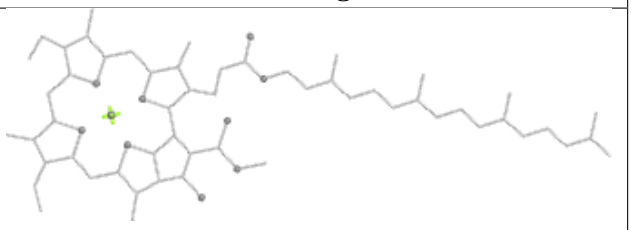
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Bond angles

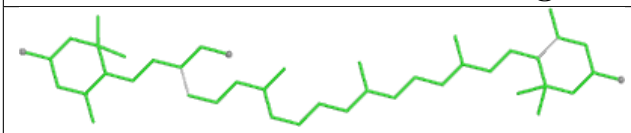


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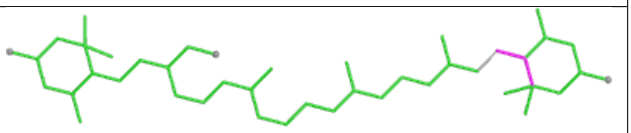


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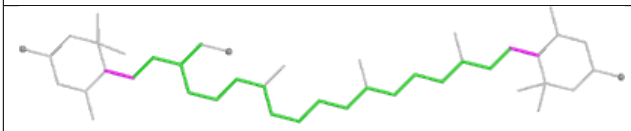
Ligand 32N I 613



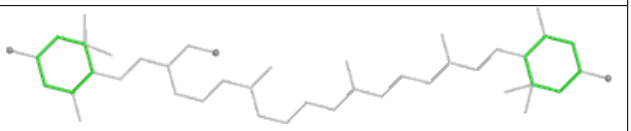
Bond lengths



Bond angles

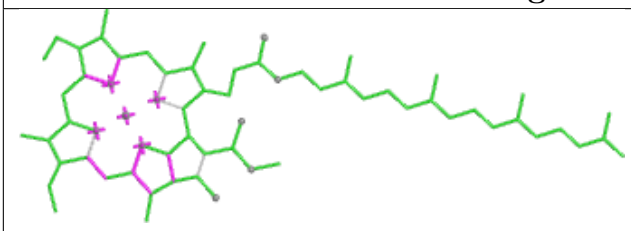


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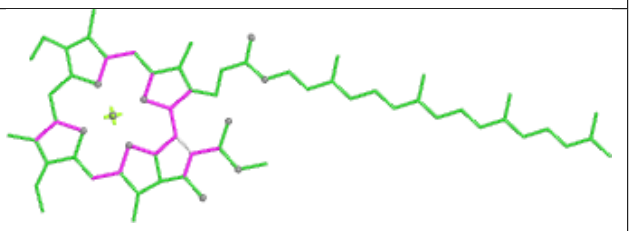


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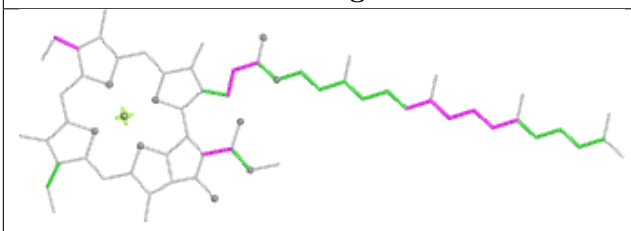
Ligand CLA J 610



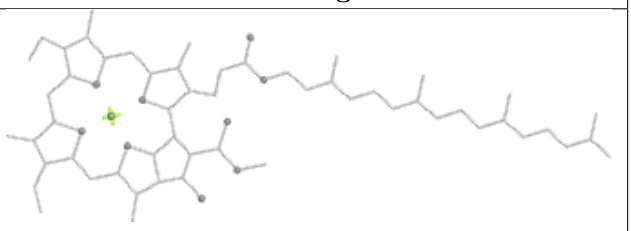
Bond lengths



Bond angles

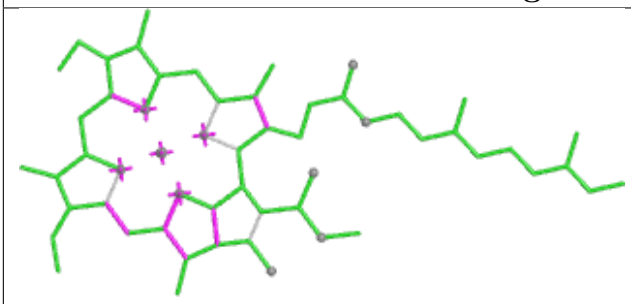


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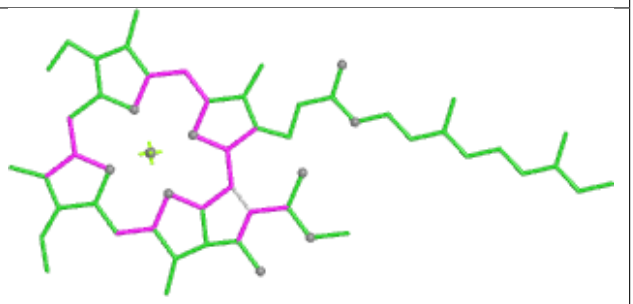


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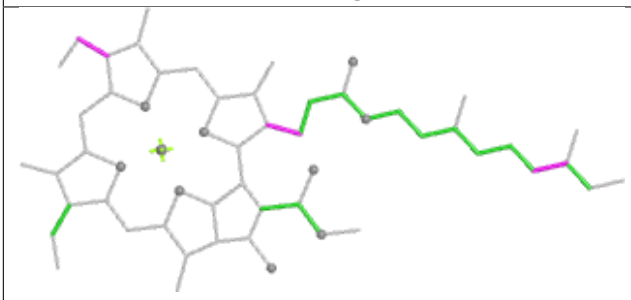
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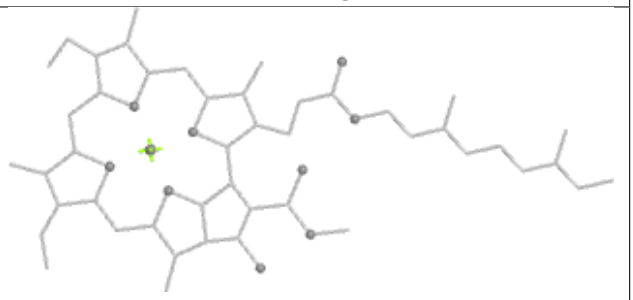
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Bond angles

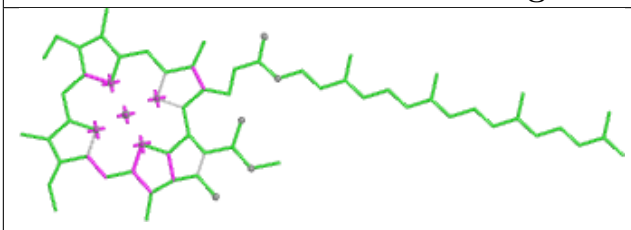


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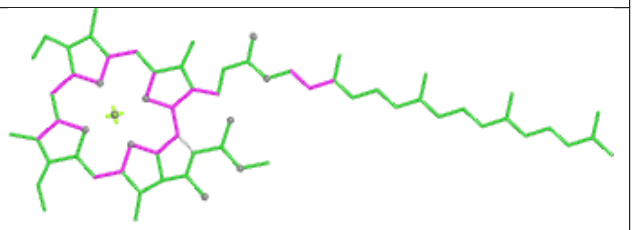


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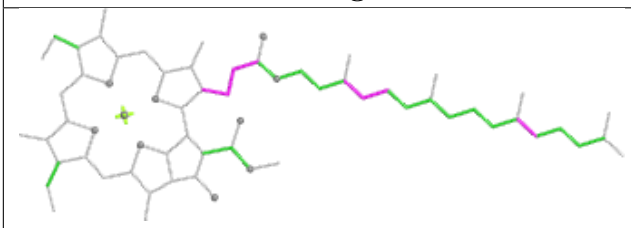
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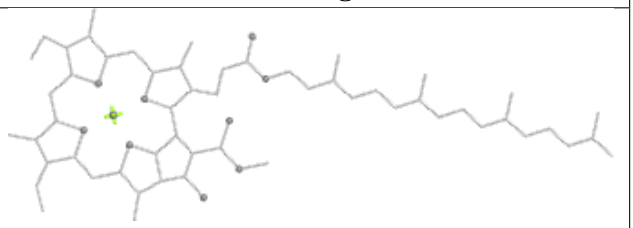
Bond lengths



Bond angles

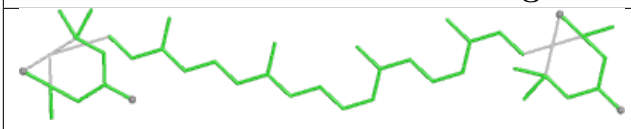


Torsions



Rings

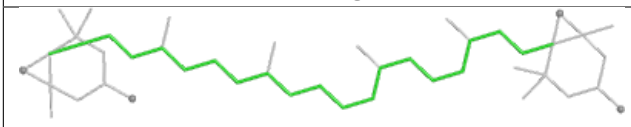
Ligand XAT H 612



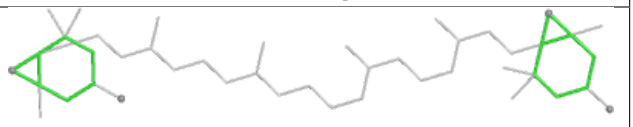
Bond lengths



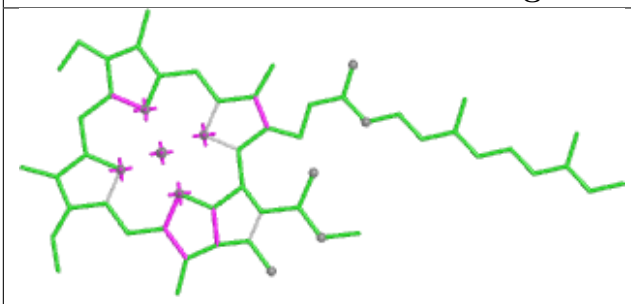
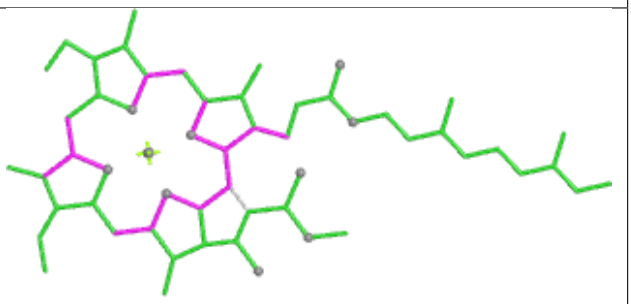
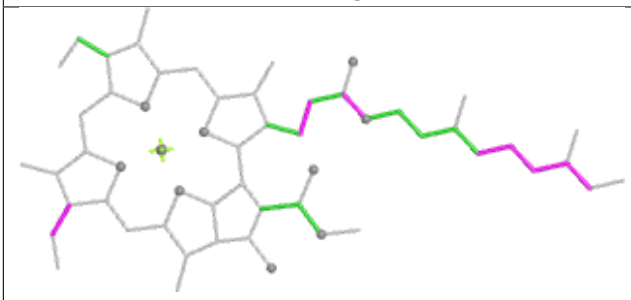
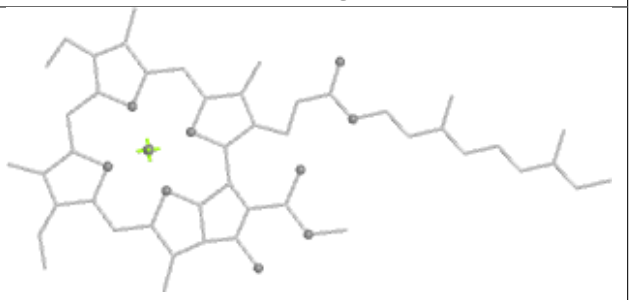
Bond angles

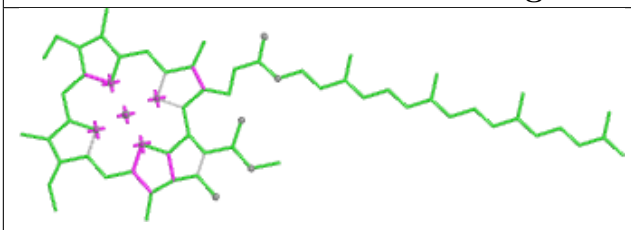
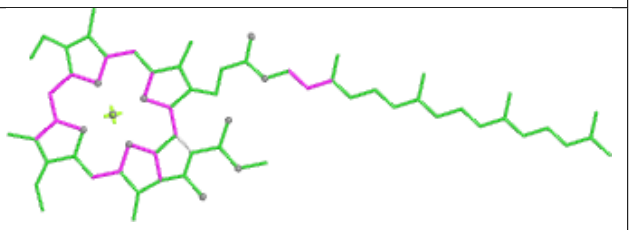
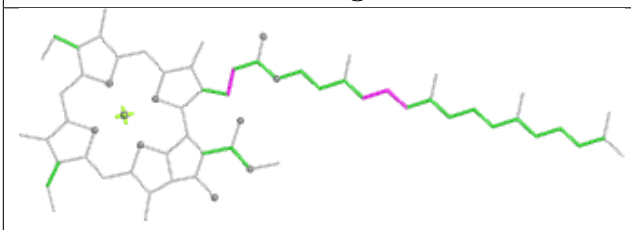
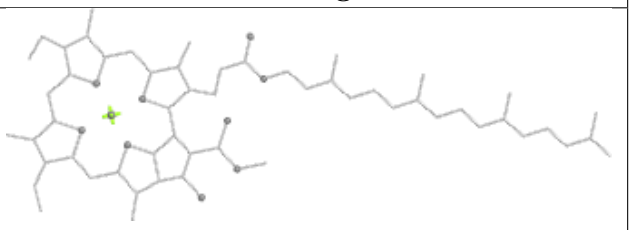


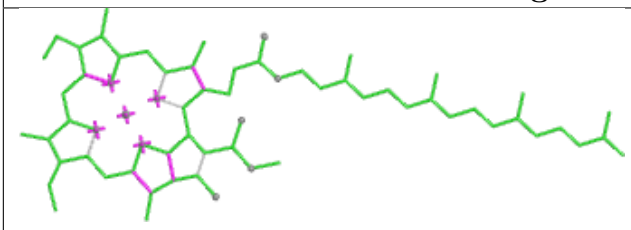
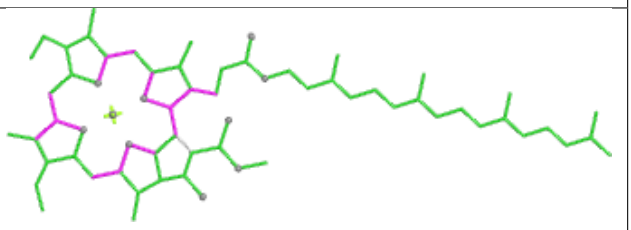
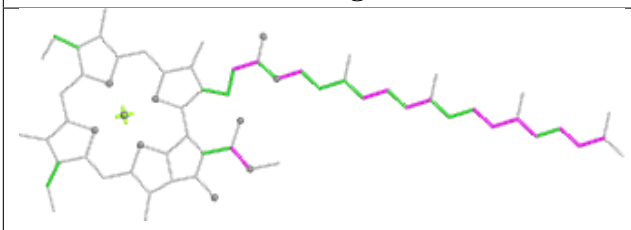
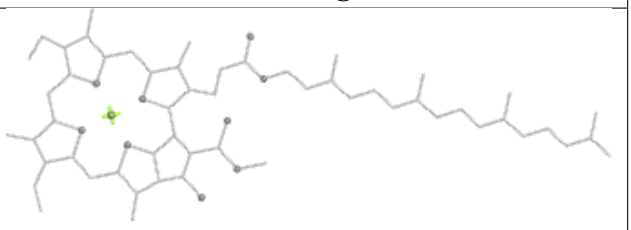
Torsions

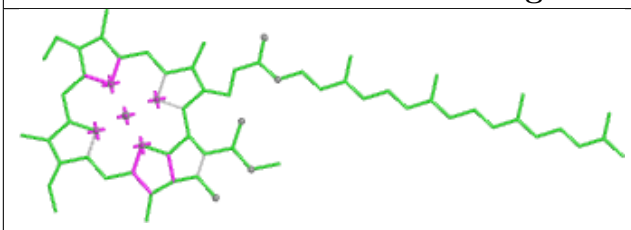
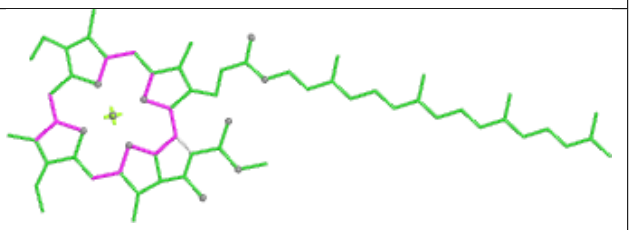
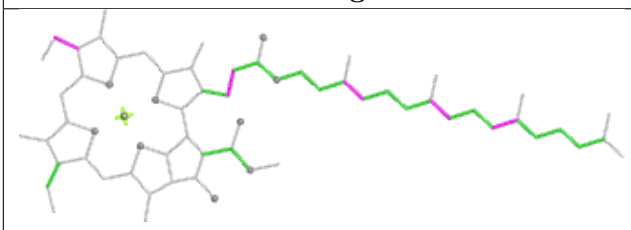
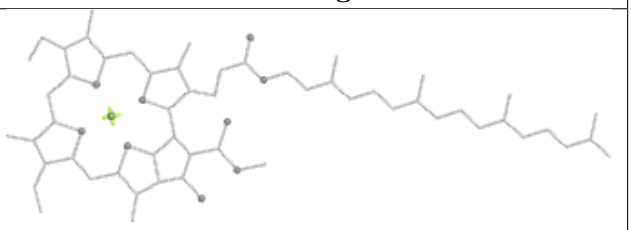


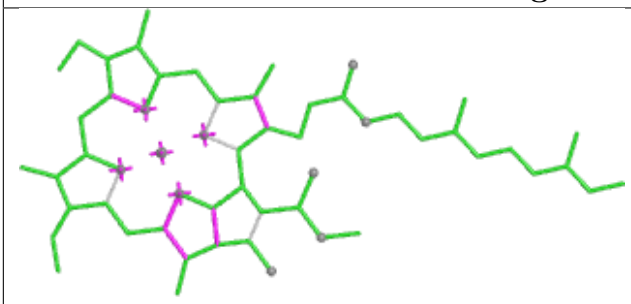
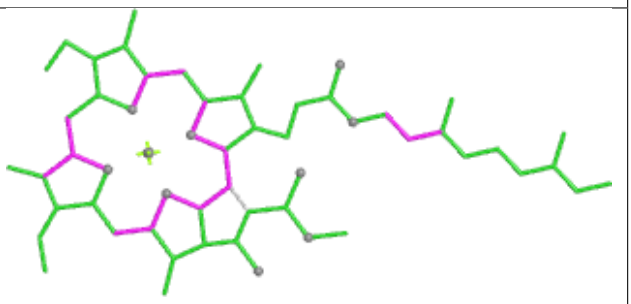
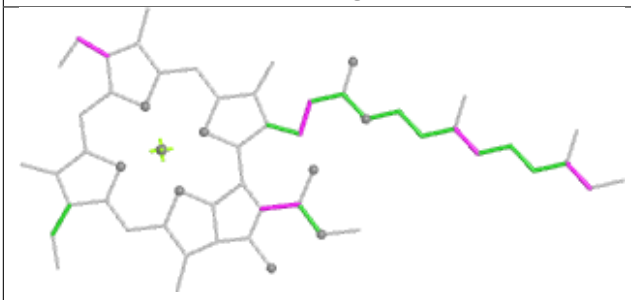
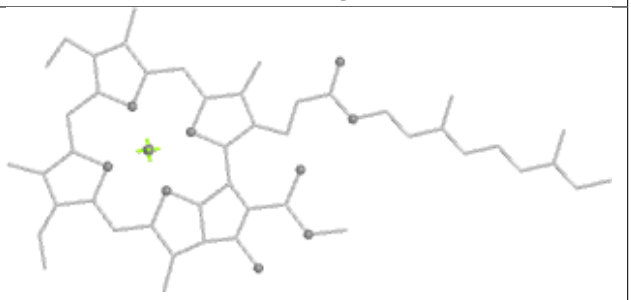
Rings

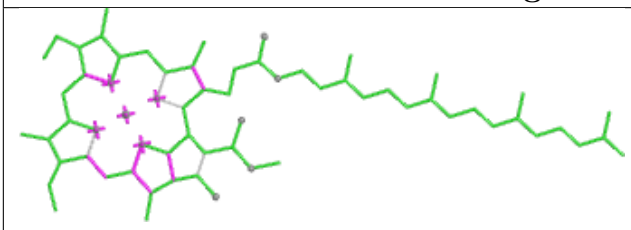
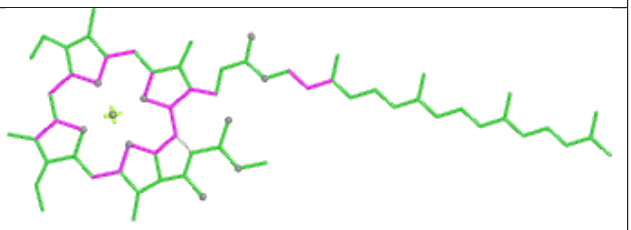
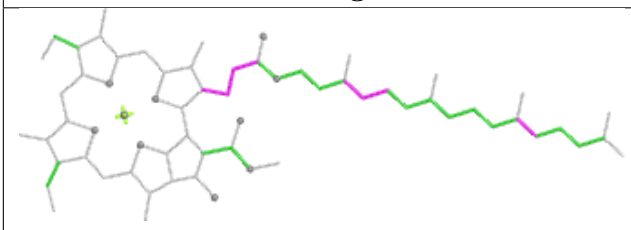
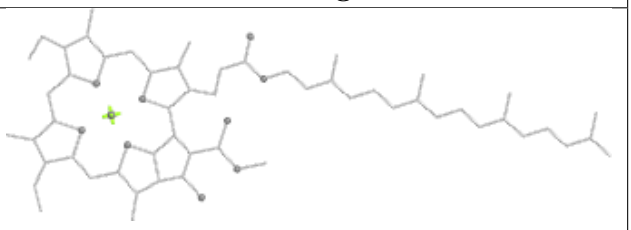
Ligand CLA A 603	
	
Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA E 609	
	
Bond lengths	Bond angles
	
Torsions	Rings

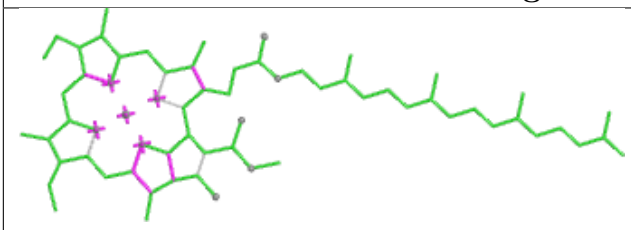
Ligand CLA I 607	
 Bond lengths	 Bond angles
 Torsions	 Rings

Ligand CLA E 604	
 Bond lengths	 Bond angles
 Torsions	 Rings

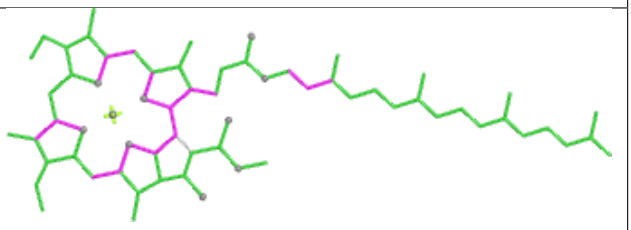
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Bond lengths	Bond angles
	
Torsions	Rings

Ligand CLA A 606	
	
Bond lengths	Bond angles
	
Torsions	Rings

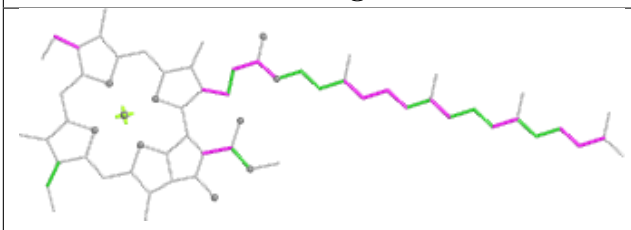
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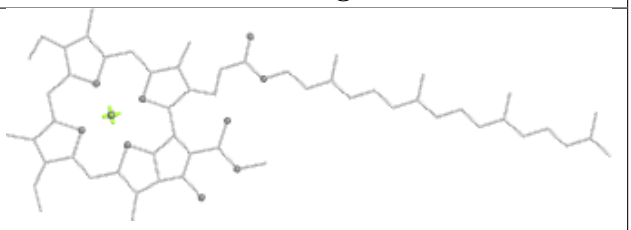
Bond lengths



Bond angles

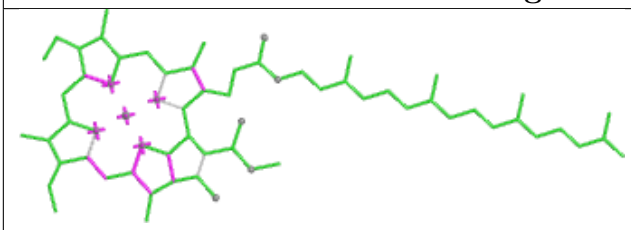


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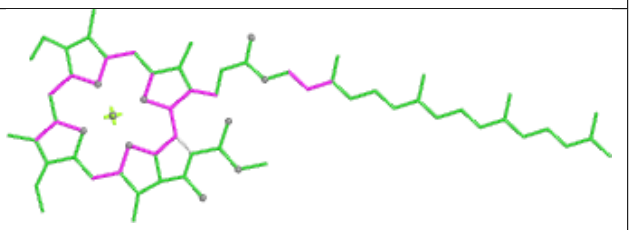


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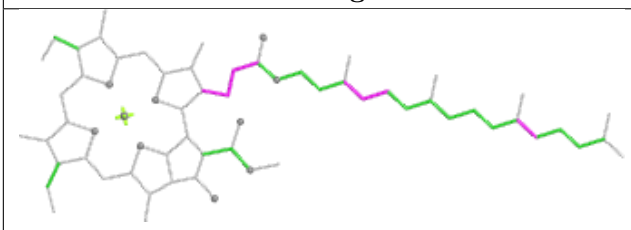
Ligand CLA B 606



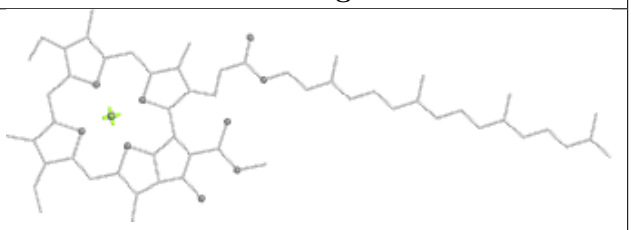
Bond lengths



Bond angles

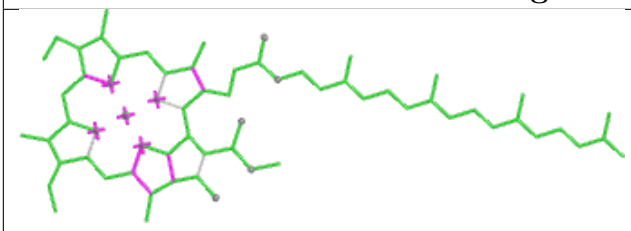


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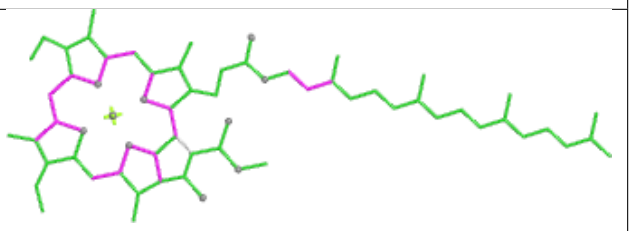


Rings

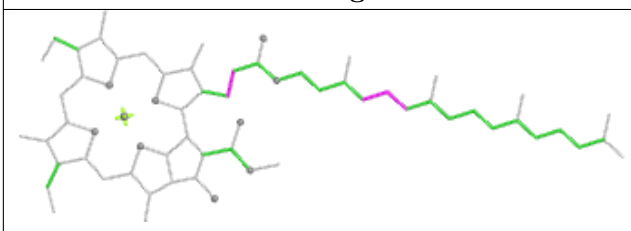
Ligand CLA I 609



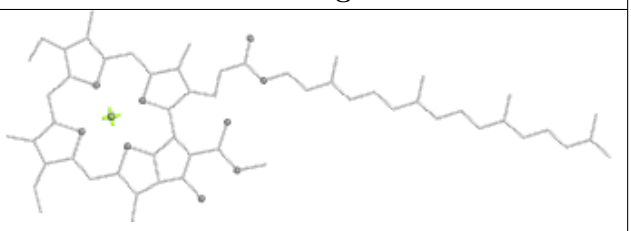
Bond lengths



Bond angles

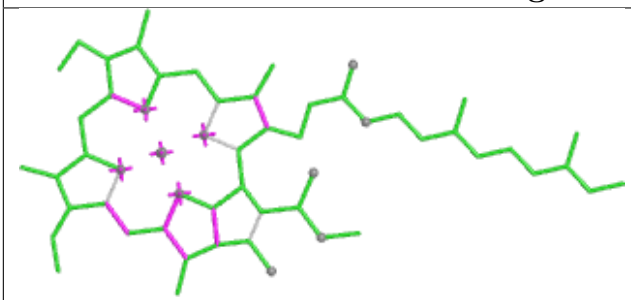


Torsions

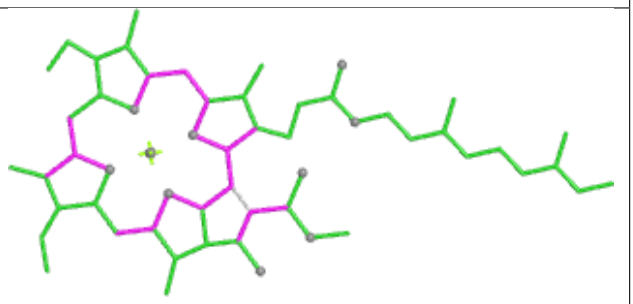


Rings

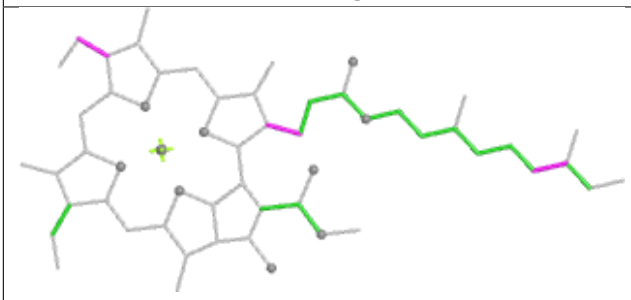
Ligand CLA J 608



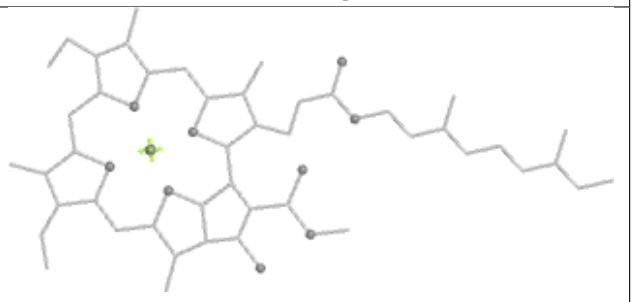
Bond lengths



Bond angles

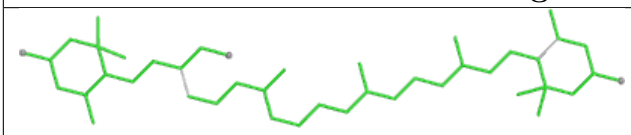


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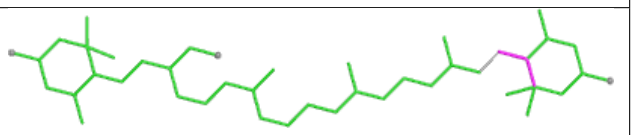


Rings

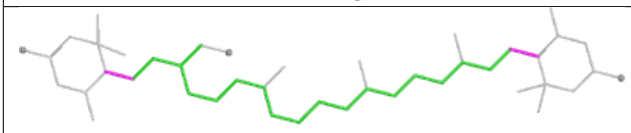
Ligand 32N E 613



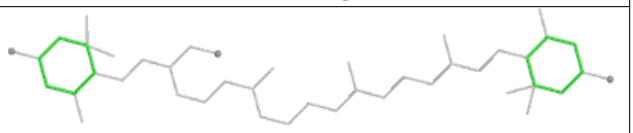
Bond lengths



Bond angles

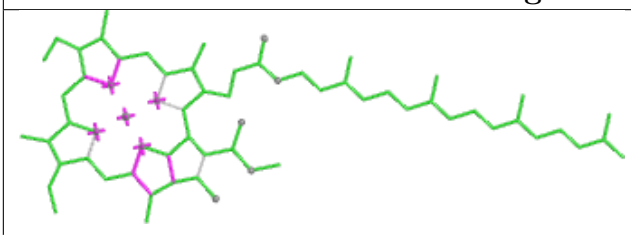


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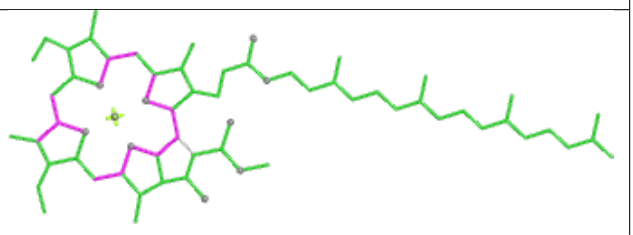


Rings

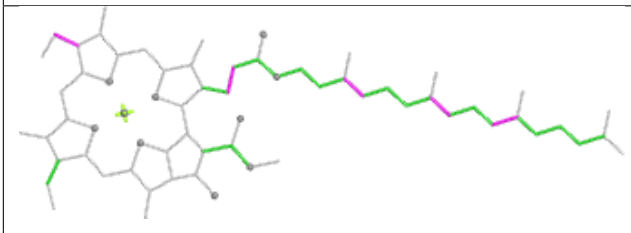
Ligand CLA C 604



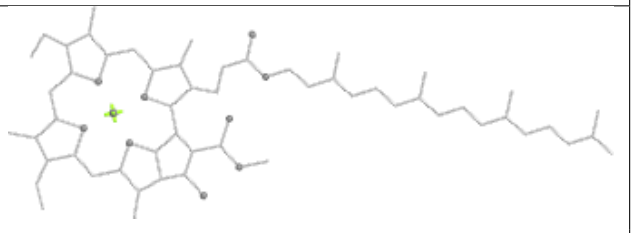
Bond lengths



Bond angles

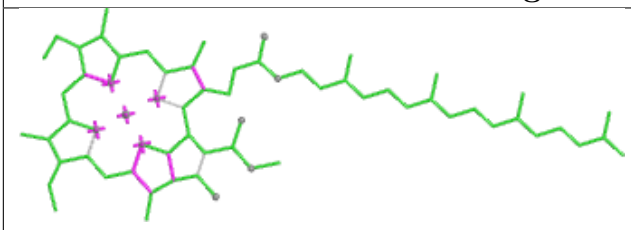


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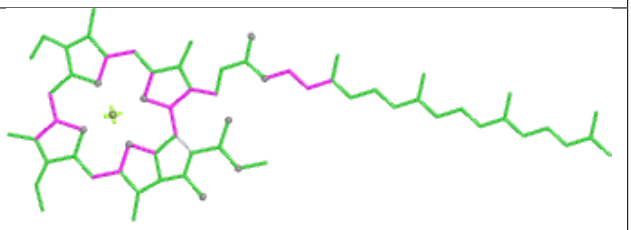


Rings

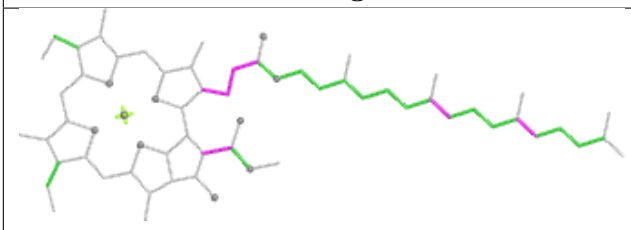
Ligand CLA D 602



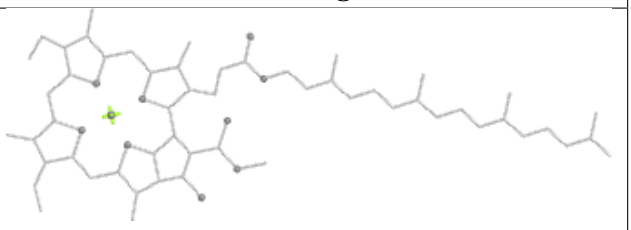
Bond lengths



Bond angles

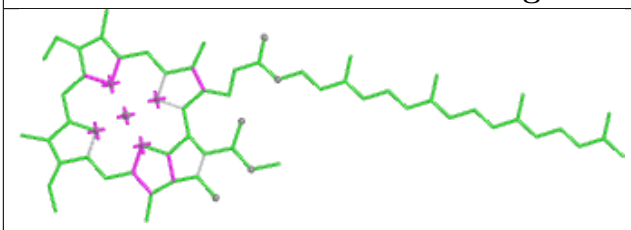


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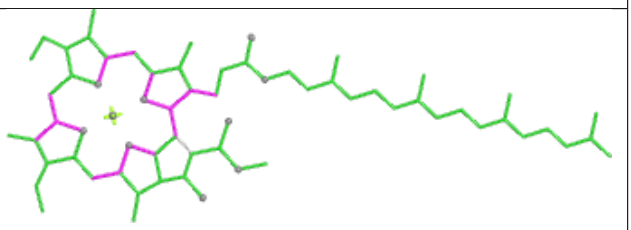


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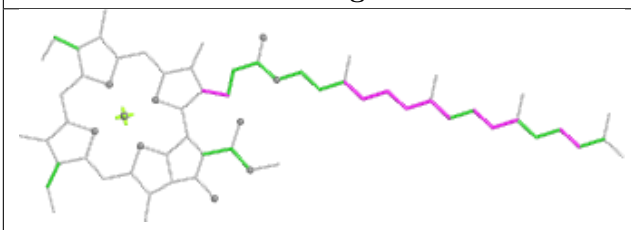
Ligand CLA G 601



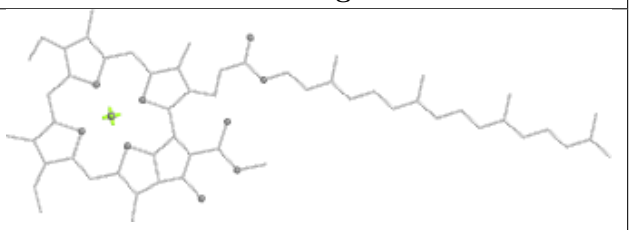
Bond lengths



Bond angles

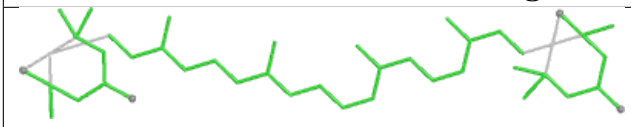


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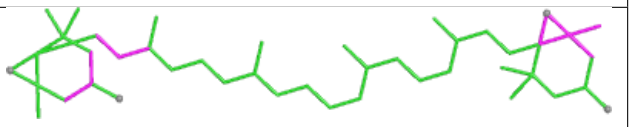


Rings

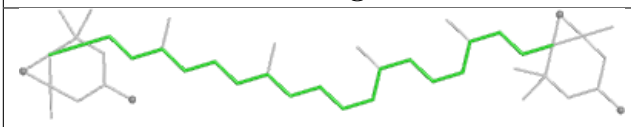
Ligand XAT B 612



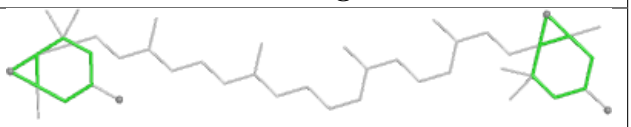
Bond lengths



Bond angles

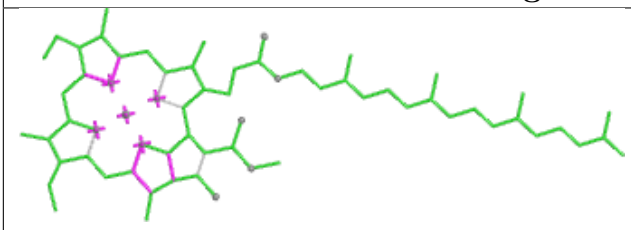


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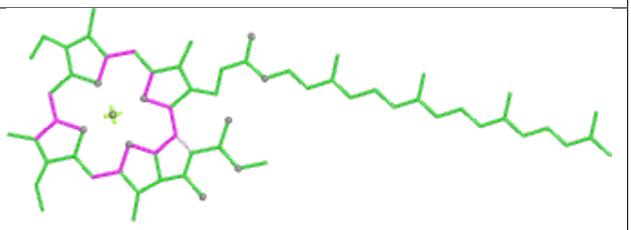


Rings

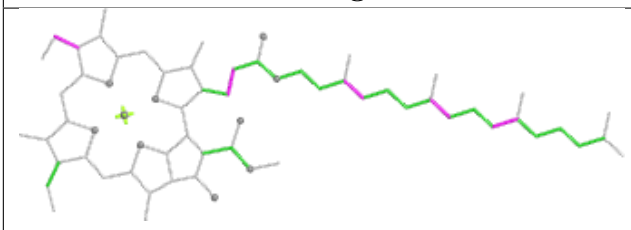
Ligand CLA H 604



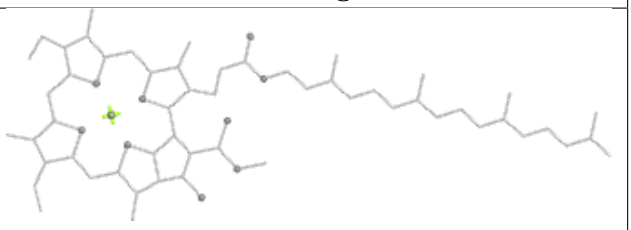
Bond lengths



Bond angles

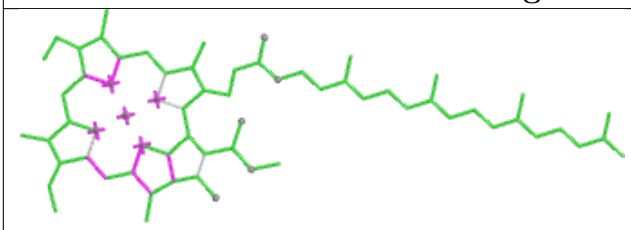


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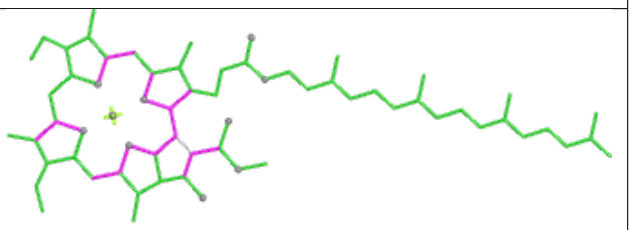


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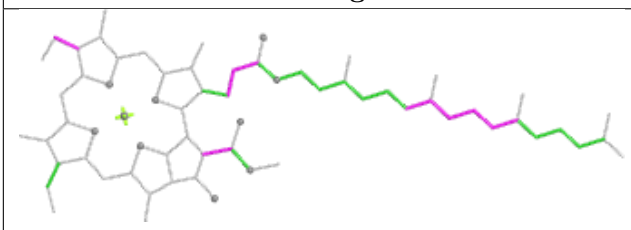
Ligand CLA H 610



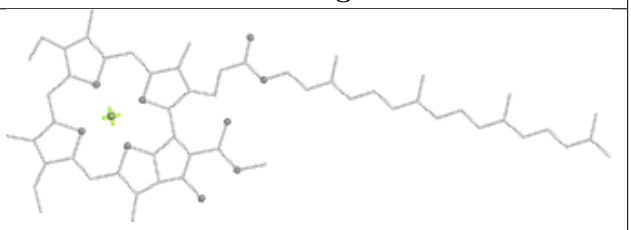
Bond lengths



Bond angles

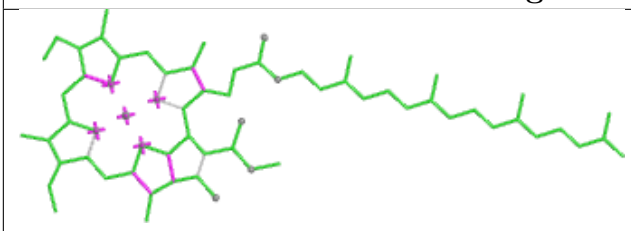


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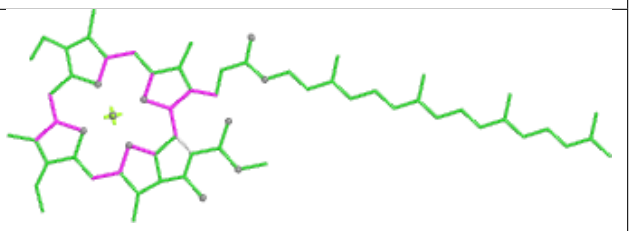


Rings

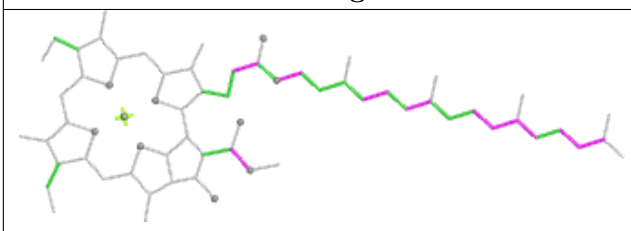
Ligand CLA F 607



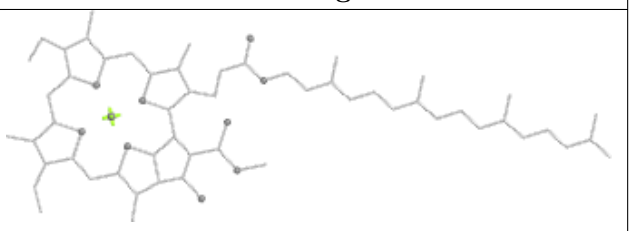
Bond lengths



Bond angles

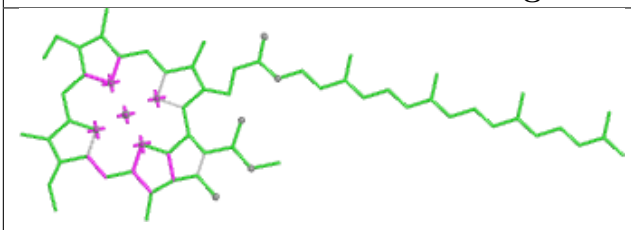


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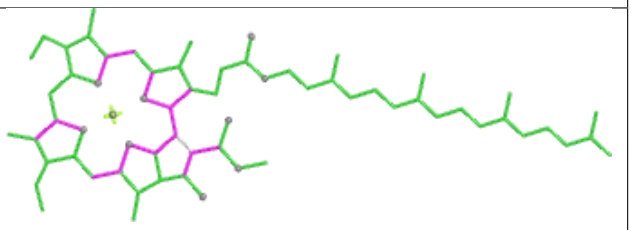


Rings

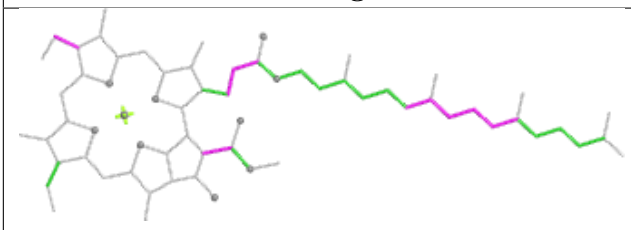
Ligand CLA A 610



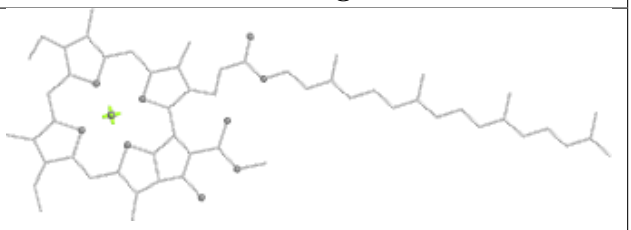
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

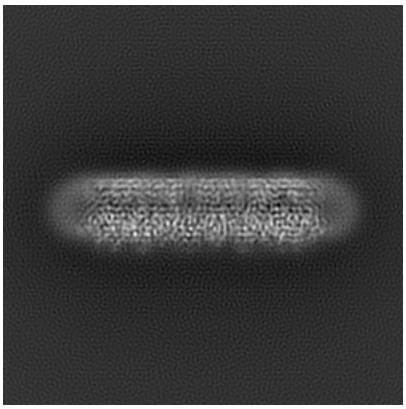
6 Map visualisation

This section contains visualisations of the EMDB entry EMD-35080. These allow visual inspection of the internal detail of the map and identification of artifacts.

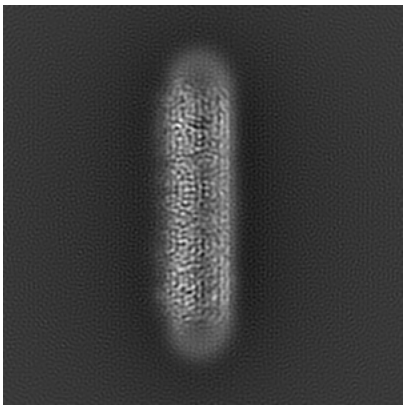
Images derived from a raw map, generated by summing the deposited half-maps, are presented below the corresponding image components of the primary map to allow further visual inspection and comparison with those of the primary map.

6.1 Orthogonal projections

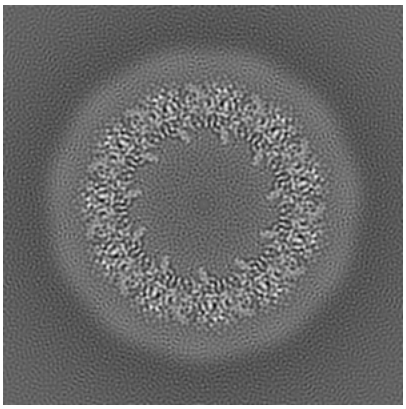
6.1.1 Primary map



X

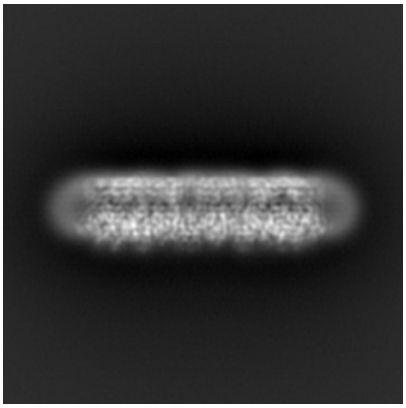


Y

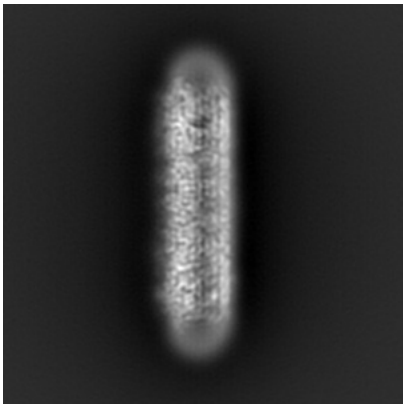


Z

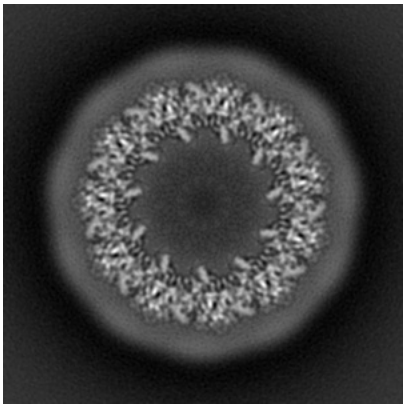
6.1.2 Raw map



X



Y

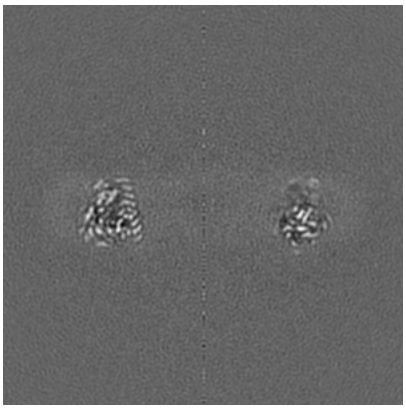


Z

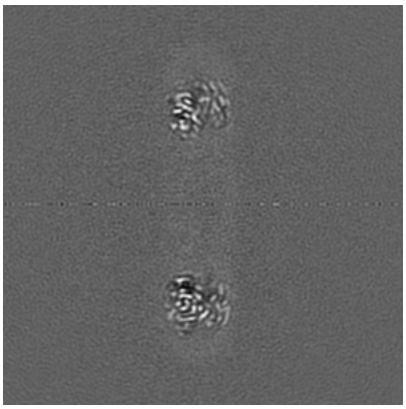
The images above show the map projected in three orthogonal directions.

6.2 Central slices (i)

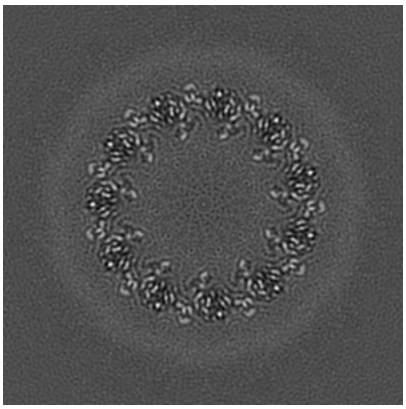
6.2.1 Primary map



X Index: 125

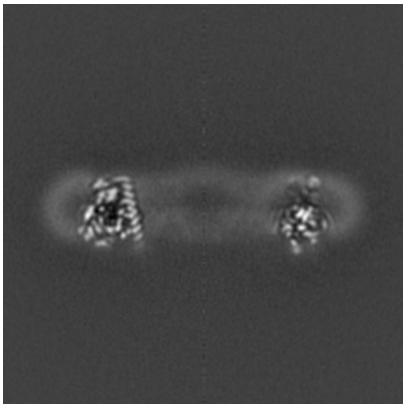


Y Index: 125

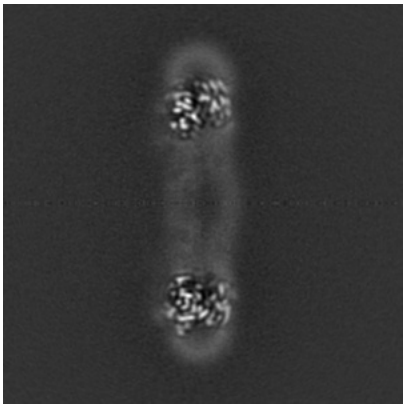


Z Index: 125

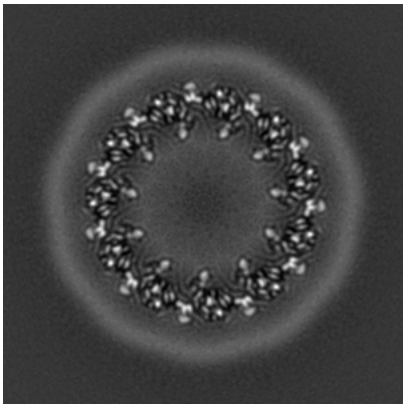
6.2.2 Raw map



X Index: 125



Y Index: 125

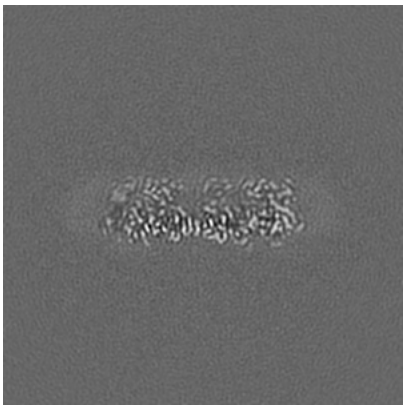


Z Index: 125

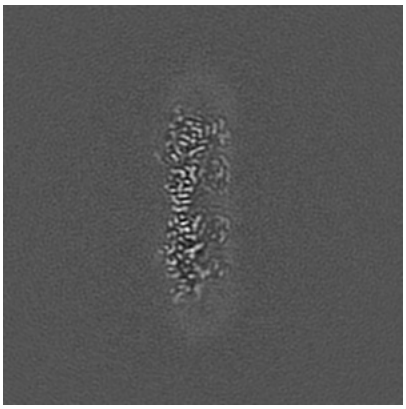
The images above show central slices of the map in three orthogonal directions.

6.3 Largest variance slices (i)

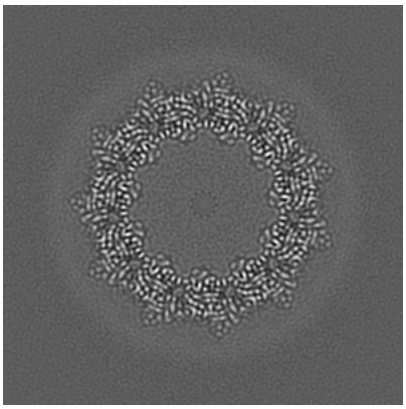
6.3.1 Primary map



X Index: 175

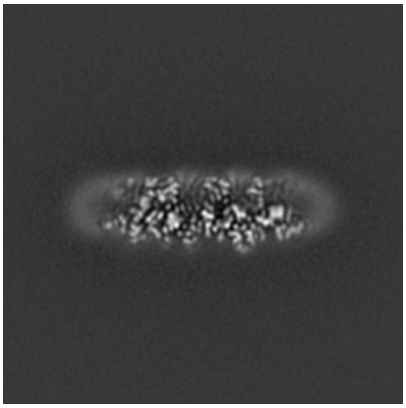


Y Index: 175

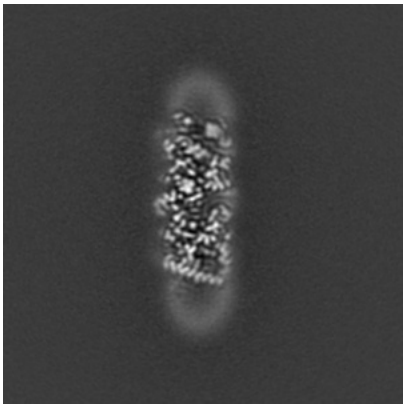


Z Index: 113

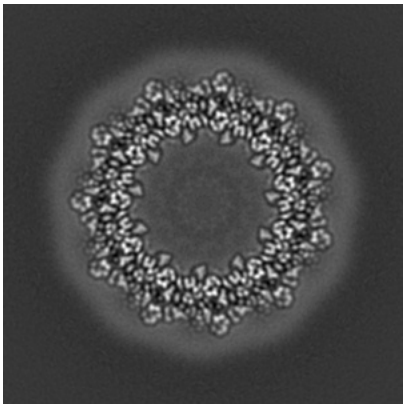
6.3.2 Raw map



X Index: 179



Y Index: 182

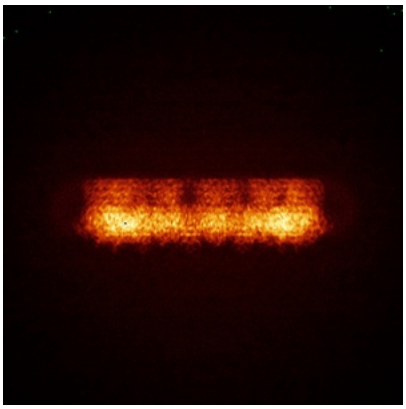


Z Index: 111

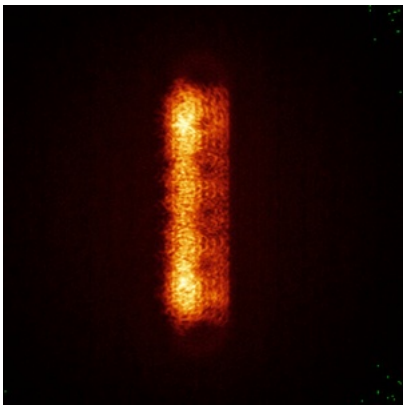
The images above show the largest variance slices of the map in three orthogonal directions.

6.4 Orthogonal standard-deviation projections (False-color) (i)

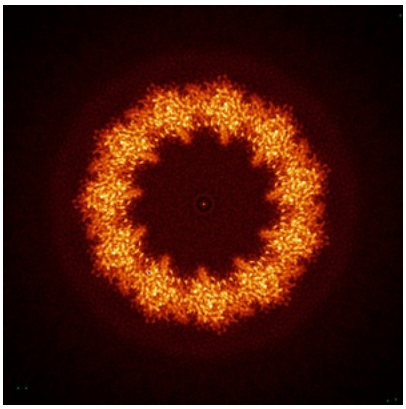
6.4.1 Primary map



X

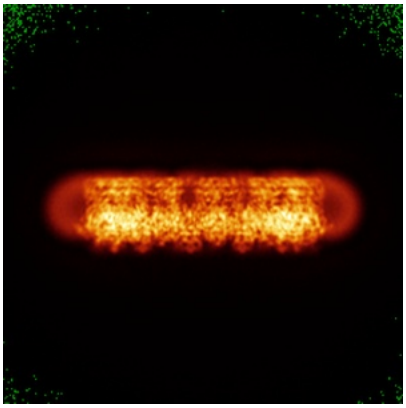


Y

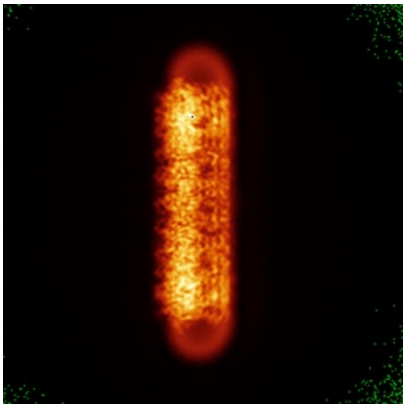


Z

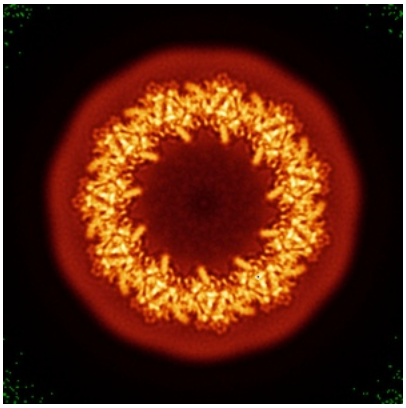
6.4.2 Raw map



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



X Y Z

The images above show the 3D surface view of the map at the recommended contour level 0.06. 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



X Y Z

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

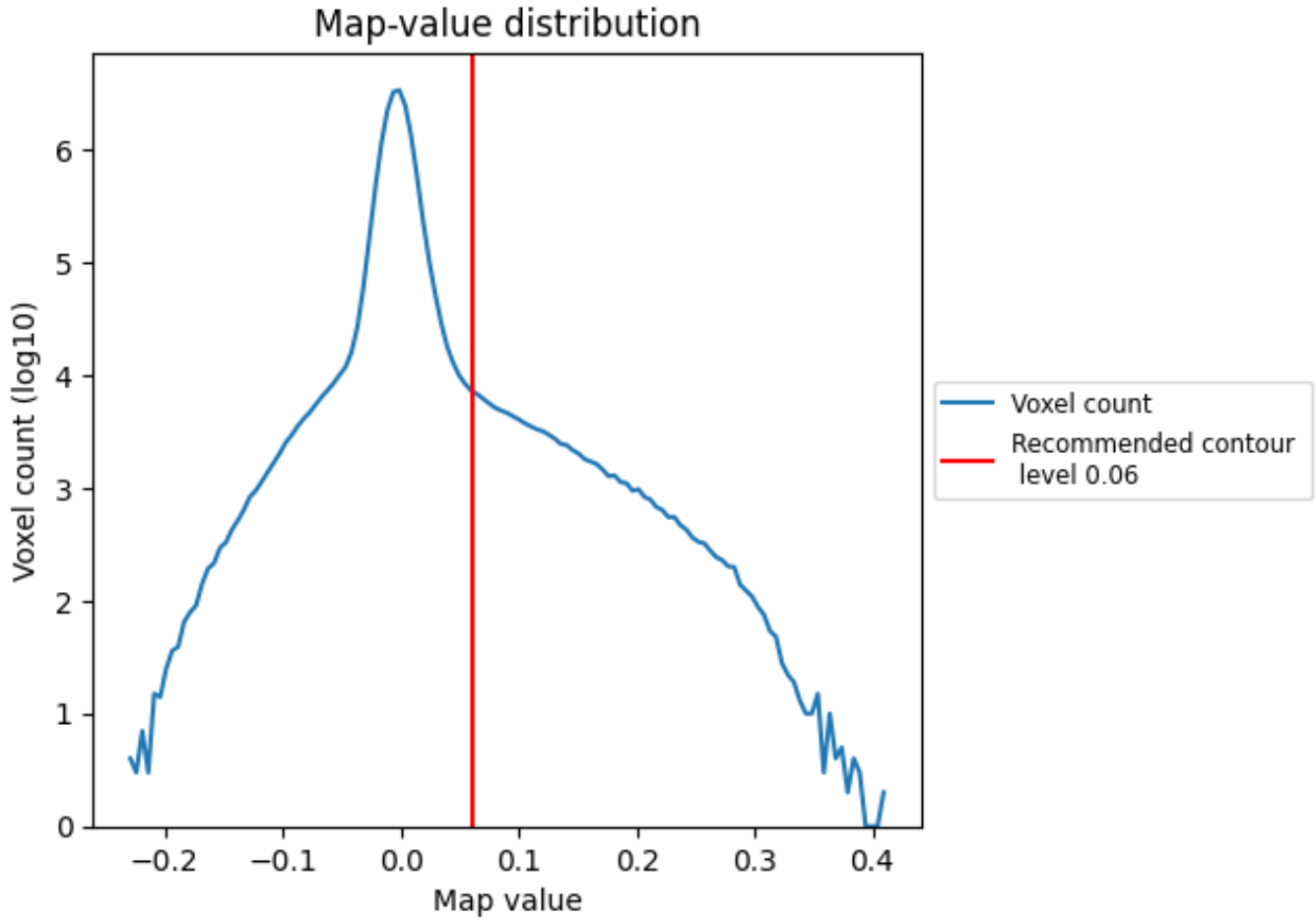
6.6 Mask visualisation (i)

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

7 Map analysis

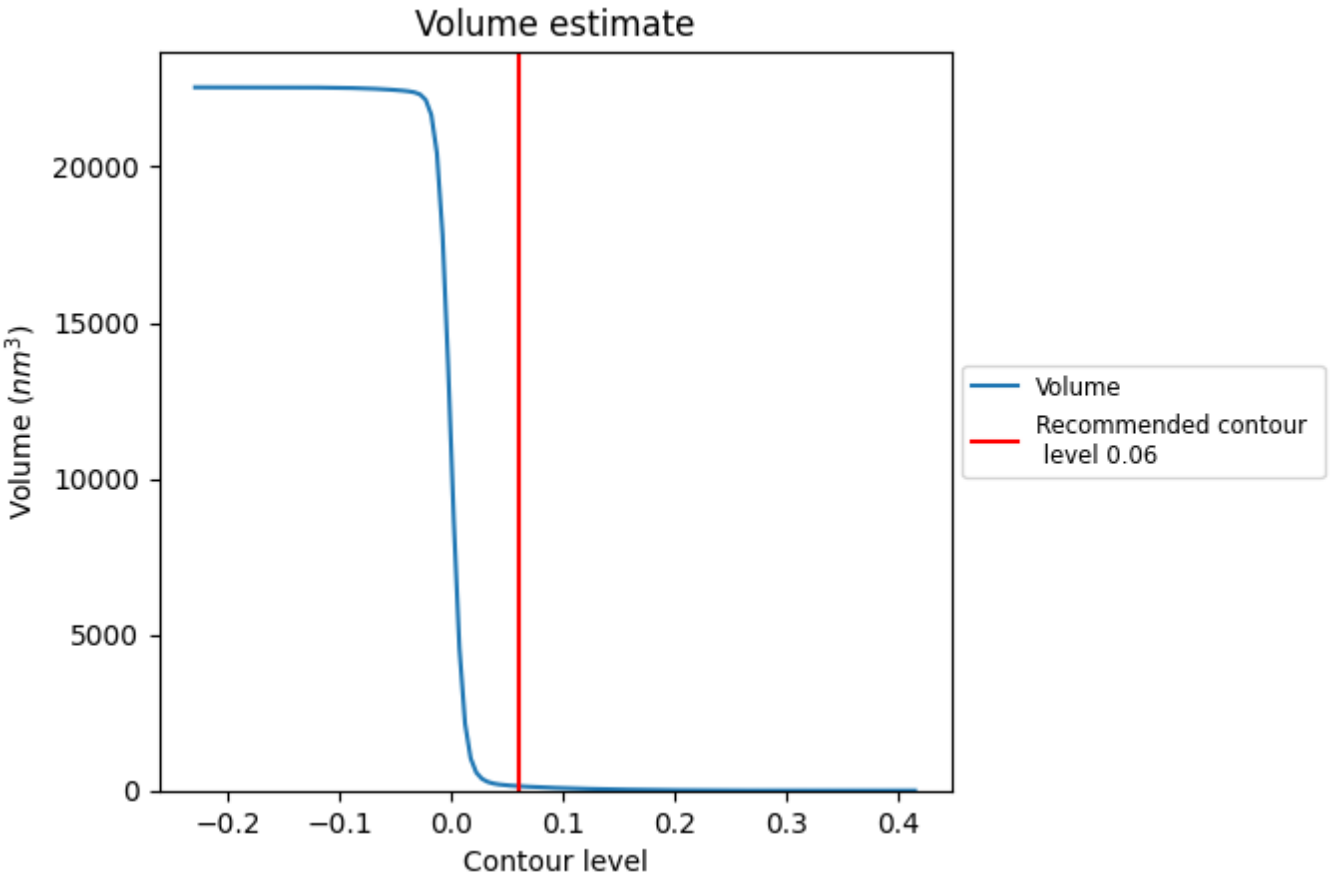
This section contains the results of statistical analysis of the map.

7.1 Map-value distribution



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.

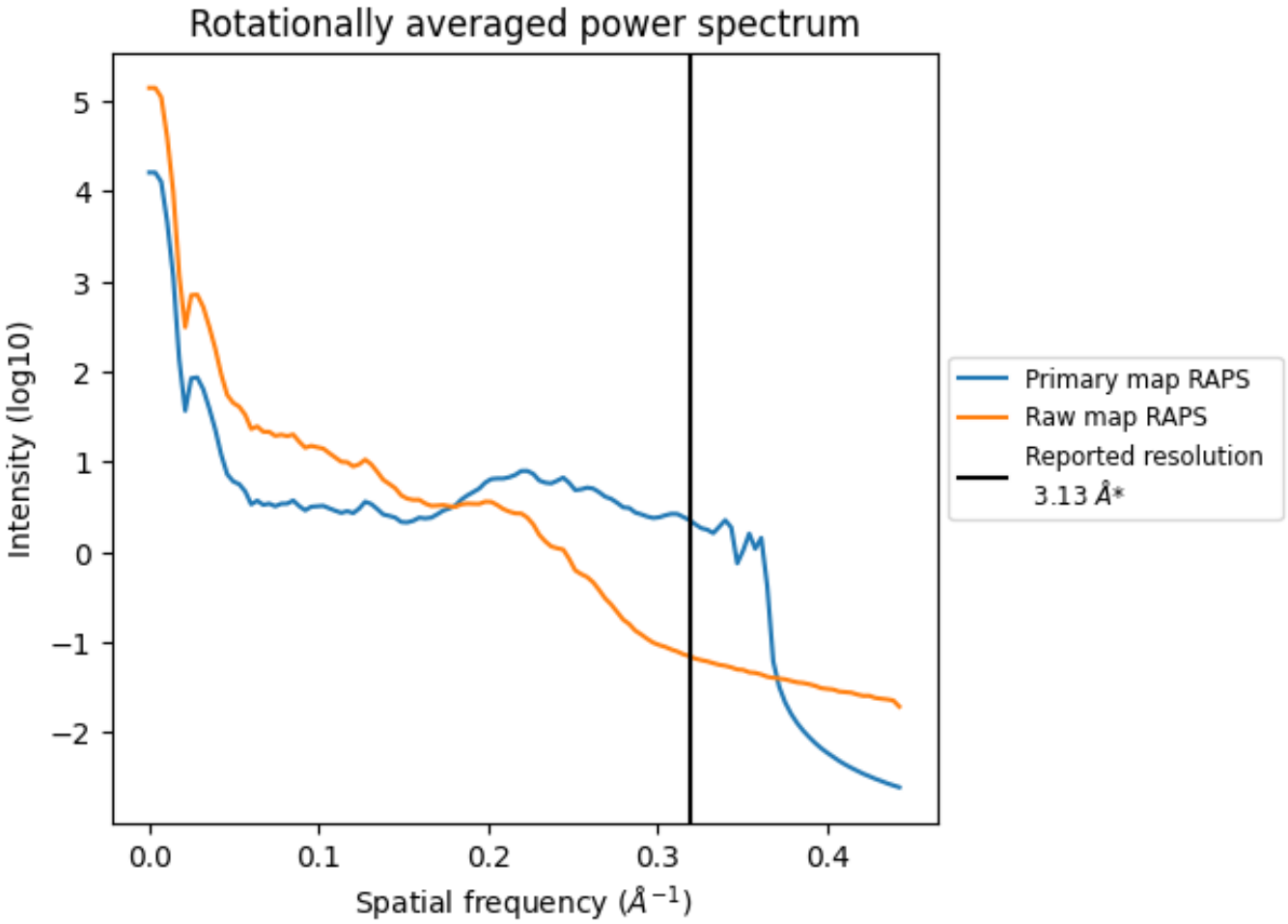
7.2 Volume estimate (i)



The volume at the recommended contour level is 142 nm^3; this corresponds to an approximate mass of 128 kDa.

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

7.3 Rotationally averaged power spectrum (i)

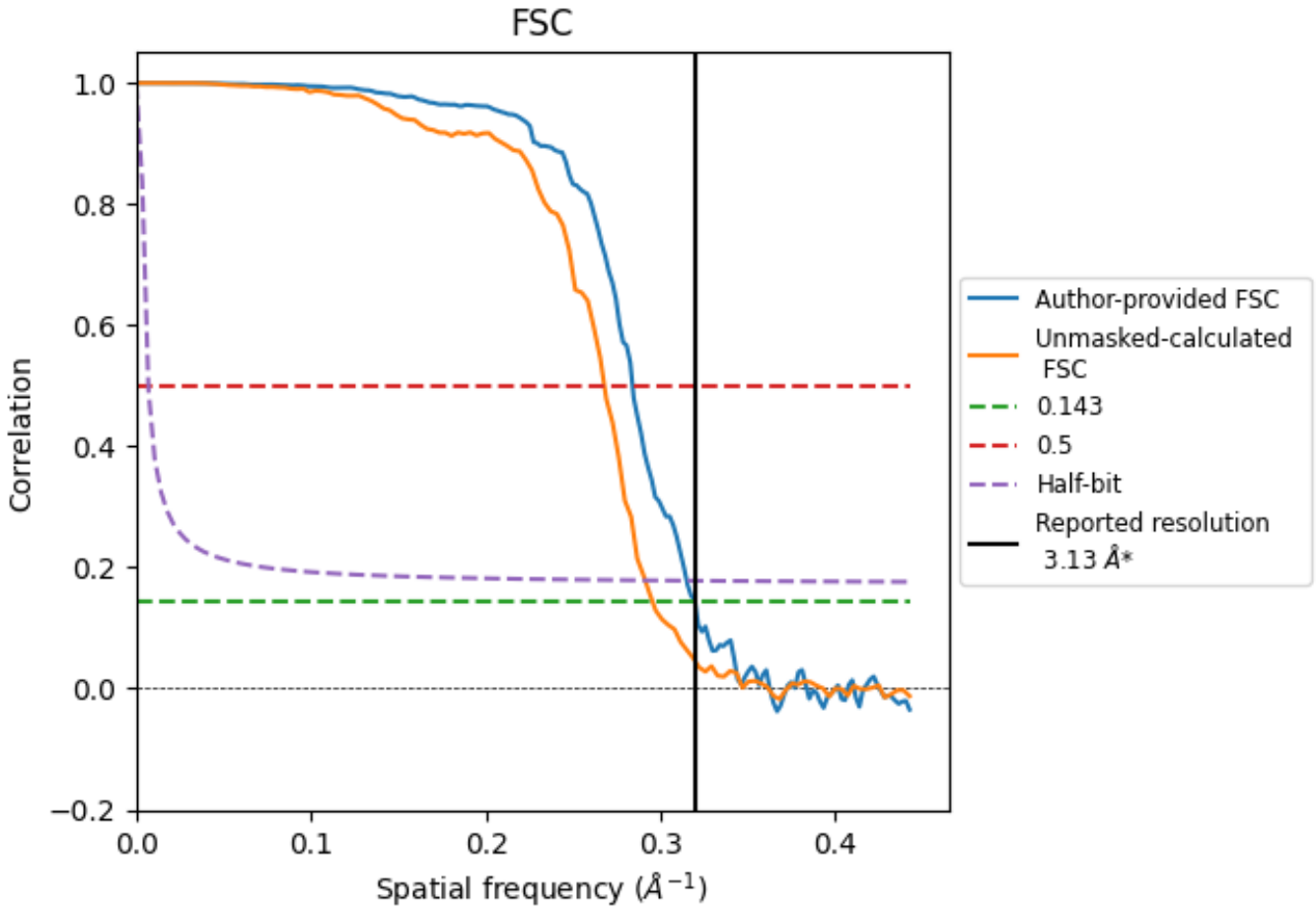


*Reported resolution corresponds to spatial frequency of 0.319 $Å^{-1}$

8 Fourier-Shell correlation

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



*Reported resolution corresponds to spatial frequency of 0.319 Å^{-1}

8.2 Resolution estimates (i)

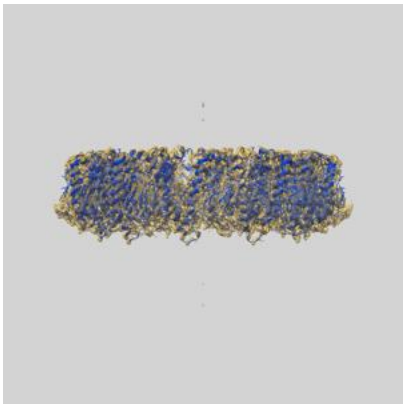
Resolution estimate (Å)	Estimation criterion (FSC cut-off)		
	0.143	0.5	Half-bit
Reported by author	3.13	-	-
Author-provided FSC curve	3.13	3.52	3.17
Unmasked-calculated*	3.38	3.73	3.43

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

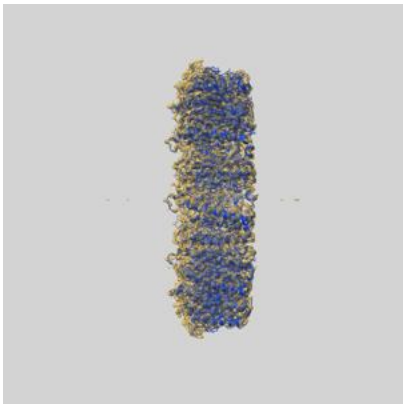
9 Map-model fit

This section contains information regarding the fit between EMDB map EMD-35080 and PDB model 8HW1. Per-residue inclusion information can be found in section 3 on page 14.

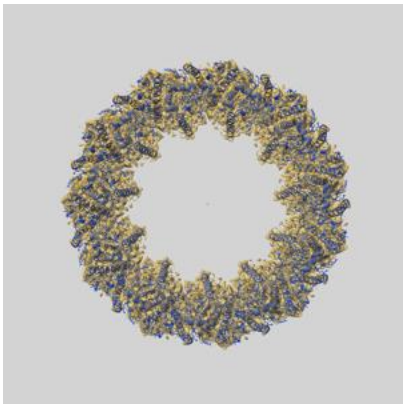
9.1 Map-model overlay



X



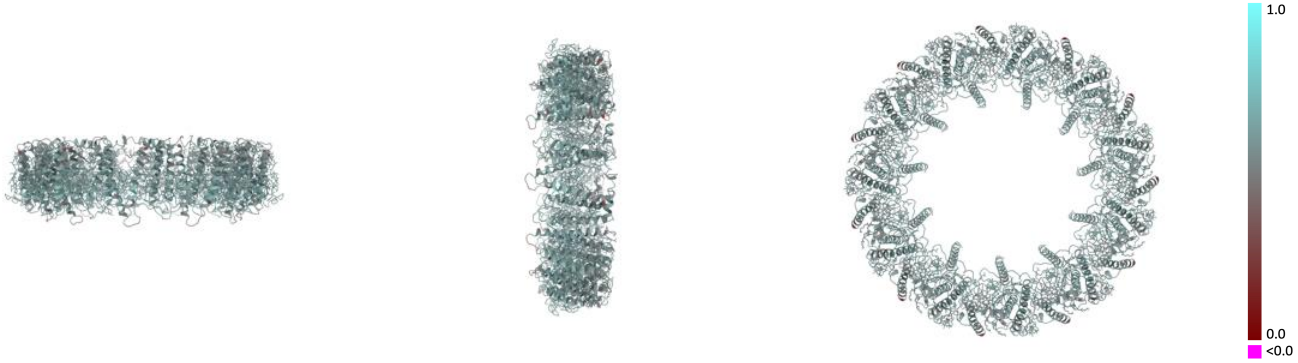
Y



Z

The images above show the 3D surface view of the map at the recommended contour level 0.06 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.

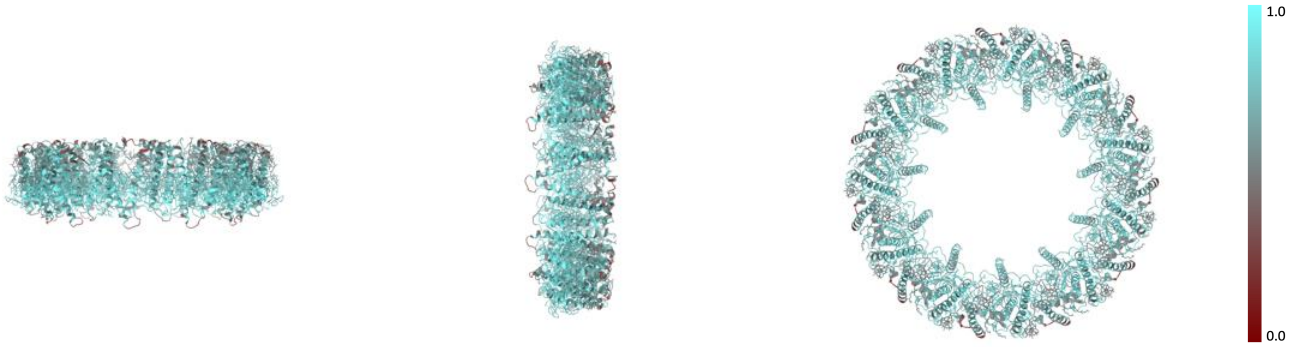
9.2 Q-score mapped to coordinate model (i)



X Y Z

The images above show the model with each residue coloured according 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.

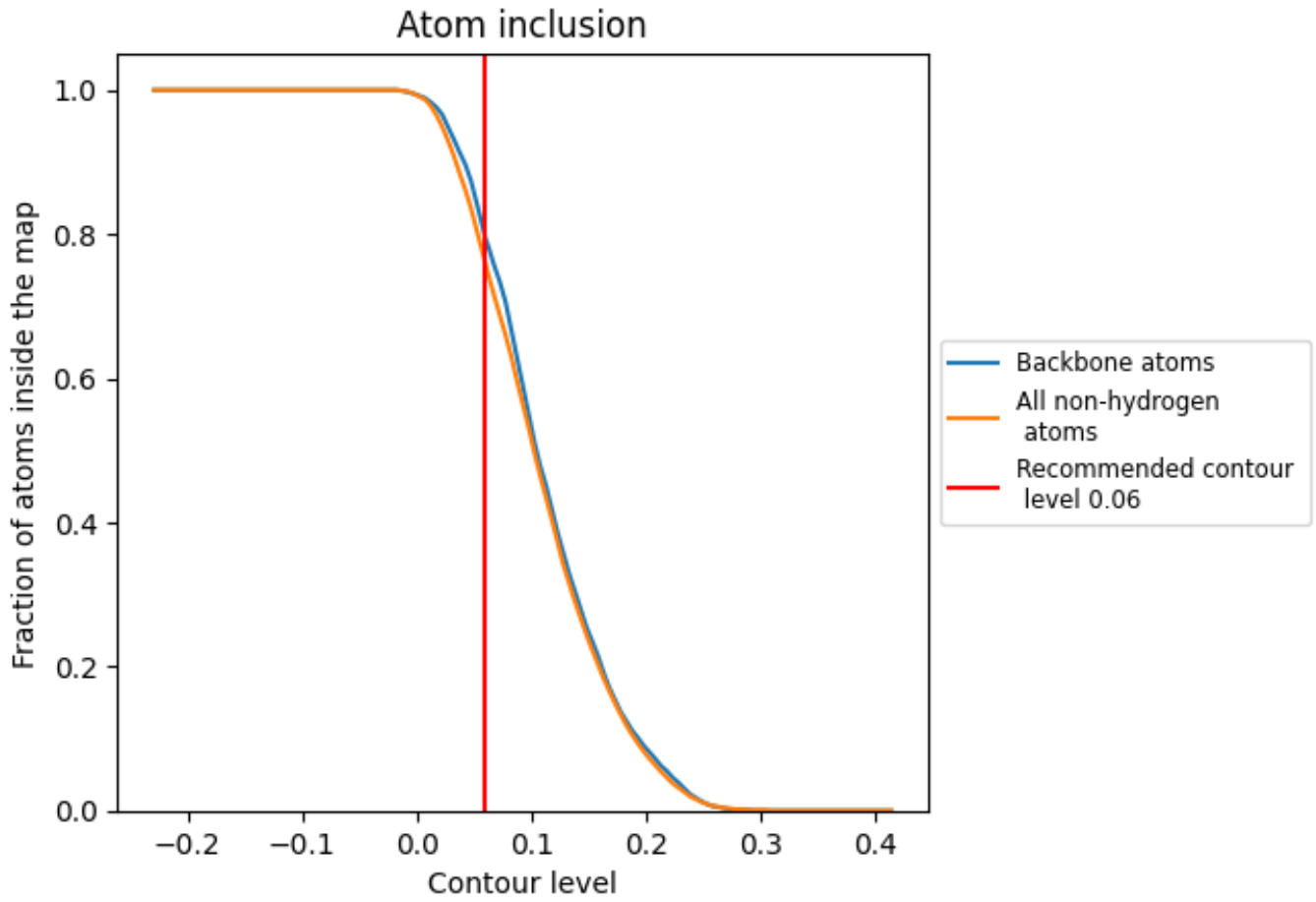
9.3 Atom inclusion mapped to coordinate model (i)



X Y Z

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 (0.06).

9.4 Atom inclusion (i)



At the recommended contour level, 79% of all backbone atoms, 76% 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 (0.06) and Q-score for the entire model and for each chain.

Chain	Atom inclusion	Q-score
All	0.7590	0.5800
A	0.7580	0.5810
B	0.7550	0.5810
C	0.7560	0.5790
D	0.7600	0.5810
E	0.7560	0.5800
F	0.7570	0.5780
G	0.7620	0.5780
H	0.7610	0.5800
I	0.7590	0.5790
J	0.7610	0.5800
K	0.7610	0.5810

