



wwPDB X-ray Structure Validation Summary Report ⓘ

Mar 8, 2026 – 03:34 PM UTC

PDB ID : 1GU6 / pdb_00001gu6
Title : Structure of the Periplasmic Cytochrome c Nitrite Reductase from Escherichia coli
Authors : Bamford, V.A.; Angove, H.C.; Seward, H.E.; Thomson, A.J.; Cole, J.A.; Butt, J.N.; Hemmings, A.M.; Richardson, D.J.
Deposited on : 2002-01-24
Resolution : 2.50 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

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

A user guide is available at

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

with specific help available everywhere you see the ⓘ symbol.

The types of validation reports are described at

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

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

MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	NOT EXECUTED
EDS	:	NOT EXECUTED
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

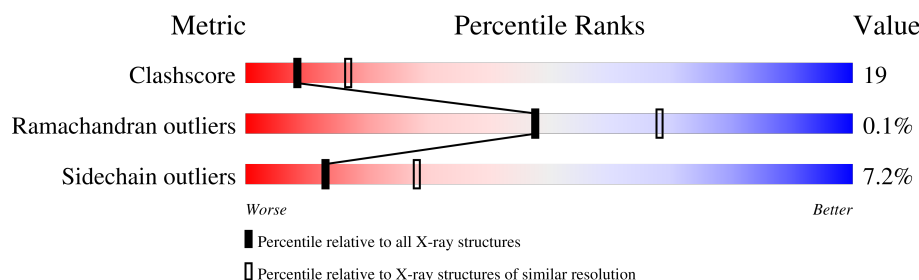
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for ≥ 3 , 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	452	 69% 25% . .
1	C	452	 66% 27% 5% .
1	E	452	 69% 25% . .
1	G	452	 67% 27% . .

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	GOL	A	1481	-	X	-	-
3	GOL	C	1481	-	X	-	-
3	GOL	E	1481	-	X	-	-
3	GOL	G	1481	-	X	-	-

2 Entry composition

There are 5 unique types of molecules in this entry. The entry contains 15486 atoms, of which 0 are hydrogens and 0 are deuteriums.

In the tables below, the ZeroOcc column contains the number of atoms modelled with zero occupancy, the AltConf column contains the number of residues with at least one atom in alternate conformation and the Trace column contains the number of residues modelled with at most 2 atoms.

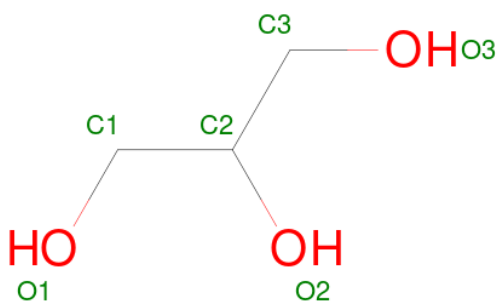
- Molecule 1 is a protein called CYTOCHROME C552.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	441	Total	C	N	O	S	0	0	0
			3479	2179	620	658	22			
1	C	441	Total	C	N	O	S	0	0	0
			3479	2179	620	658	22			
1	E	441	Total	C	N	O	S	0	0	0
			3479	2179	620	658	22			
1	G	441	Total	C	N	O	S	0	0	0
			3479	2179	620	658	22			

- Molecule 2 is CALCIUM ION (CCD ID: CA) (formula: Ca).

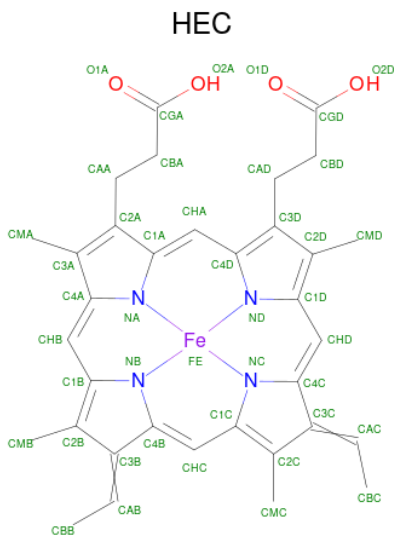
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Ca	0	0
			2	2		
2	C	2	Total	Ca	0	0
			2	2		
2	E	2	Total	Ca	0	0
			2	2		
2	G	2	Total	Ca	0	0
			2	2		

- Molecule 3 is GLYCEROL (CCD ID: GOL) (formula: C₃H₈O₃).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
3	A	1	Total 6	C 3	O 3	0	0
3	C	1	Total 6	C 3	O 3	0	0
3	E	1	Total 6	C 3	O 3	0	0
3	G	1	Total 6	C 3	O 3	0	0

- Molecule 4 is HEME C (CCD ID: HEC) (formula: $\text{C}_{34}\text{H}_{34}\text{FeN}_4\text{O}_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
4	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	C	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	E	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	E	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	E	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	E	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	E	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	G	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	G	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	G	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	G	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
4	G	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 5 is water.

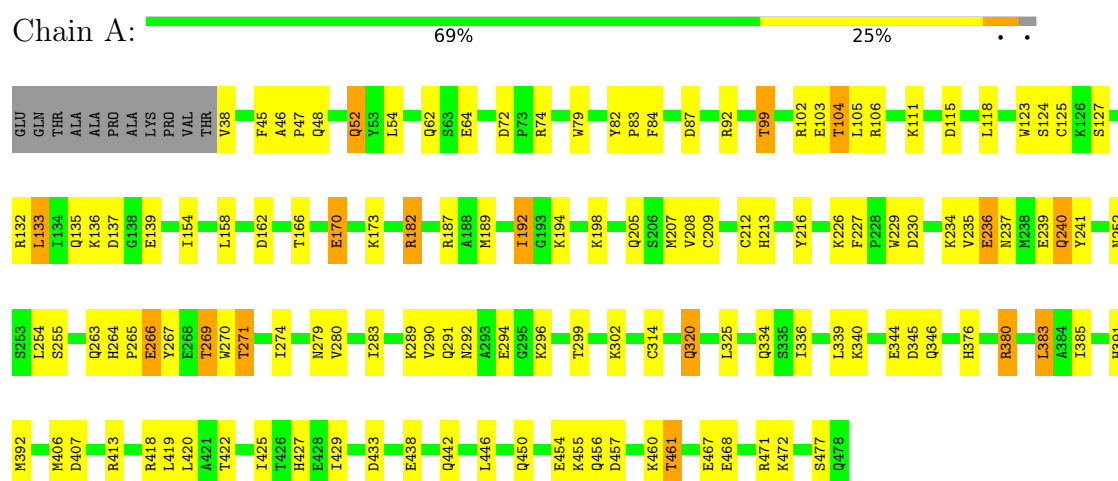
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	163	Total 163	O 163	0	0
5	C	121	Total 121	O 121	0	0
5	E	191	Total 191	O 191	0	0
5	G	203	Total 203	O 203	0	0

3 Residue-property plots

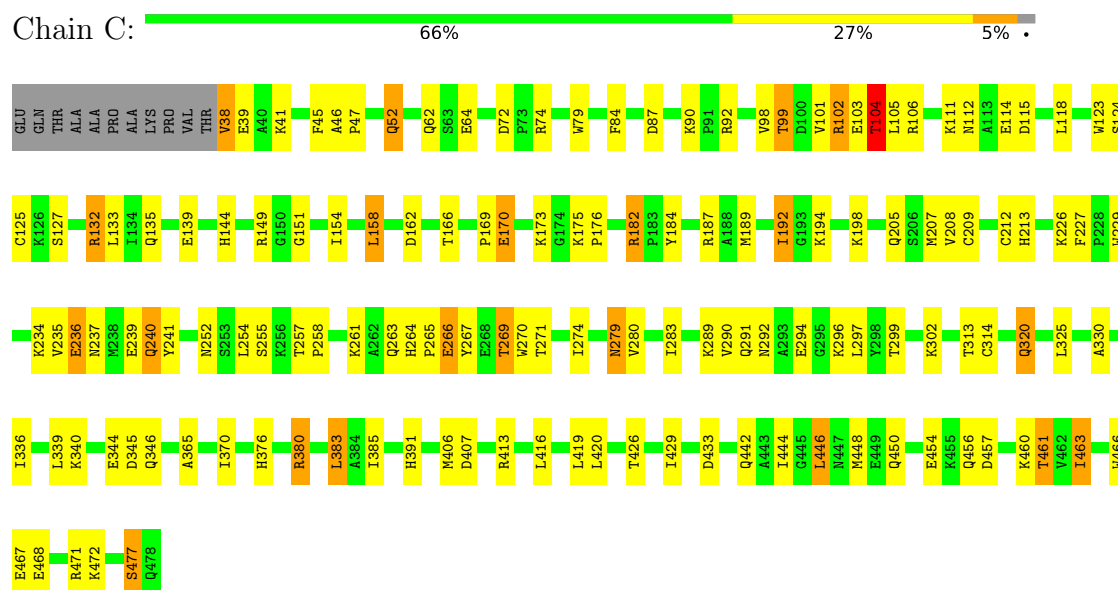
These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry. Residues are color-coded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

Note EDS was not executed.

• Molecule 1: CYTOCHROME C552

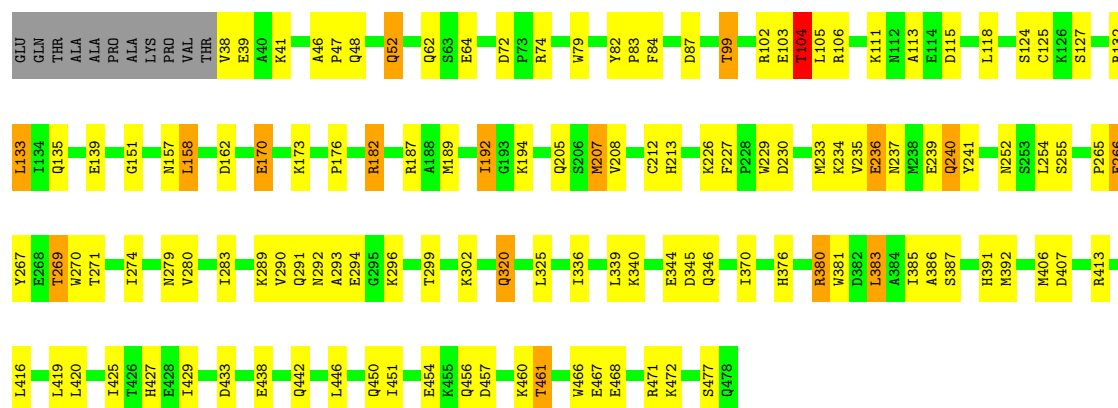


• Molecule 1: CYTOCHROME C552



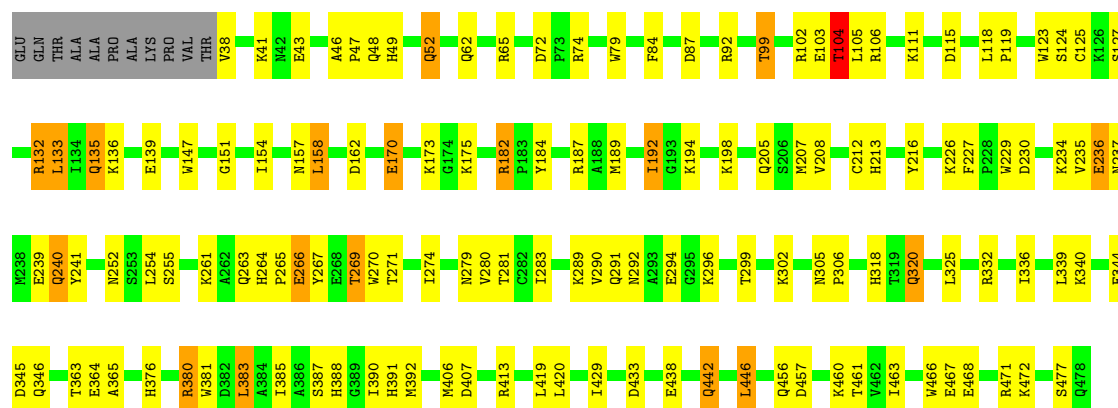
• Molecule 1: CYTOCHROME C552

Chain E:  69% 25% . .



• Molecule 1: CYTOCHROME C552

Chain G:  67% 27% . .



4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	81.47Å 90.84Å 293.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.50	Depositor
% Data completeness (in resolution range)	91.5 (20.00-2.50)	Depositor
R_{merge}	0.09	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	CNS 1.0	Depositor
R, R_{free}	0.203 , 0.243	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	15486	wwPDB-VP
Average B, all atoms (Å ²)	35.0	wwPDB-VP

5 Model quality

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEC, GOL, CA

The Z score for a bond length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 5$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Chain	Bond lengths		Bond angles	
		RMSZ	# Z >5	RMSZ	# Z >5
1	A	0.85	0/3563	1.01	3/4817 (0.1%)
1	C	0.89	1/3563 (0.0%)	1.02	11/4817 (0.2%)
1	E	0.90	2/3563 (0.1%)	1.02	5/4817 (0.1%)
1	G	0.88	0/3563	1.02	10/4817 (0.2%)
All	All	0.88	3/14252 (0.0%)	1.02	29/19268 (0.2%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1
1	G	0	1
All	All	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	E	207	MET	SD-CE	-6.44	1.63	1.79
1	E	233	MET	SD-CE	5.94	1.94	1.79
1	C	313	THR	CA-C	5.86	1.55	1.52

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	G	127	SER	N-CA-C	6.89	117.00	108.11
1	G	365	ALA	N-CA-C	-6.49	104.30	111.82
1	A	127	SER	N-CA-C	6.15	116.04	108.11
1	A	118	LEU	CA-C-N	6.13	126.59	120.52

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Mol	Chain	Res	Type	Atoms	Z	Observed(^o)	Ideal(^o)
1	A	118	LEU	C-N-CA	6.13	126.59	120.52

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	216	TYR	Sidechain
1	G	216	TYR	Sidechain

5.2 Too-close contacts

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3479	0	3365	134	0
1	C	3479	0	3365	139	0
1	E	3479	0	3365	132	0
1	G	3479	0	3365	143	0
2	A	2	0	0	0	0
2	C	2	0	0	0	0
2	E	2	0	0	0	0
2	G	2	0	0	0	0
3	A	6	0	5	3	0
3	C	6	0	5	3	0
3	E	6	0	4	3	0
3	G	6	0	5	3	0
4	A	215	0	150	12	0
4	C	215	0	150	12	0
4	E	215	0	150	8	0
4	G	215	0	150	9	0
5	A	163	0	0	16	0
5	C	121	0	0	20	0
5	E	191	0	0	24	0
5	G	203	0	0	28	0
All	All	15486	0	14079	550	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 19.

The worst 5 of 550 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)
3:A:1481:GOL:O1	3:A:1481:GOL:C1	1.66	1.44
3:G:1481:GOL:C1	3:G:1481:GOL:O1	1.65	1.43
3:E:1481:GOL:O1	3:E:1481:GOL:C1	1.65	1.43
3:C:1481:GOL:O1	3:C:1481:GOL:C1	1.68	1.42
1:E:240:GLN:HB2	5:E:2102:HOH:O	1.54	1.07

There are no symmetry-related clashes.

5.3 Torsion angles [i](#)

5.3.1 Protein backbone [i](#)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	439/452 (97%)	421 (96%)	18 (4%)	0	100	100
1	C	439/452 (97%)	421 (96%)	18 (4%)	0	100	100
1	E	439/452 (97%)	421 (96%)	17 (4%)	1 (0%)	43	63
1	G	439/452 (97%)	419 (95%)	20 (5%)	0	100	100
All	All	1756/1808 (97%)	1682 (96%)	73 (4%)	1 (0%)	48	68

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	E	113	ALA

5.3.2 Protein sidechains [i](#)

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

The Analysed column shows the number of residues for which the sidechain conformation was

analysed, and the total number of residues.

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	362/370 (98%)	335 (92%)	27 (8%)	12	26
1	C	362/370 (98%)	336 (93%)	26 (7%)	13	28
1	E	362/370 (98%)	336 (93%)	26 (7%)	13	28
1	G	362/370 (98%)	337 (93%)	25 (7%)	14	30
All	All	1448/1480 (98%)	1344 (93%)	104 (7%)	13	28

5 of 104 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	E	104	THR
1	E	325	LEU
1	G	380	ARG
1	E	133	LEU
1	E	236	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 72 such sidechains are listed below:

Mol	Chain	Res	Type
1	G	205	GLN
1	G	456	GLN
1	G	252	ASN
1	G	326	GLN
1	C	252	ASN

5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

5.5 Carbohydrates ⓘ

There are no oligosaccharides in this entry.

5.6 Ligand geometry

Of 32 ligands modelled in this entry, 8 are monoatomic - leaving 24 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
3	GOL	E	1481	-	5,5,5	4.57	4 (80%)	5,5,5	6.19	3 (60%)
4	HEC	A	1486	1	46,50,50	1.66	7 (15%)	58,82,82	1.57	6 (10%)
3	GOL	C	1481	-	5,5,5	4.66	3 (60%)	5,5,5	6.21	3 (60%)
4	HEC	G	1486	1	46,50,50	1.81	9 (19%)	58,82,82	1.59	5 (8%)
4	HEC	E	1486	1	46,50,50	1.86	9 (19%)	58,82,82	1.61	5 (8%)
4	HEC	C	1485	2,1	46,50,50	1.70	9 (19%)	58,82,82	1.79	7 (12%)
4	HEC	A	1483	1	46,50,50	1.79	8 (17%)	58,82,82	1.66	7 (12%)
4	HEC	C	1483	1	46,50,50	1.90	9 (19%)	58,82,82	1.56	6 (10%)
4	HEC	E	1483	1	46,50,50	1.77	9 (19%)	58,82,82	1.63	6 (10%)
4	HEC	E	1482	5,1	46,50,50	1.78	7 (15%)	58,82,82	1.73	6 (10%)
3	GOL	A	1481	-	5,5,5	4.60	5 (100%)	5,5,5	6.23	3 (60%)
4	HEC	A	1485	2,1	46,50,50	1.69	7 (15%)	58,82,82	1.80	7 (12%)
4	HEC	C	1486	1	46,50,50	1.77	8 (17%)	58,82,82	1.61	6 (10%)
4	HEC	G	1482	5,1	46,50,50	1.70	7 (15%)	58,82,82	1.73	5 (8%)
4	HEC	G	1483	1	46,50,50	1.77	8 (17%)	58,82,82	1.61	6 (10%)
4	HEC	G	1484	2,1	46,50,50	1.69	12 (26%)	58,82,82	1.65	5 (8%)
4	HEC	C	1482	5,1	46,50,50	1.72	7 (15%)	58,82,82	1.68	7 (12%)
4	HEC	A	1484	2,1	46,50,50	1.68	7 (15%)	58,82,82	1.68	6 (10%)
3	GOL	G	1481	-	5,5,5	4.62	4 (80%)	5,5,5	6.19	3 (60%)
4	HEC	E	1484	2,1	46,50,50	1.70	7 (15%)	58,82,82	1.66	5 (8%)
4	HEC	G	1485	2,1	46,50,50	1.76	9 (19%)	58,82,82	1.77	6 (10%)
4	HEC	E	1485	2,1	46,50,50	1.81	8 (17%)	58,82,82	1.77	7 (12%)
4	HEC	C	1484	2,1	46,50,50	1.73	8 (17%)	58,82,82	1.69	5 (8%)
4	HEC	A	1482	5,1	46,50,50	1.64	8 (17%)	58,82,82	1.65	7 (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
3	GOL	E	1481	-	-	2/4/4/4	-
4	HEC	A	1486	1	-	8/14/54/54	-
3	GOL	C	1481	-	-	2/4/4/4	-
4	HEC	G	1486	1	-	8/14/54/54	-
4	HEC	E	1486	1	-	8/14/54/54	-
4	HEC	C	1485	2,1	-	8/14/54/54	-
4	HEC	A	1483	1	-	6/14/54/54	-
4	HEC	C	1483	1	-	6/14/54/54	-
4	HEC	E	1483	1	-	6/14/54/54	-
4	HEC	E	1482	5,1	-	6/14/54/54	-
3	GOL	A	1481	-	-	2/4/4/4	-
4	HEC	A	1485	2,1	-	8/14/54/54	-
4	HEC	C	1486	1	-	8/14/54/54	-
4	HEC	G	1482	5,1	-	6/14/54/54	-
4	HEC	G	1483	1	-	6/14/54/54	-
4	HEC	G	1484	2,1	-	4/14/54/54	-
4	HEC	C	1482	5,1	-	6/14/54/54	-
4	HEC	A	1484	2,1	-	6/14/54/54	-
3	GOL	G	1481	-	-	2/4/4/4	-
4	HEC	E	1484	2,1	-	4/14/54/54	-
4	HEC	G	1485	2,1	-	8/14/54/54	-
4	HEC	E	1485	2,1	-	8/14/54/54	-
4	HEC	C	1484	2,1	-	4/14/54/54	-
4	HEC	A	1482	5,1	-	6/14/54/54	-

The worst 5 of 179 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	1481	GOL	C3-C2	-7.43	1.23	1.51
3	G	1481	GOL	C3-C2	-7.43	1.23	1.51
3	A	1481	GOL	C3-C2	-7.32	1.23	1.51
3	E	1481	GOL	C3-C2	-7.17	1.24	1.51
4	E	1485	HEC	CAC-C3C	6.43	1.55	1.35

The worst 5 of 132 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed($^{\circ}$)	Ideal($^{\circ}$)
3	A	1481	GOL	O3-C3-C2	11.66	162.85	110.38
3	E	1481	GOL	O3-C3-C2	11.63	162.75	110.38
3	C	1481	GOL	O3-C3-C2	11.55	162.35	110.38
3	G	1481	GOL	O3-C3-C2	11.52	162.24	110.38
4	E	1482	HEC	CBB-CAB-C3B	-9.16	109.14	127.43

There are no chirality outliers.

5 of 138 torsion outliers are listed below:

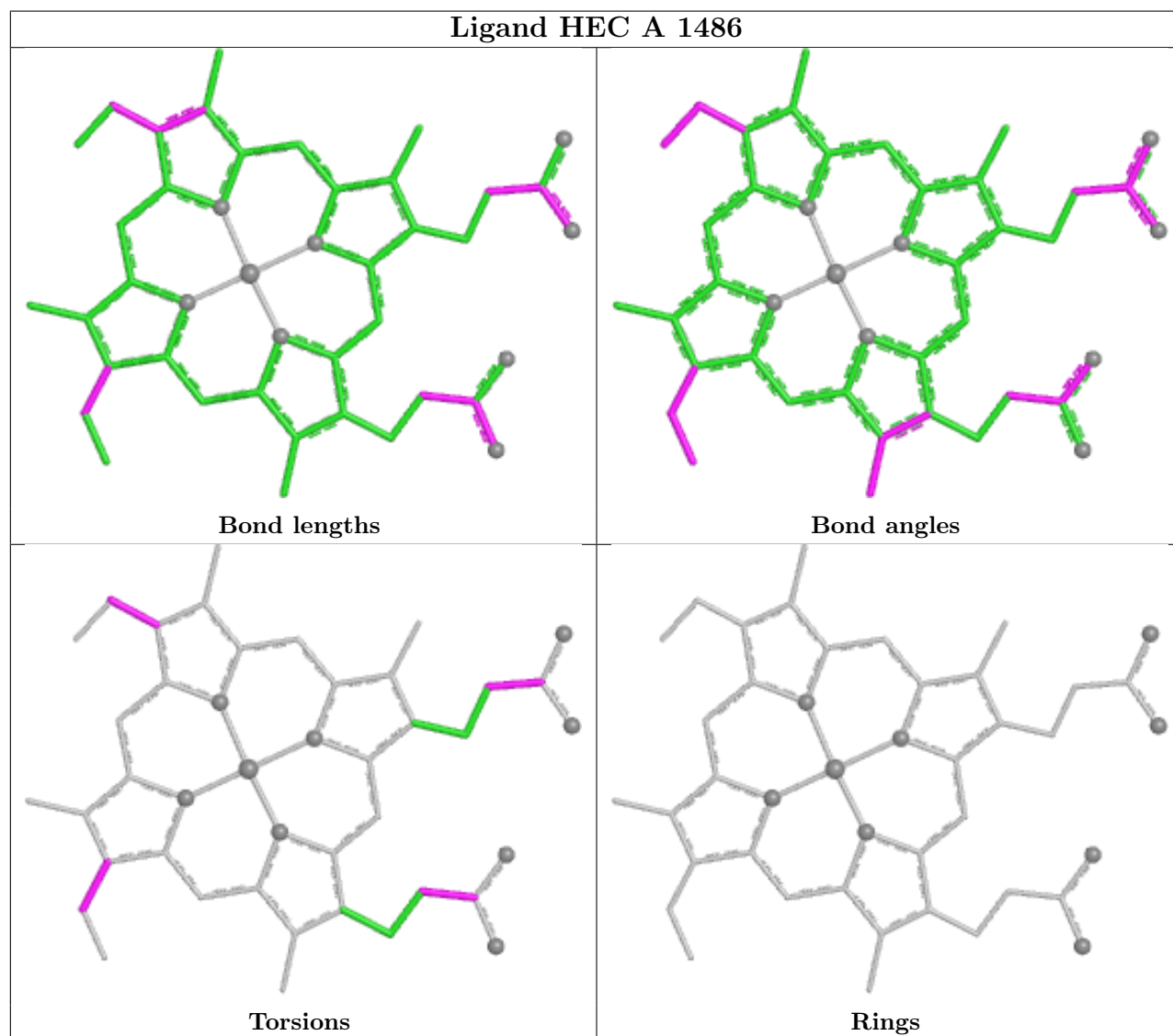
Mol	Chain	Res	Type	Atoms
3	A	1481	GOL	C1-C2-C3-O3
3	C	1481	GOL	C1-C2-C3-O3
3	E	1481	GOL	C1-C2-C3-O3
3	G	1481	GOL	C1-C2-C3-O3
4	A	1482	HEC	C2C-C3C-CAC-CBC

There are no ring outliers.

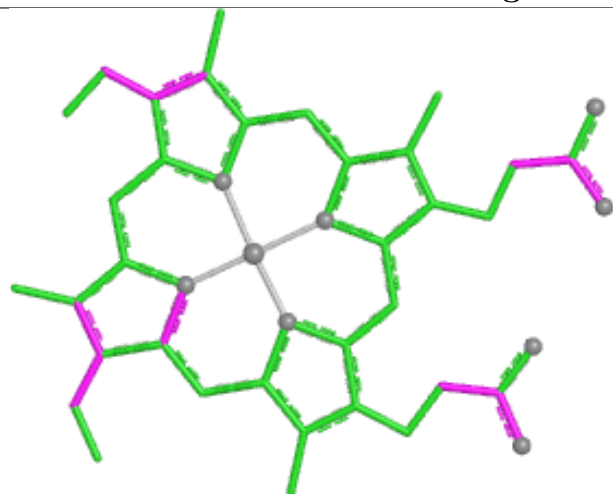
20 monomers are involved in 49 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	E	1481	GOL	3	0
4	A	1486	HEC	6	0
3	C	1481	GOL	3	0
4	G	1486	HEC	4	0
4	E	1486	HEC	4	0
4	C	1485	HEC	1	0
4	A	1483	HEC	2	0
4	C	1483	HEC	2	0
4	E	1483	HEC	2	0
3	A	1481	GOL	3	0
4	A	1485	HEC	2	0
4	C	1486	HEC	7	0
4	G	1483	HEC	2	0
4	G	1484	HEC	1	0
4	A	1484	HEC	1	0
3	G	1481	GOL	3	0
4	G	1485	HEC	2	0
4	E	1485	HEC	2	0
4	C	1484	HEC	2	0
4	A	1482	HEC	1	0

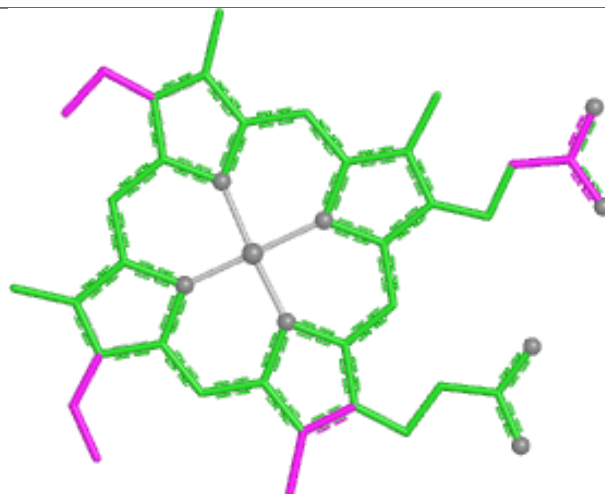
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included.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



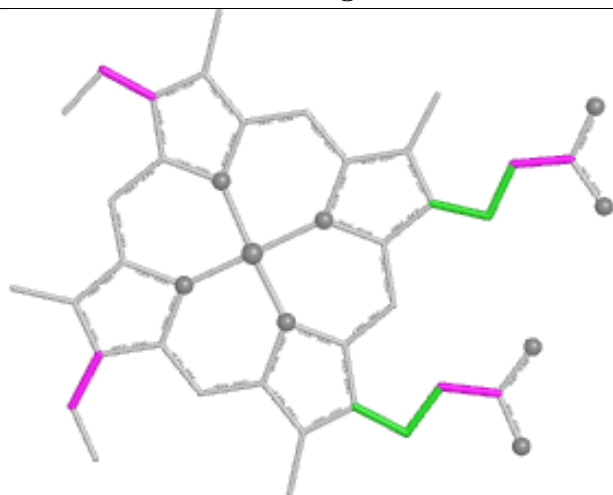
Ligand HEC G 1486



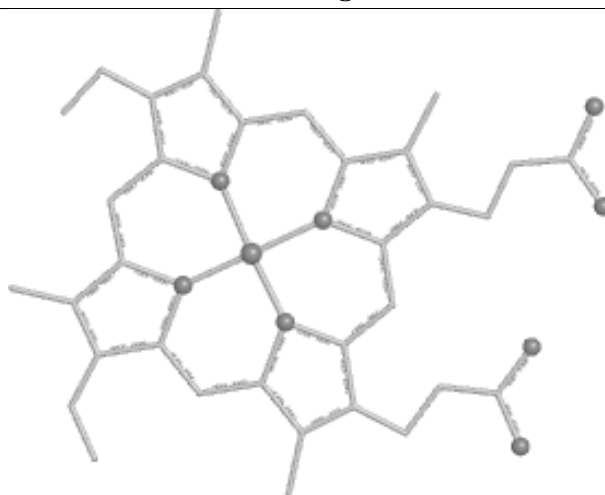
Bond lengths



Bond angles

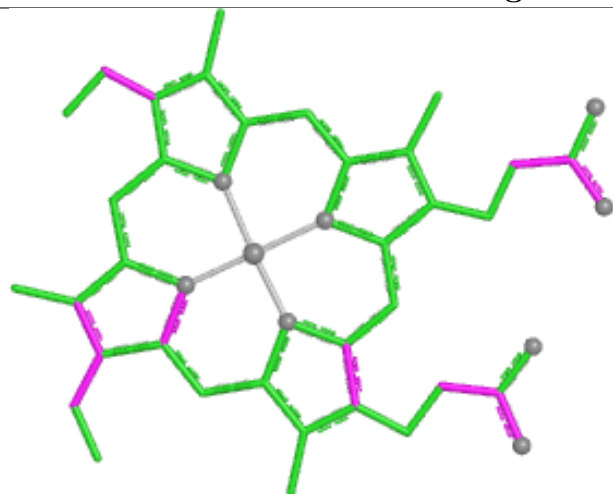


Torsions

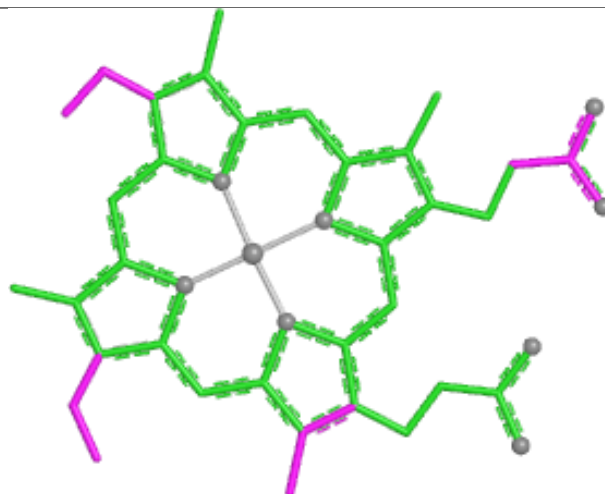


Rings

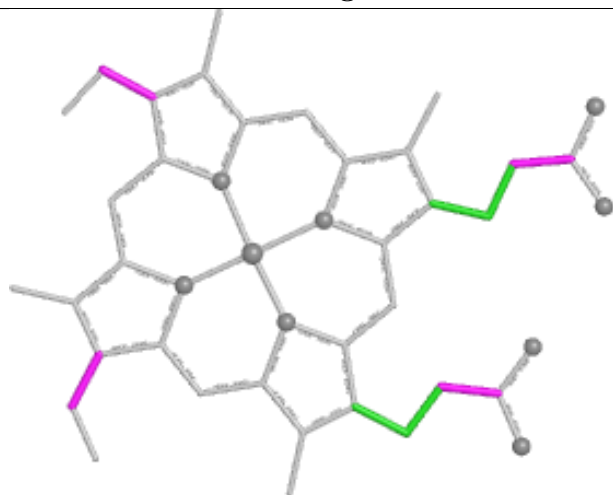
Ligand HEC E 1486



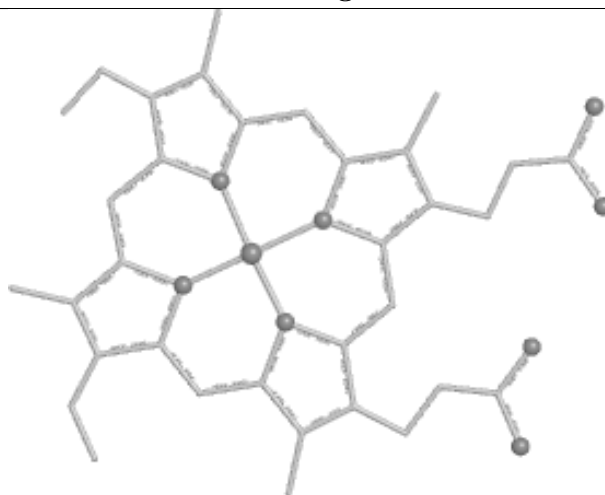
Bond lengths



Bond angles

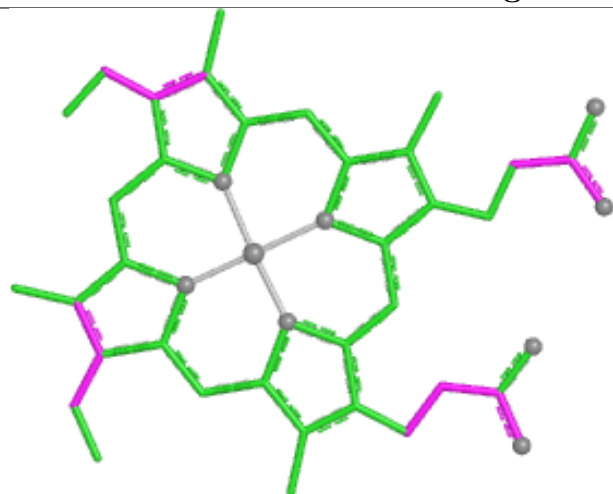


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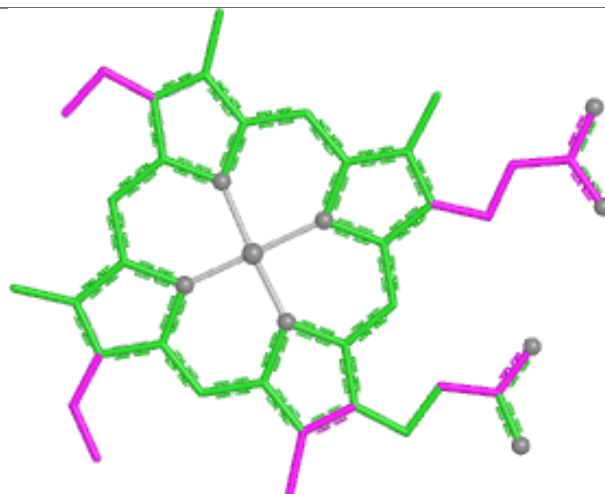


Rings

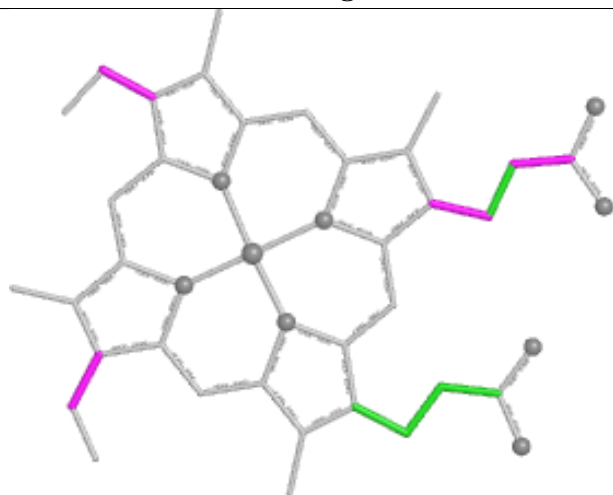
Ligand HEC C 1485



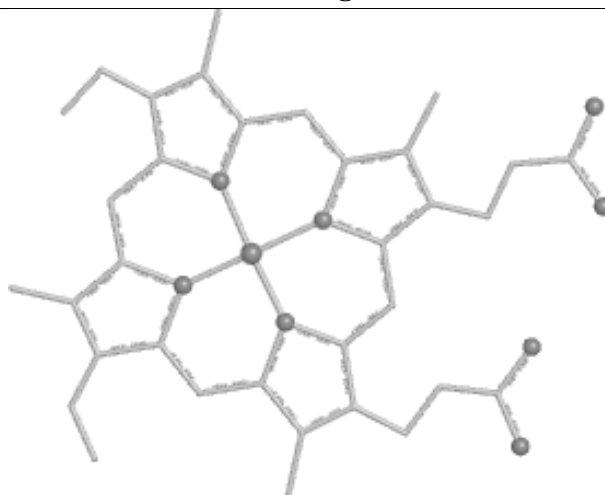
Bond lengths



Bond angles

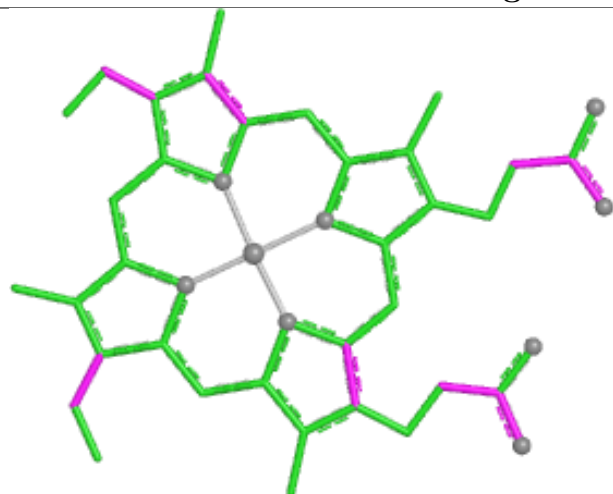


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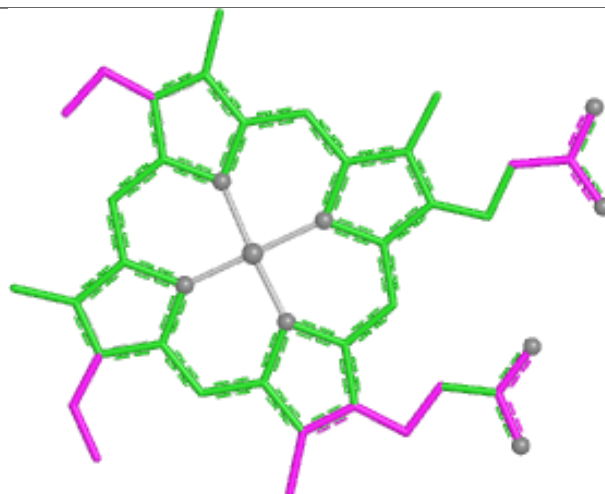


Rings

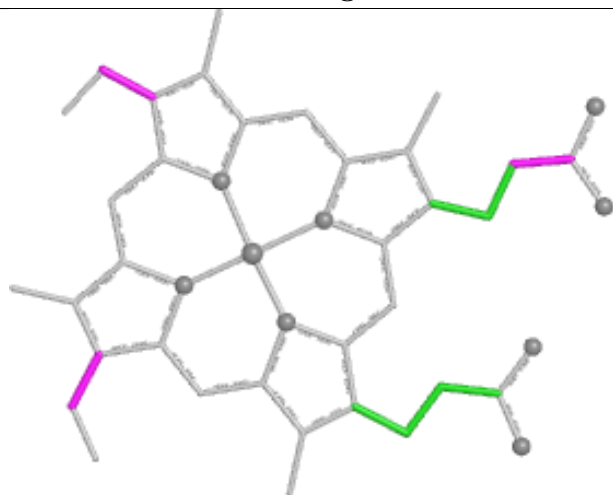
Ligand HEC A 1483



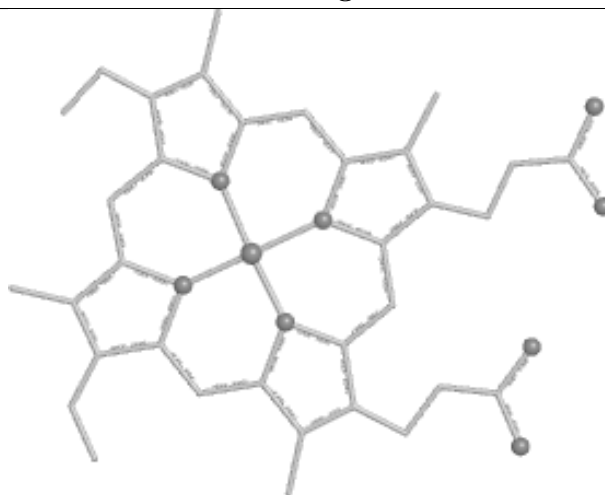
Bond lengths



Bond angles

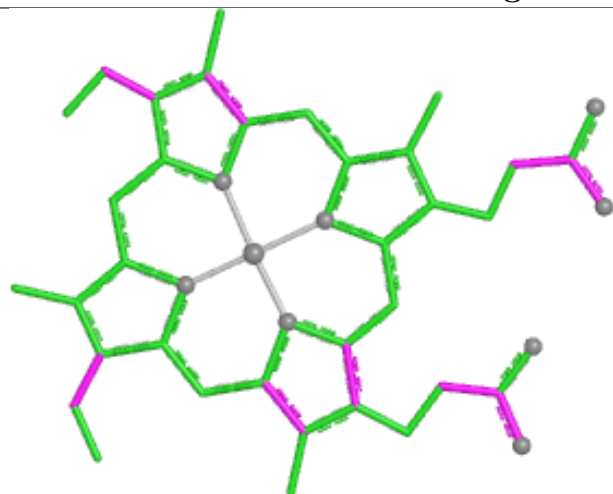


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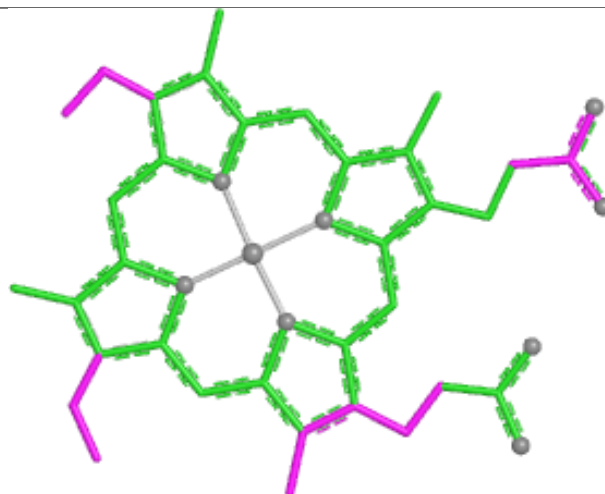


Rings

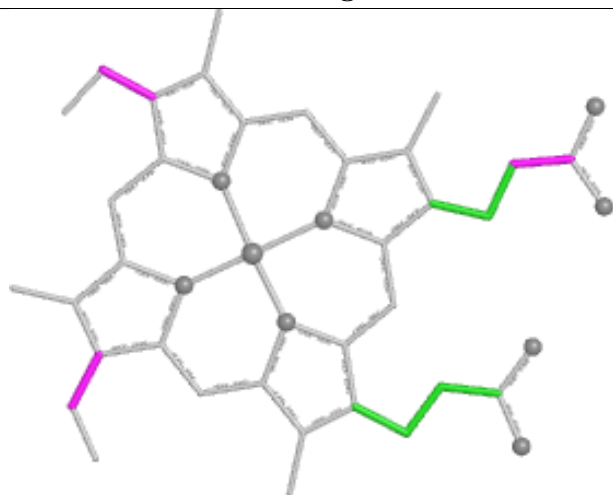
Ligand HEC C 1483



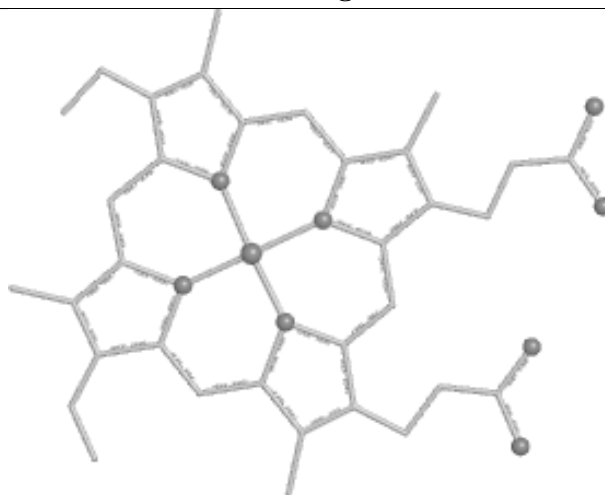
Bond lengths



Bond angles

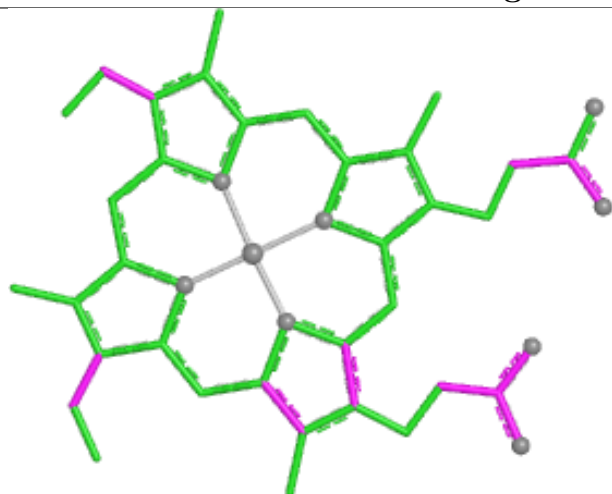


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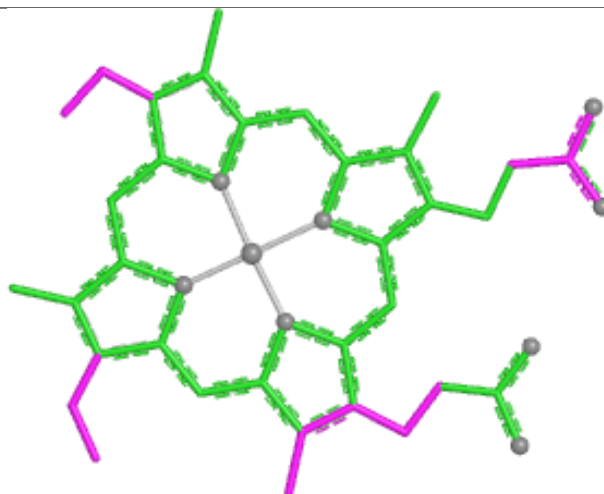


Rings

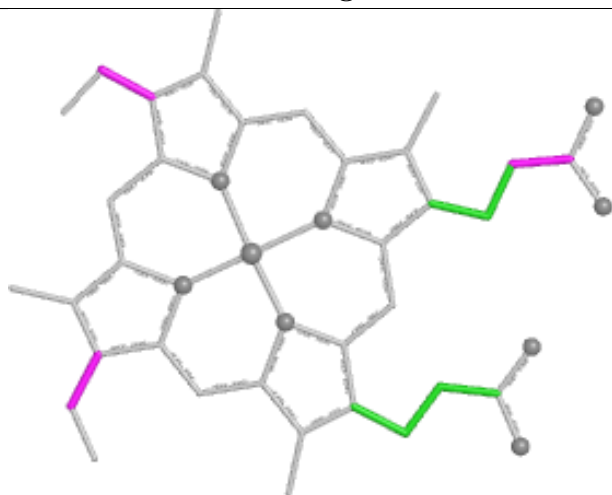
Ligand HEC E 1483



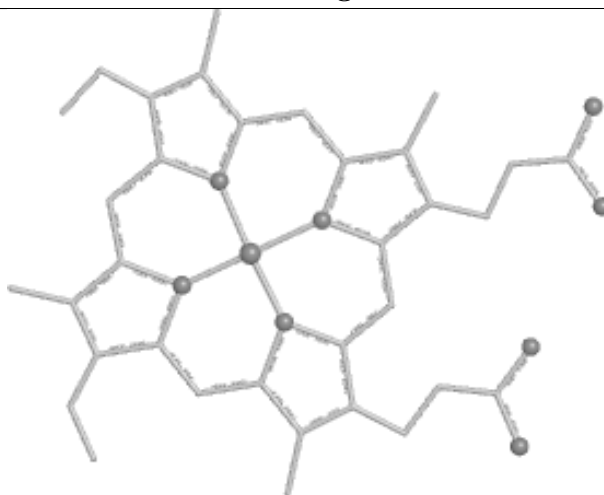
Bond lengths



Bond angles

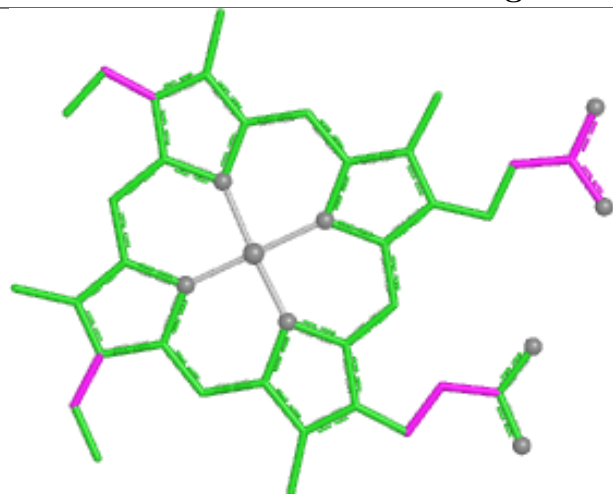


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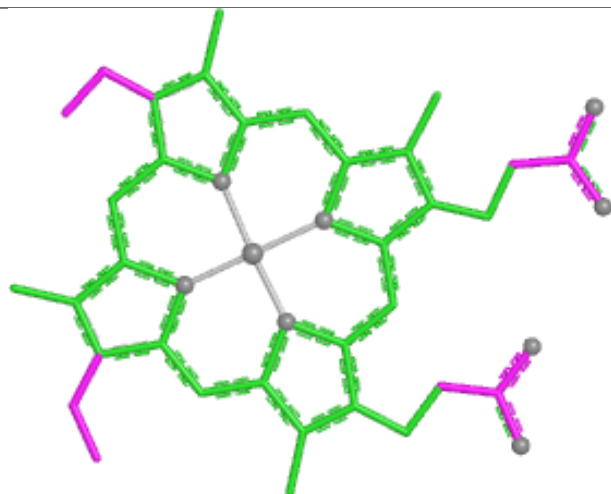


Rings

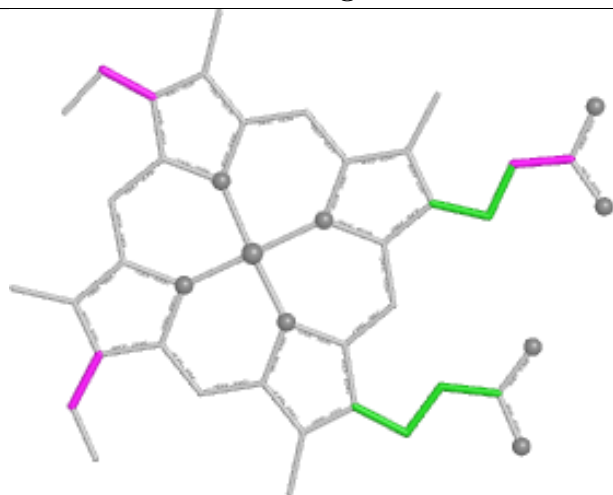
Ligand HEC E 1482



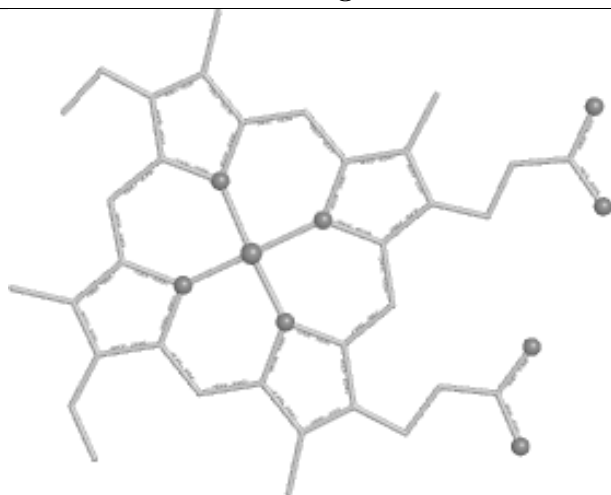
Bond lengths



Bond angles

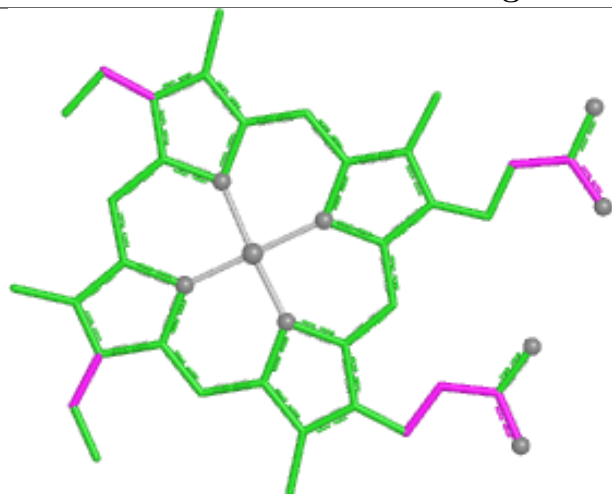


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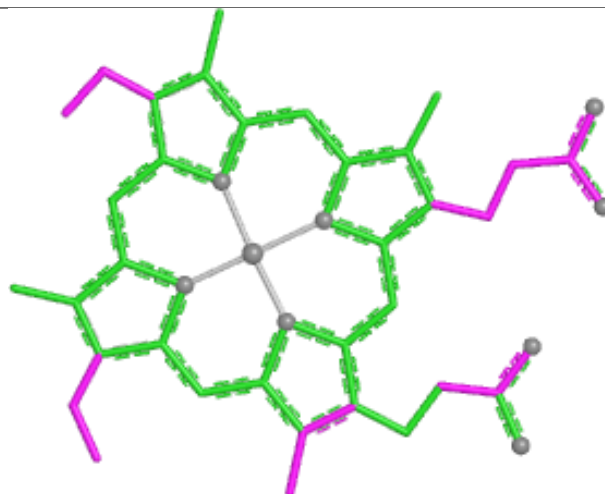


Rings

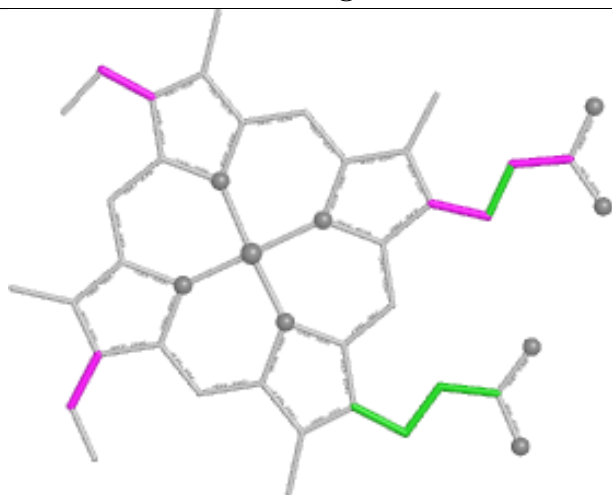
Ligand HEC A 1485



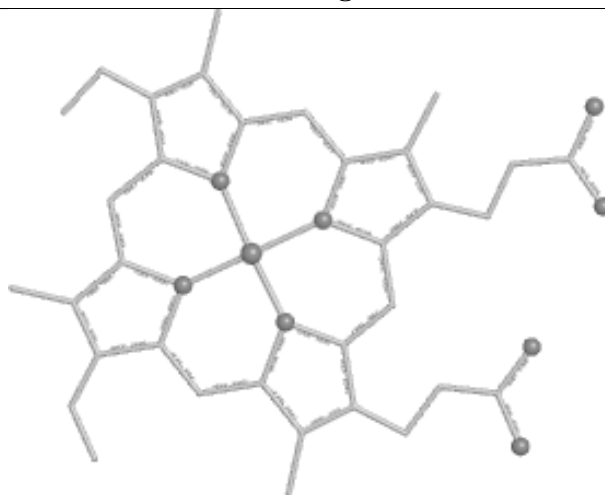
Bond lengths



Bond angles

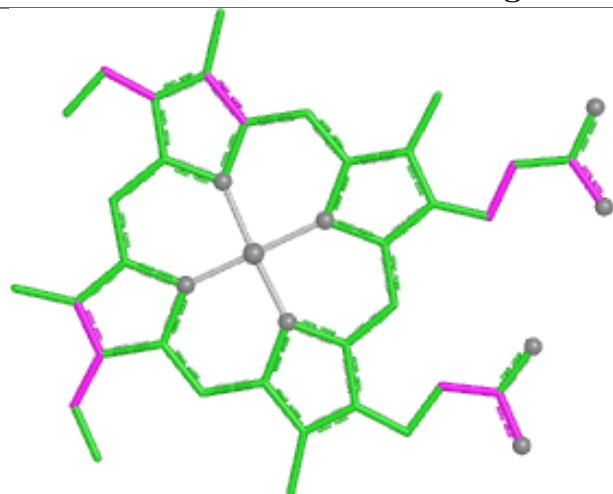


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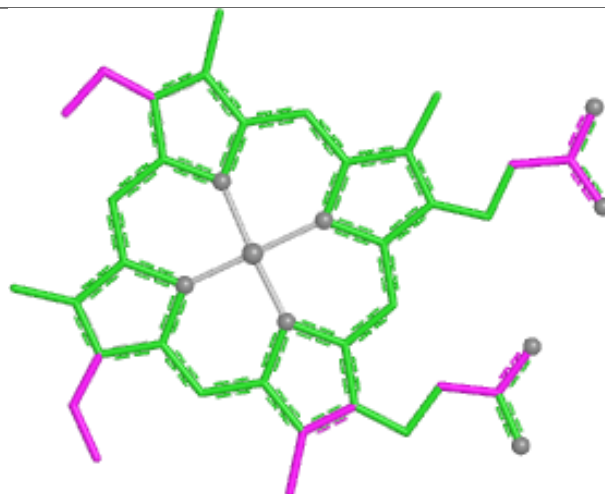


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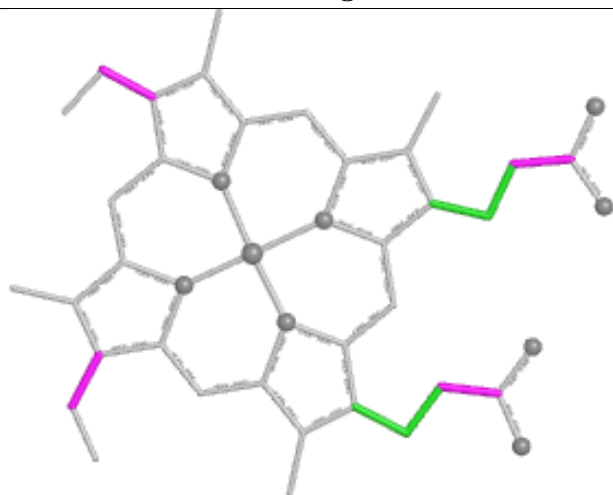
Ligand HEC C 1486



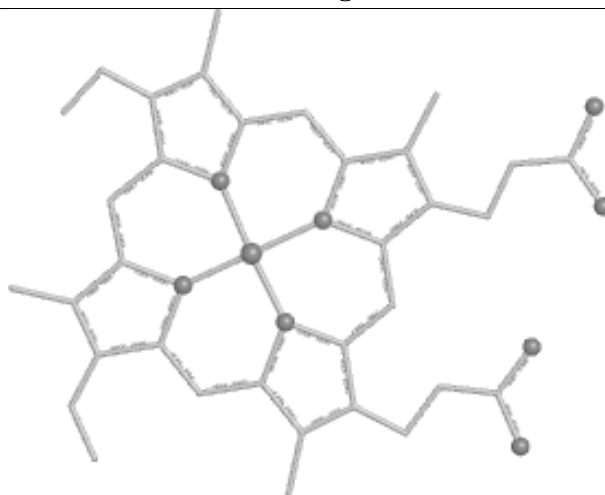
Bond lengths



Bond angles

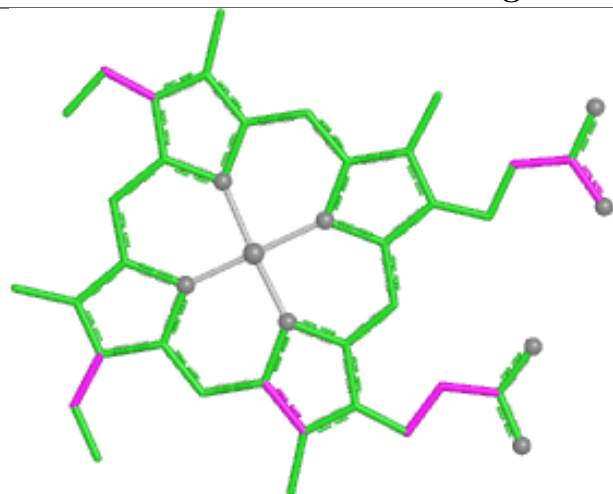


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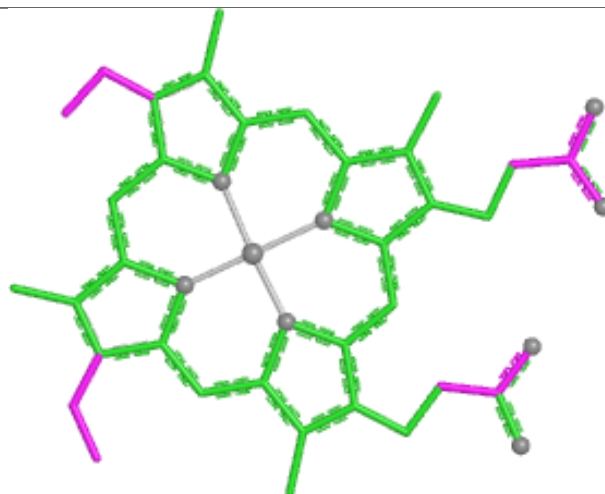


Rings

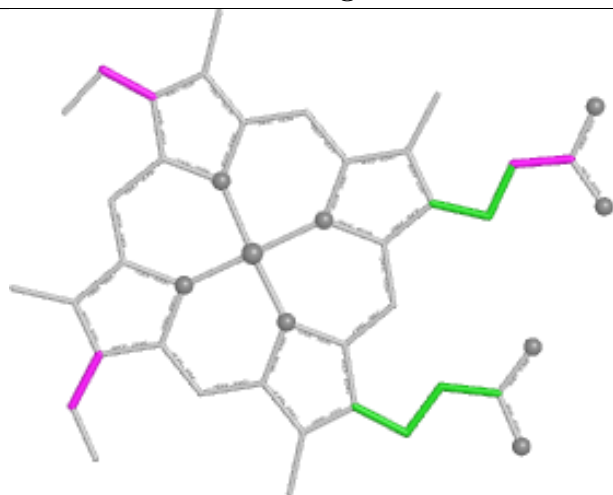
Ligand HEC G 1482



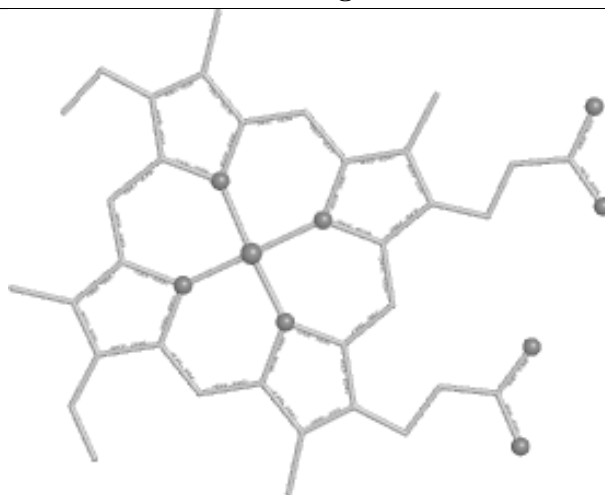
Bond lengths



Bond angles

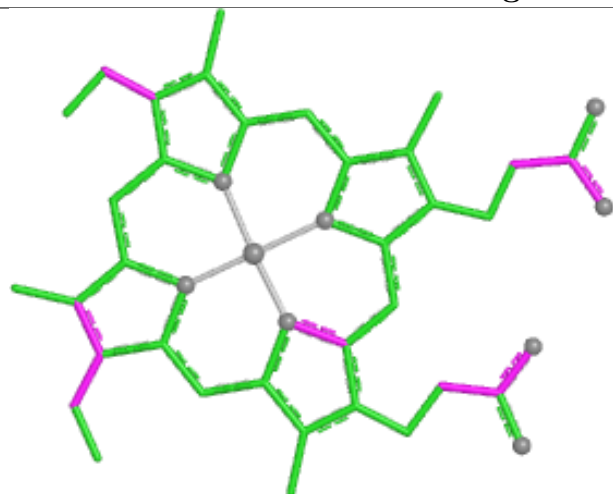


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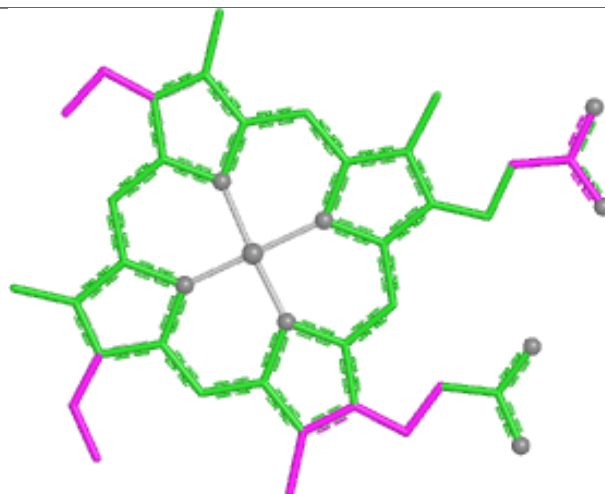


Rings

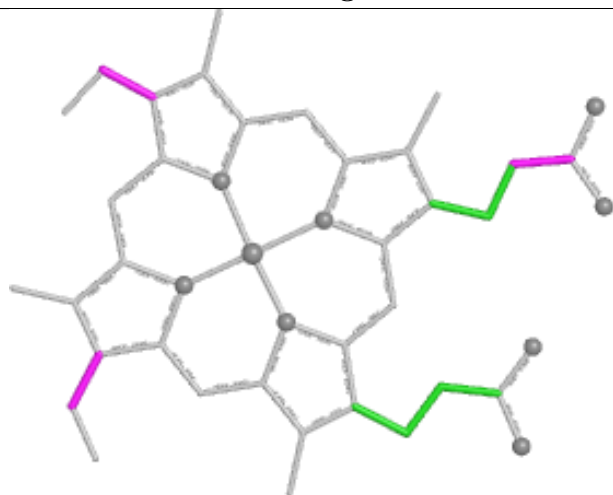
Ligand HEC G 1483



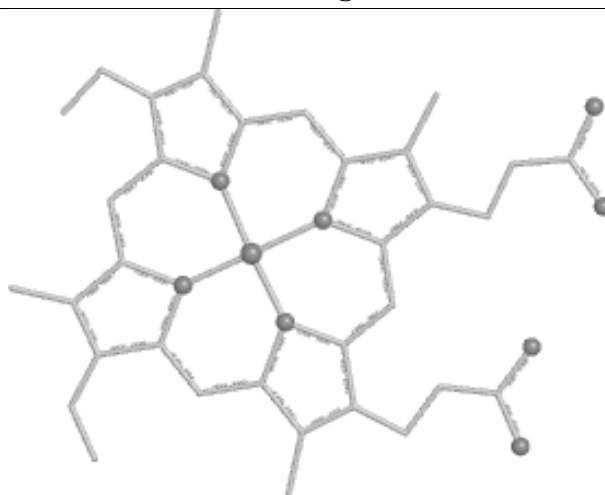
Bond lengths



Bond angles

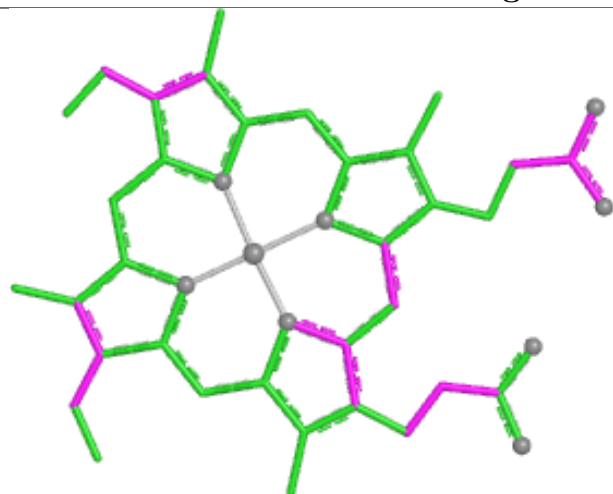


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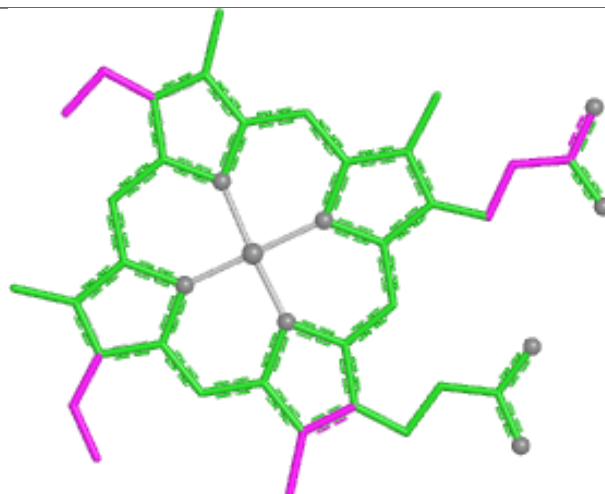


Rings

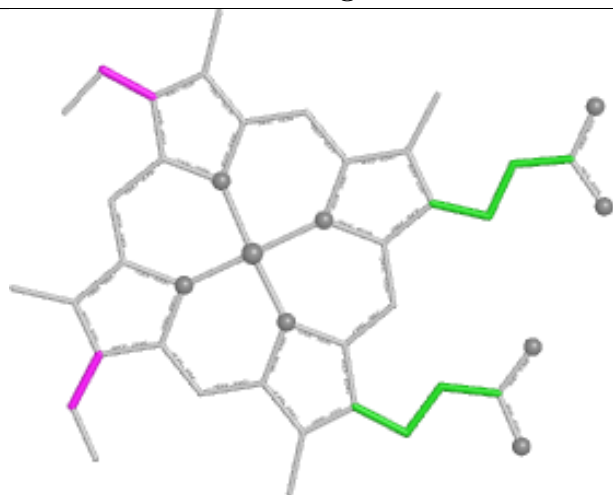
Ligand HEC G 1484



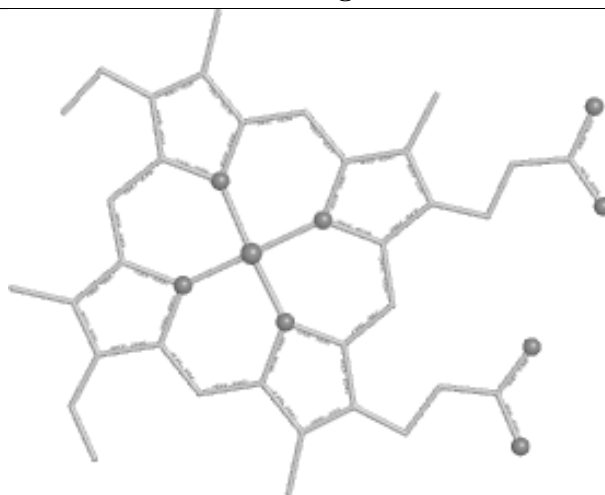
Bond lengths



Bond angles

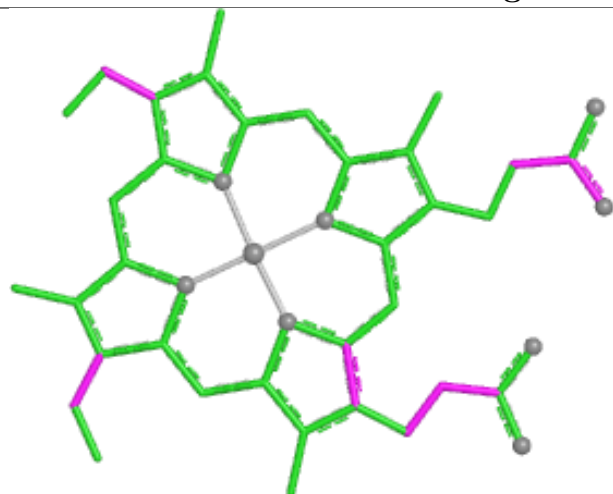


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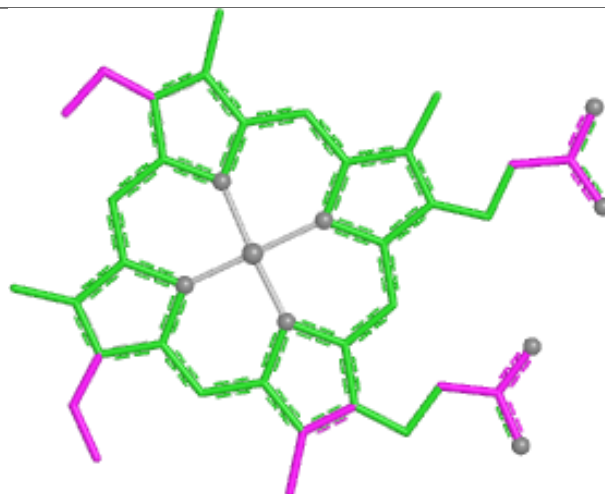


Rings

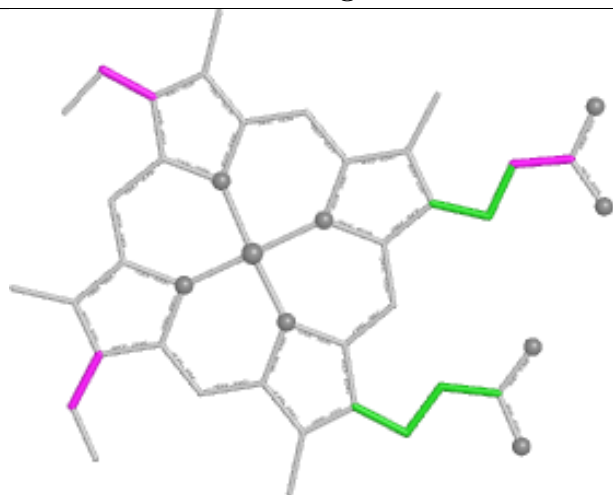
Ligand HEC C 1482



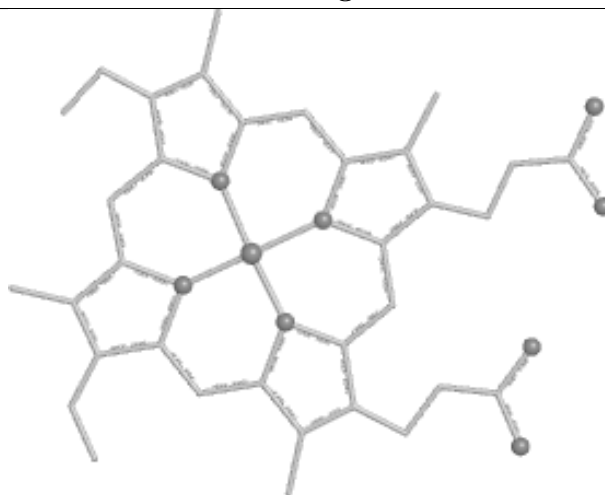
Bond lengths



Bond angles

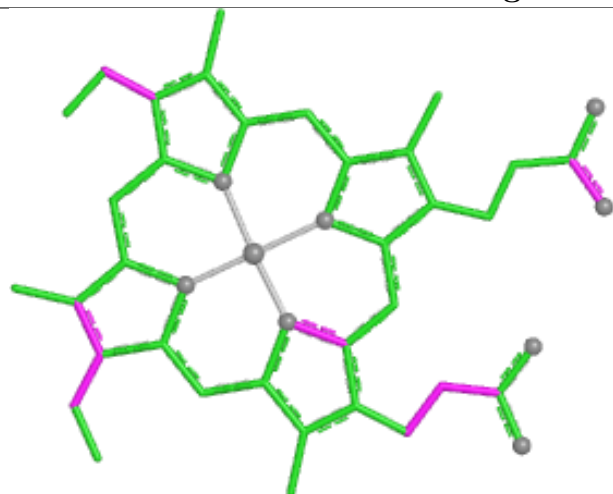


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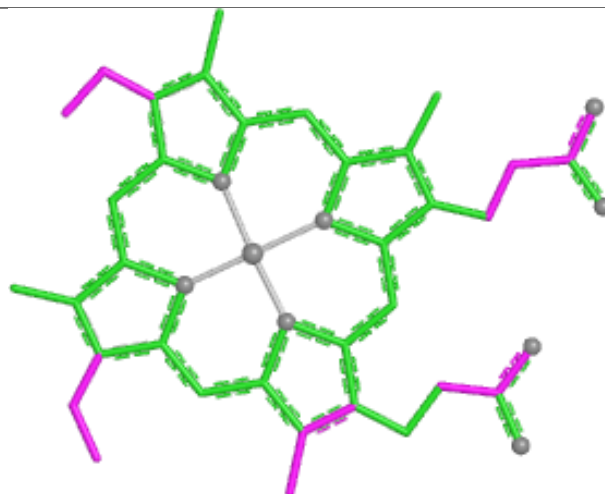


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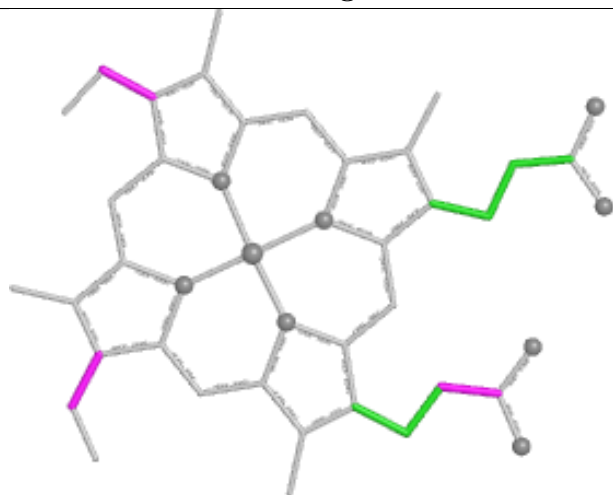
Ligand HEC A 1484



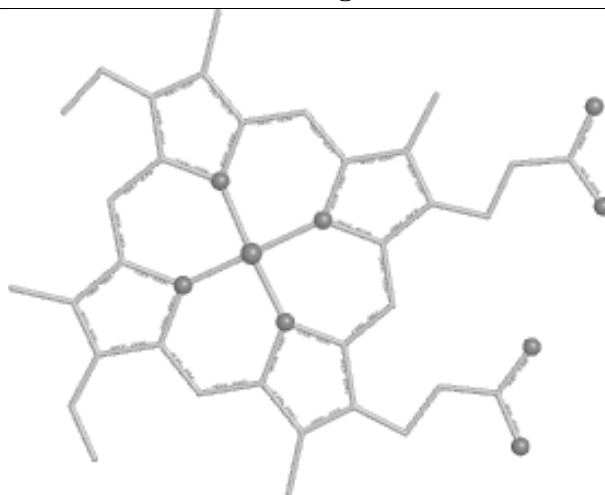
Bond lengths



Bond angles

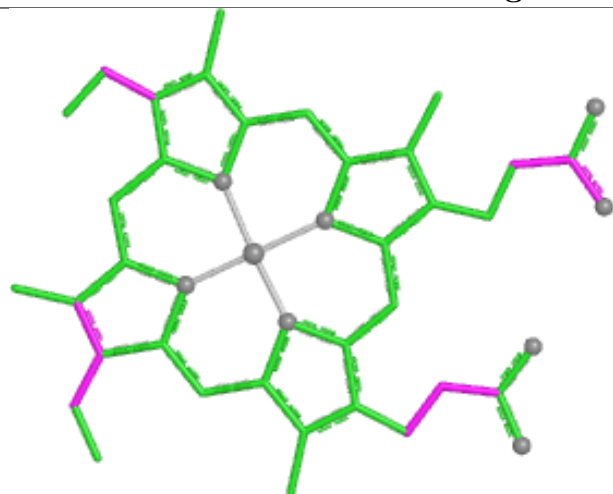


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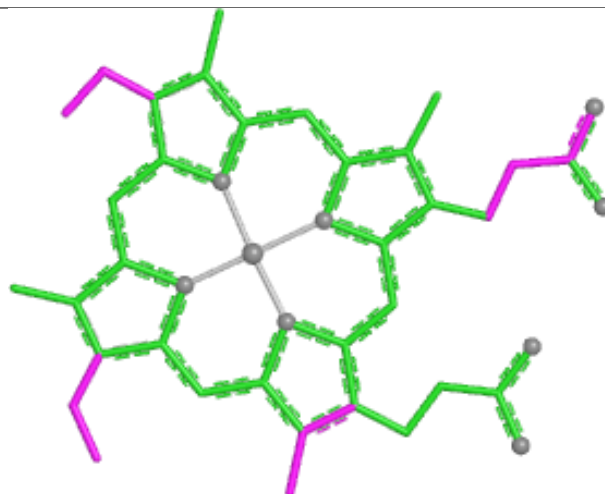


Rings

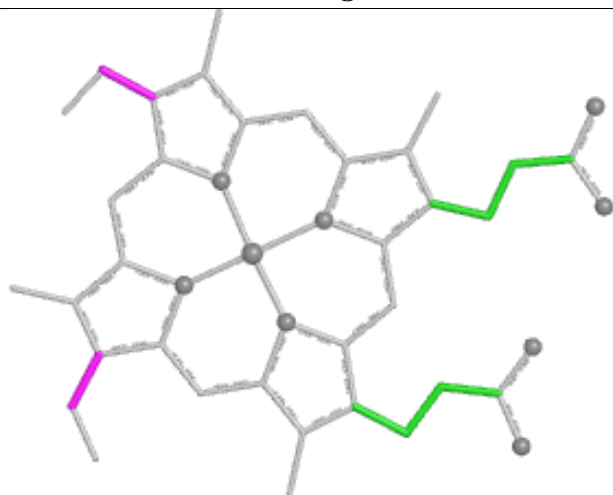
Ligand HEC E 1484



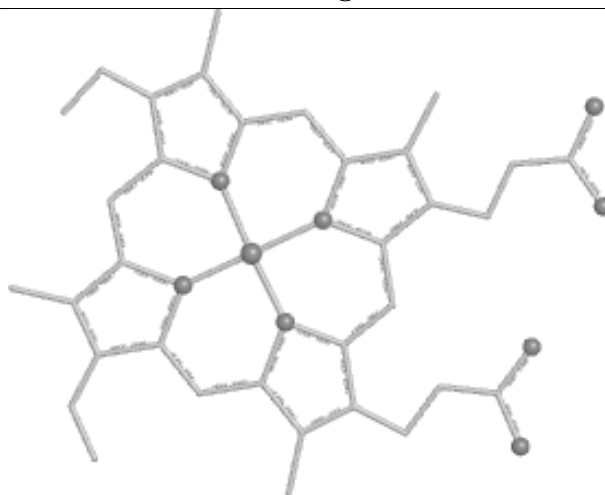
Bond lengths



Bond angles

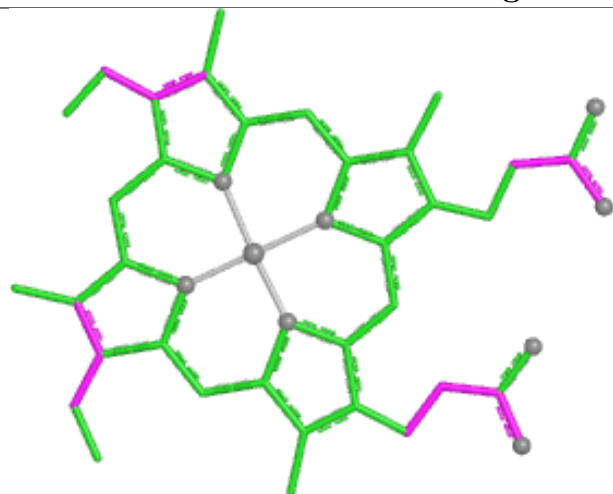


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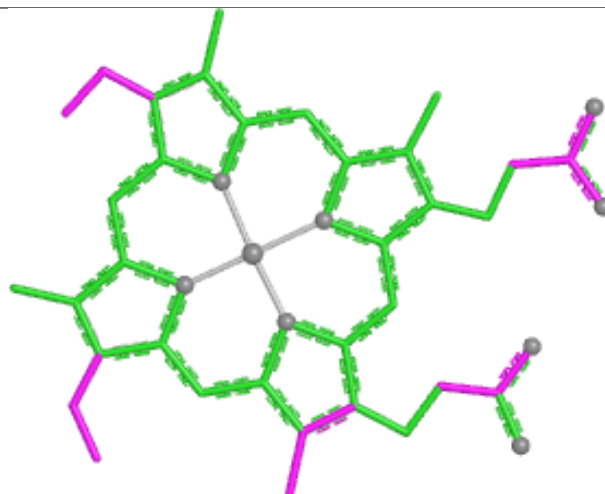


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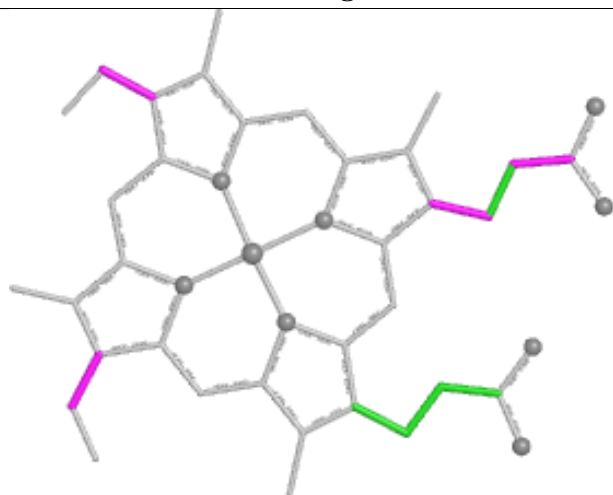
Ligand HEC G 1485



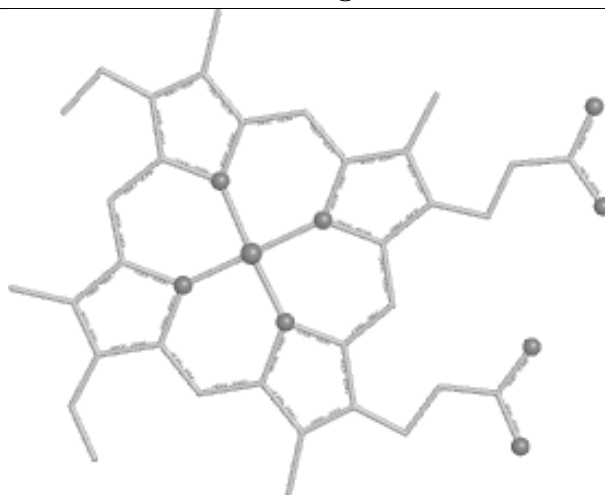
Bond lengths



Bond angles

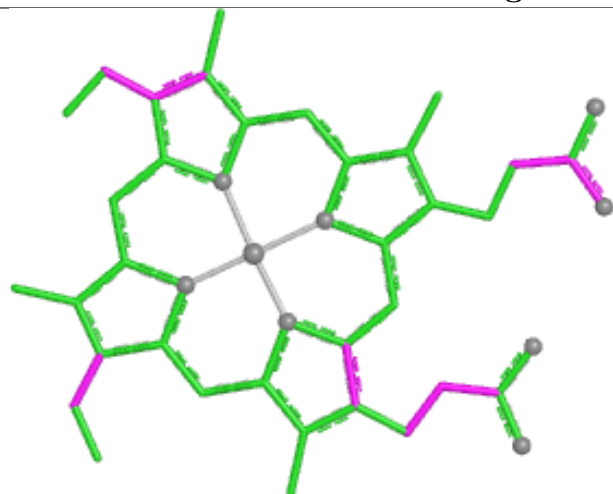


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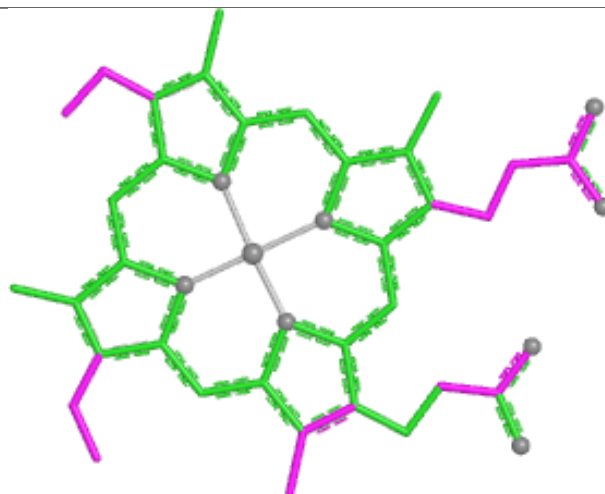


Rings

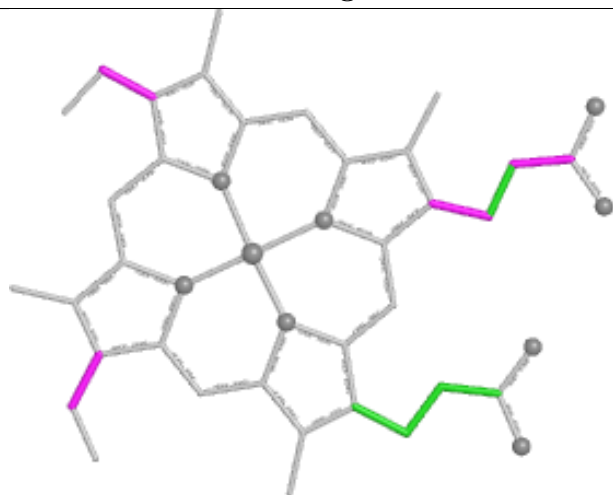
Ligand HEC E 1485



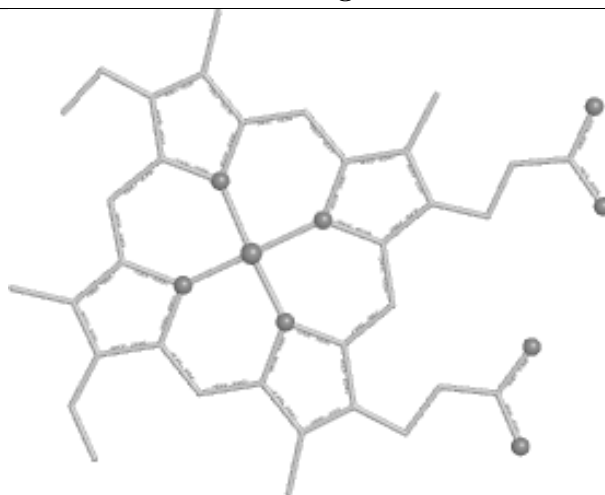
Bond lengths



Bond angles

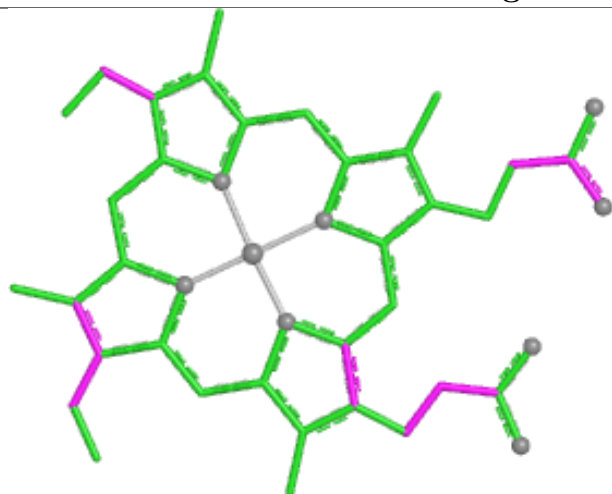


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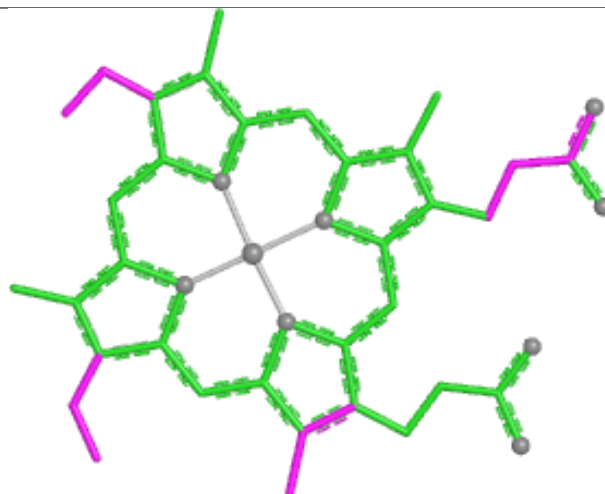


Rings

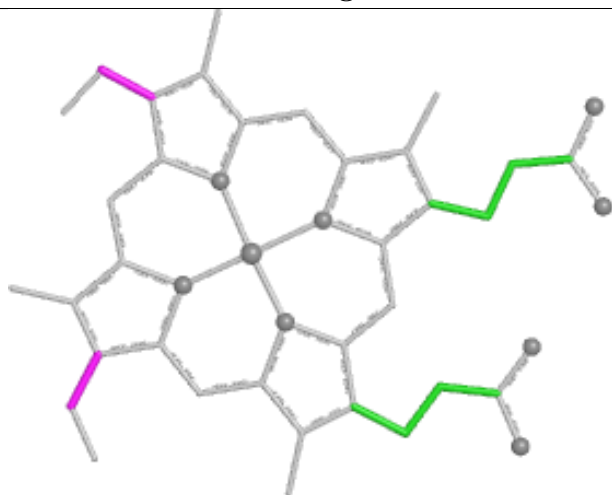
Ligand HEC C 1484



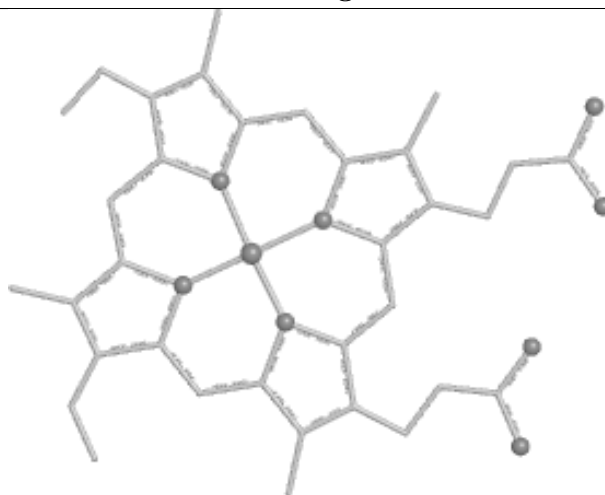
Bond lengths



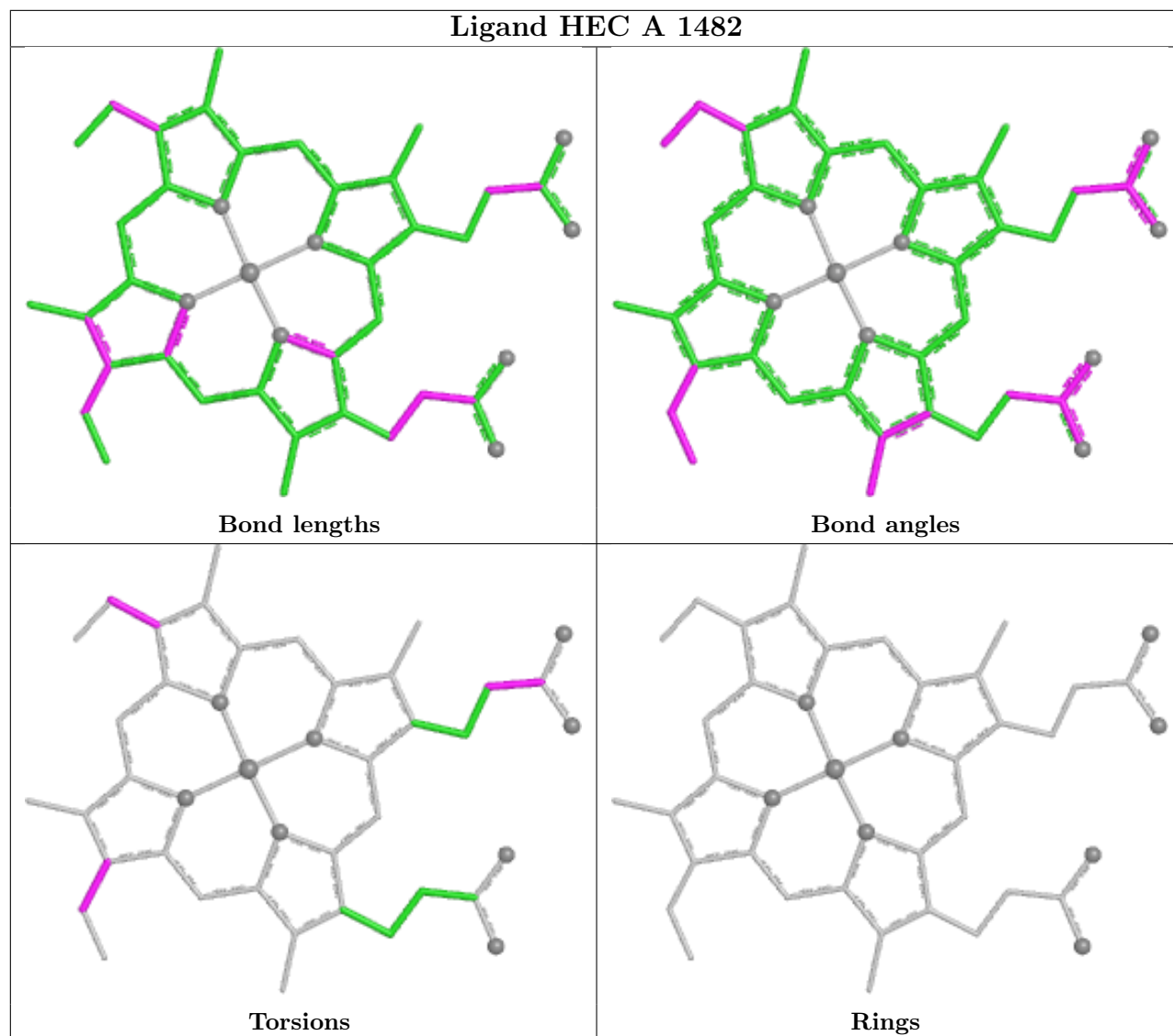
Bond angles



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 Fit of model and data [i](#)

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.