Full wwPDB X-ray Structure Validation Report (i)

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PDB ID	:	4N4J
Title	:	Kuenenia stuttgartiensis hydroxylamine oxidoreductase
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Deposited on	:	2013-10-08
Resolution	:	1.80 Å(reported)

This is a Full wwPDB X-ray Structure Validation 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 (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:

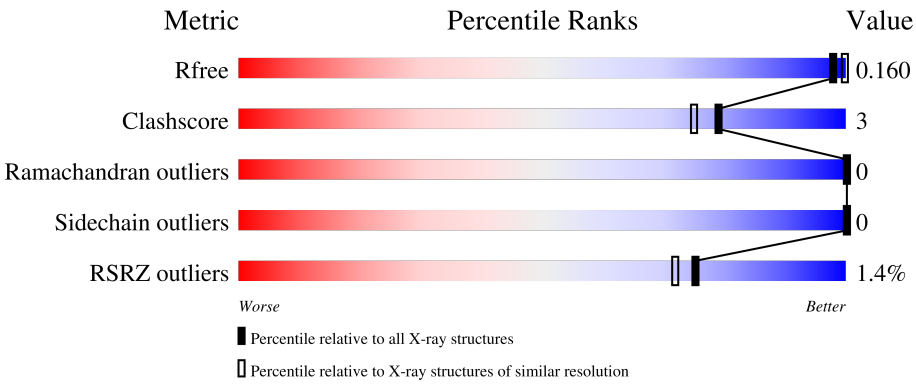
MolProbity	:	4.02b-467
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	1.20.1
EDS	:	2.37.1
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.37.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:
X-RAY DIFFRACTION

The reported resolution of this entry is 1.80 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 <=5% The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	500	% 97%

2 Entry composition (i)

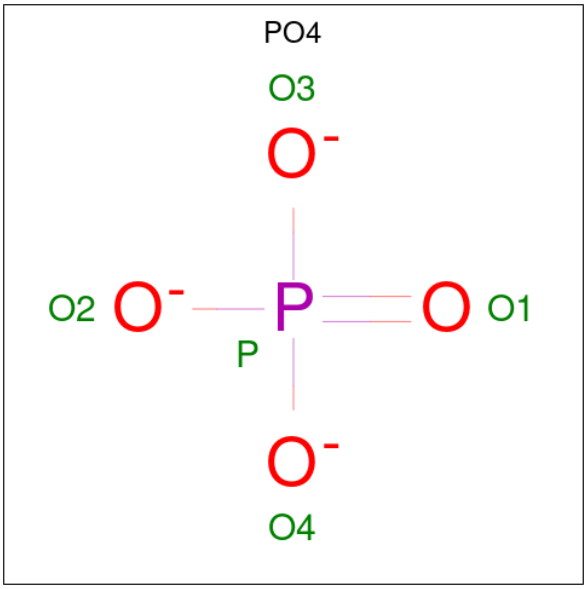
There are 5 unique types of molecules in this entry. The entry contains 4872 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 hydroxylamine oxidoreductase.

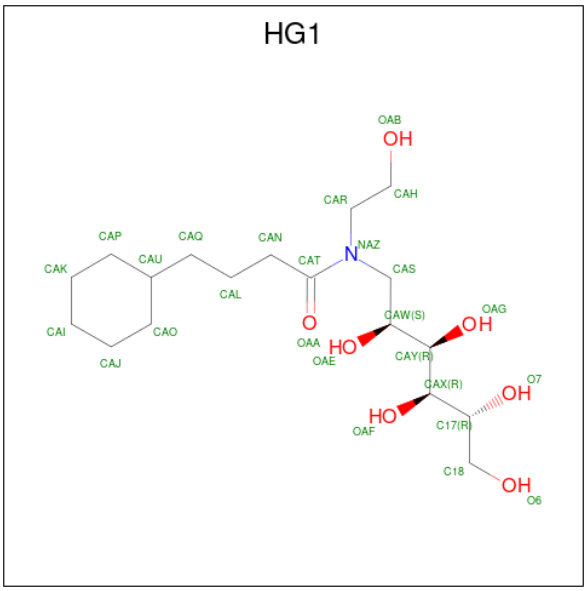
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	497	3979	2499	708	741	31	0	6	0

- Molecule 2 is PHOSPHATE ION (three-letter code: PO4) (formula: O_4P).



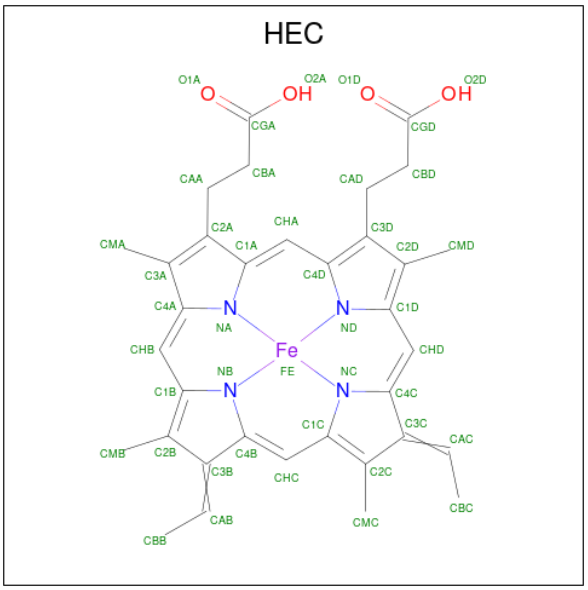
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	P		
2	A	1	5	4	1	0	0
2	A	1	5	4	1	0	0
2	A	1	5	4	1	0	0
2	A	1	5	4	1	0	0

- Molecule 3 is 1-[(4-cyclohexylbutanoyl)(2-hydroxyethyl)amino]-1-deoxy-D-glucitol (three-letter code: HG1) (formula: $C_{18}H_{35}NO_7$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total C N O 12 10 1 1	0	0
3	A	1	Total C N O 21 16 1 4	0	0
3	A	1	Total C N O 15 12 1 2	0	0
3	A	1	Total C 9 9	0	0

- Molecule 4 is HEME C (three-letter code: HEC) (formula: $C_{34}H_{34}FeN_4O_4$).



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

- Molecule 5 is water.

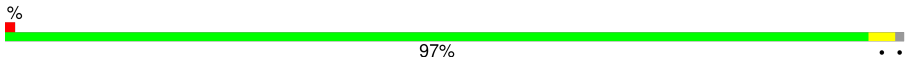
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	472	472	472	0	0

3 Residue-property plots (i)

These plots are drawn for all protein, RNA, DNA and oligosaccharide chains in the entry. The first graphic for a chain summarises the proportions of the various outlier classes displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. Residues are color-coded according to the number of geometric quality criteria for which they contain at least one outlier: green = 0, yellow = 1, orange = 2 and red = 3 or more. A red dot above a residue indicates a poor fit to the electron density (RSRZ > 2). Stretches of 2 or more consecutive residues without any outlier are shown as a green connector. Residues present in the sample, but not in the model, are shown in grey.

• Molecule 1: hydroxylamine oxidoreductase

Chain A:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α, β, γ	130.00Å 130.00Å 130.00Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.96 – 1.80 45.96 – 1.80	Depositor EDS
% Data completeness (in resolution range)	100.0 (45.96-1.80) 100.0 (45.96-1.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	4.30 (at 1.81Å)	Xtriage
Refinement program	REFMAC 5.5.010	Depositor
R, R_{free}	0.143 , 0.157 0.150 , 0.160	Depositor DCC
R_{free} test set	3340 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	15.7	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.40 , 42.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.31$	Xtriage
Estimated twinning fraction	0.035 for l,-k,h	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	4872	wwPDB-VP
Average B, all atoms (Å^2)	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.43% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

[2]Theoretical values of $< |L| >$, $< L^2 >$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

5 Model quality (i)

5.1 Standard geometry (i)

Bond lengths and bond angles in the following residue types are not validated in this section: HG1, PO4, HEC

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	0.33	0/4097	0.53	0/5545

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

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

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3979	0	3758	10	2
2	A	20	0	0	0	0
3	A	57	0	81	1	0
4	A	344	0	240	20	2
5	A	472	0	0	0	0
All	All	4872	0	4079	24	2

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 3.

All (24) 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)
4:A:612:HEC:HMC1	4:A:612:HEC:HBC3	1.69	0.75
4:A:609:HEC:HBC3	4:A:609:HEC:HMC1	1.71	0.71
1:A:98:GLU:OE1	1:A:177:LYS:HD2	1.98	0.62
4:A:613:HEC:HMC1	4:A:613:HEC:HBC3	1.84	0.59
4:A:609:HEC:HBB3	4:A:609:HEC:HMB1	1.85	0.58
4:A:616:HEC:HMC1	4:A:616:HEC:HBC3	1.90	0.54
1:A:444:MET:CE	4:A:616:HEC:HMB2	2.37	0.54
4:A:612:HEC:HMC1	4:A:612:HEC:CBC	2.38	0.54
4:A:609:HEC:HMC1	4:A:609:HEC:CBC	2.39	0.53
4:A:614:HEC:HMB1	4:A:614:HEC:HBB3	1.92	0.50
4:A:612:HEC:HAD2	4:A:614:HEC:HBD2	1.96	0.47
1:A:180:ASN:HB2	1:A:187:THR:HG21	1.98	0.46
4:A:614:HEC:HMC1	4:A:614:HEC:HBC3	1.98	0.45
1:A:300:THR:C	4:A:615:HEC:HMC3	2.36	0.45
1:A:234:CYS:HB2	4:A:611:HEC:HMB1	1.98	0.45
4:A:610:HEC:HMC1	4:A:610:HEC:HBC3	1.99	0.44
1:A:205:ASN:O	4:A:612:HEC:HHB	2.19	0.43
1:A:273:ILE:HD12	4:A:616:HEC:HAA2	2.01	0.42
1:A:62:TYR:CG	1:A:68:ALA:HB2	2.54	0.42
1:A:414:PRO:HB2	3:A:606:HG1:H4	2.01	0.42
4:A:613:HEC:HMB1	4:A:613:HEC:HBB3	2.02	0.42
4:A:615:HEC:HMB1	4:A:615:HEC:HBB3	2.02	0.41
1:A:198:HIS:CD2	4:A:614:HEC:ND	2.88	0.41
4:A:615:HEC:HBC3	4:A:615:HEC:HMC1	2.02	0.40

All (2) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:451:TYR:OH	4:A:612:HEC:C4B[12_455]	1.49	0.71
1:A:451:TYR:CE2	4:A:612:HEC:CHC[12_455]	1.51	0.69

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	501/500 (100%)	491 (98%)	10 (2%)	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 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	419/416 (101%)	419 (100%)	0	100 100

There are no protein residues with a non-rotameric sidechain to report.

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 monosaccharides in this entry.

5.6 Ligand geometry (i)

16 ligands are modelled in this entry.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	HEC	A	611	1	32,50,50	2.19	5 (15%)	30,82,82	2.03	9 (30%)
4	HEC	A	613	1	32,50,50	2.30	5 (15%)	30,82,82	2.12	7 (23%)
2	PO4	A	604	-	4,4,4	0.86	0	6,6,6	0.51	0
4	HEC	A	615	1	32,50,50	2.17	5 (15%)	30,82,82	2.02	8 (26%)
4	HEC	A	609	1	32,50,50	2.22	5 (15%)	30,82,82	1.76	5 (16%)
3	HG1	A	608	-	9,9,26	0.47	0	10,10,33	0.58	0
4	HEC	A	614	1	32,50,50	2.25	5 (15%)	30,82,82	1.75	6 (20%)
4	HEC	A	610	1	32,50,50	2.27	5 (15%)	30,82,82	1.60	6 (20%)
2	PO4	A	602	-	4,4,4	0.99	0	6,6,6	0.68	0
2	PO4	A	601	-	4,4,4	0.97	0	6,6,6	0.66	0
3	HG1	A	605	-	12,12,26	0.24	0	14,14,33	0.76	0
3	HG1	A	607	-	15,15,26	0.57	0	17,17,33	0.47	0
4	HEC	A	612	1,5	32,50,50	2.50	12 (37%)	30,82,82	2.77	15 (50%)
2	PO4	A	603	-	4,4,4	1.15	0	6,6,6	0.33	0
4	HEC	A	616	1	32,50,50	2.26	5 (15%)	30,82,82	1.82	7 (23%)
3	HG1	A	606	-	21,21,26	0.61	0	23,25,33	0.88	1 (4%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	HEC	A	613	1	-	0/10/54/54	-
4	HEC	A	615	1	-	1/10/54/54	-
3	HG1	A	608	-	-	1/3/11/39	0/1/1/1
4	HEC	A	614	1	-	4/10/54/54	-
4	HEC	A	610	1	-	0/10/54/54	-
4	HEC	A	611	1	-	3/10/54/54	-
3	HG1	A	605	-	-	0/6/14/39	0/1/1/1
3	HG1	A	607	-	-	3/10/18/39	0/1/1/1
4	HEC	A	612	1,5	-	3/10/54/54	-
4	HEC	A	609	1	-	3/10/54/54	-
4	HEC	A	616	1	-	3/10/54/54	-

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Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HG1	A	606	-	-	6/20/28/39	0/1/1/1

All (47) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	612	HEC	C4B-C3B	6.61	1.55	1.43
4	A	616	HEC	C3C-C2C	-6.27	1.33	1.40
4	A	613	HEC	C2B-C3B	-6.21	1.33	1.40
4	A	610	HEC	C2B-C3B	-6.21	1.33	1.40
4	A	613	HEC	C3C-C2C	-6.12	1.33	1.40
4	A	610	HEC	C3C-C2C	-6.06	1.33	1.40
4	A	615	HEC	C2B-C3B	-5.95	1.34	1.40
4	A	611	HEC	C3C-C2C	-5.94	1.34	1.40
4	A	609	HEC	C2B-C3B	-5.90	1.34	1.40
4	A	614	HEC	C2B-C3B	-5.90	1.34	1.40
4	A	614	HEC	C3C-C2C	-5.78	1.34	1.40
4	A	616	HEC	C2B-C3B	-5.75	1.34	1.40
4	A	611	HEC	C2B-C3B	-5.64	1.34	1.40
4	A	615	HEC	C3C-C2C	-5.61	1.34	1.40
4	A	612	HEC	C2A-C1A	5.33	1.54	1.42
4	A	609	HEC	C3C-C2C	-5.25	1.34	1.40
4	A	613	HEC	C3D-C2D	5.06	1.52	1.37
4	A	609	HEC	C3D-C2D	5.04	1.52	1.37
4	A	614	HEC	C3D-C2D	4.97	1.52	1.37
4	A	616	HEC	C3D-C2D	4.89	1.52	1.37
4	A	615	HEC	C3D-C2D	4.87	1.52	1.37
4	A	611	HEC	C3D-C2D	4.82	1.52	1.37
4	A	610	HEC	C3D-C2D	4.77	1.51	1.37
4	A	612	HEC	C3C-C4C	4.63	1.51	1.43
4	A	609	HEC	CBC-CAC	-4.07	1.34	1.49
4	A	609	HEC	CBB-CAB	-4.06	1.34	1.49
4	A	616	HEC	CBB-CAB	-4.02	1.34	1.49
4	A	614	HEC	CBC-CAC	-4.01	1.34	1.49
4	A	616	HEC	CBC-CAC	-3.99	1.34	1.49
4	A	613	HEC	CBB-CAB	-3.99	1.34	1.49
4	A	610	HEC	CBB-CAB	-3.96	1.34	1.49
4	A	613	HEC	CBC-CAC	-3.95	1.34	1.49
4	A	610	HEC	CBC-CAC	-3.94	1.34	1.49
4	A	611	HEC	CBC-CAC	-3.93	1.34	1.49
4	A	614	HEC	CBB-CAB	-3.92	1.34	1.49
4	A	612	HEC	C2B-C3B	-3.92	1.36	1.40
4	A	615	HEC	CBC-CAC	-3.91	1.35	1.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	615	HEC	CBB-CAB	-3.87	1.35	1.49
4	A	611	HEC	CBB-CAB	-3.82	1.35	1.49
4	A	612	HEC	C3C-C2C	-3.58	1.36	1.40
4	A	612	HEC	CBD-CGD	2.96	1.57	1.50
4	A	612	HEC	C4D-ND	2.82	1.42	1.36
4	A	612	HEC	O1A-CGA	2.78	1.31	1.22
4	A	612	HEC	C1C-CHC	2.66	1.48	1.41
4	A	612	HEC	C1D-ND	2.57	1.41	1.36
4	A	612	HEC	O2A-CGA	-2.54	1.22	1.30
4	A	612	HEC	C3A-C4A	2.18	1.47	1.42

All (64) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	612	HEC	O2A-CGA-CBA	6.16	133.46	114.00
4	A	613	HEC	CBD-CAD-C3D	-5.84	102.72	112.54
4	A	612	HEC	CBC-CAC-C3C	-5.69	114.18	127.49
4	A	616	HEC	CBB-CAB-C3B	-5.29	115.11	127.49
4	A	613	HEC	CBB-CAB-C3B	-5.07	115.62	127.49
4	A	615	HEC	CBB-CAB-C3B	-4.98	115.84	127.49
4	A	612	HEC	CBB-CAB-C3B	-4.47	117.03	127.49
4	A	611	HEC	CBA-CAA-C2A	-4.47	105.19	112.55
4	A	611	HEC	CBB-CAB-C3B	-4.45	117.08	127.49
4	A	610	HEC	CBB-CAB-C3B	-4.18	117.70	127.49
4	A	611	HEC	CBD-CAD-C3D	-4.15	105.56	112.54
4	A	613	HEC	CBC-CAC-C3C	-4.11	117.88	127.49
4	A	612	HEC	CAA-CBA-CGA	-4.10	102.79	113.83
4	A	609	HEC	CBB-CAB-C3B	-4.07	117.97	127.49
4	A	615	HEC	CBD-CAD-C3D	-3.99	105.83	112.54
4	A	609	HEC	CBA-CAA-C2A	-3.95	106.04	112.55
4	A	615	HEC	CBA-CAA-C2A	-3.93	106.07	112.55
4	A	614	HEC	C1D-C2D-C3D	-3.78	104.37	107.00
4	A	614	HEC	CBC-CAC-C3C	-3.75	118.71	127.49
4	A	612	HEC	C3B-C4B-NB	-3.70	103.96	110.94
4	A	615	HEC	CBC-CAC-C3C	-3.69	118.86	127.49
4	A	612	HEC	CMC-C2C-C1C	-3.63	123.14	128.46
4	A	612	HEC	O2A-CGA-O1A	-3.60	114.06	123.33
4	A	610	HEC	CBC-CAC-C3C	-3.59	119.09	127.49
4	A	614	HEC	CBB-CAB-C3B	-3.46	119.39	127.49
4	A	612	HEC	O1A-CGA-CBA	-3.41	112.27	123.09
4	A	616	HEC	CMB-C2B-C1B	-3.41	123.46	128.46
4	A	612	HEC	C3C-C4C-NC	-3.40	104.52	110.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	A	613	HEC	CBA-CAA-C2A	-3.17	107.33	112.55
4	A	611	HEC	CMB-C2B-C1B	-3.04	124.00	128.46
4	A	616	HEC	CBC-CAC-C3C	-3.03	120.41	127.49
4	A	611	HEC	CBC-CAC-C3C	-3.00	120.47	127.49
4	A	611	HEC	CMC-C2C-C1C	-2.97	124.10	128.46
4	A	616	HEC	CMB-C2B-C3B	2.97	129.31	125.82
4	A	609	HEC	CBC-CAC-C3C	-2.97	120.55	127.49
4	A	612	HEC	CMC-C2C-C3C	2.97	129.31	125.82
4	A	613	HEC	C1D-C2D-C3D	-2.92	104.96	107.00
4	A	614	HEC	CMB-C2B-C1B	-2.91	124.19	128.46
4	A	609	HEC	C1D-C2D-C3D	-2.87	105.00	107.00
4	A	613	HEC	CMB-C2B-C1B	-2.86	124.27	128.46
4	A	612	HEC	CMB-C2B-C1B	-2.85	124.28	128.46
4	A	609	HEC	CMC-C2C-C1C	-2.85	124.29	128.46
4	A	614	HEC	CMC-C2C-C1C	-2.84	124.29	128.46
4	A	612	HEC	C1C-CHC-C4B	-2.84	115.36	123.67
4	A	614	HEC	CAD-CBD-CGD	-2.80	106.30	113.83
4	A	611	HEC	C1D-C2D-C3D	-2.78	105.06	107.00
4	A	615	HEC	CMC-C2C-C1C	-2.75	124.43	128.46
4	A	615	HEC	CMB-C2B-C1B	-2.68	124.53	128.46
4	A	615	HEC	C1D-C2D-C3D	-2.57	105.21	107.00
4	A	616	HEC	CMC-C2C-C1C	-2.53	124.75	128.46
4	A	612	HEC	CBA-CAA-C2A	-2.48	108.47	112.55
4	A	616	HEC	C1D-C2D-C3D	-2.45	105.29	107.00
4	A	613	HEC	CMC-C2C-C1C	-2.39	124.95	128.46
4	A	615	HEC	C2B-C3B-C4B	2.25	108.78	106.35
3	A	606	HG1	CAN-CAT-NAZ	-2.24	114.70	118.09
4	A	611	HEC	O2A-CGA-CBA	2.18	120.90	114.00
4	A	611	HEC	O1A-CGA-CBA	-2.18	116.18	123.09
4	A	610	HEC	CMB-C2B-C1B	-2.17	125.28	128.46
4	A	612	HEC	CMA-C3A-C2A	2.17	129.03	124.94
4	A	610	HEC	C1D-C2D-C3D	-2.15	105.50	107.00
4	A	612	HEC	CMB-C2B-C3B	2.15	128.34	125.82
4	A	616	HEC	C2B-C3B-C4B	2.11	108.63	106.35
4	A	610	HEC	C2B-C3B-C4B	2.09	108.61	106.35
4	A	610	HEC	CMC-C2C-C1C	-2.07	125.42	128.46

There are no chirality outliers.

All (27) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	607	HG1	CAQ-CAL-CAN-CAT

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Mol	Chain	Res	Type	Atoms
3	A	606	HG1	CAW-CAS-NAZ-CAR
3	A	606	HG1	CAW-CAS-NAZ-CAT
3	A	608	HG1	CAN-CAL-CAQ-CAU
3	A	606	HG1	OAF-CAX-CAY-CAW
3	A	607	HG1	CAN-CAL-CAQ-CAU
4	A	616	HEC	C3D-CAD-CBD-CGD
3	A	607	HG1	OAB-CAH-CAR-NAZ
4	A	615	HEC	C3D-CAD-CBD-CGD
3	A	606	HG1	OAF-CAX-CAY-OAG
4	A	612	HEC	CAA-CBA-CGA-O1A
4	A	612	HEC	CAA-CBA-CGA-O2A
4	A	609	HEC	CAD-CBD-CGD-O1D
3	A	606	HG1	CAQ-CAL-CAN-CAT
4	A	609	HEC	CAD-CBD-CGD-O2D
4	A	611	HEC	CAD-CBD-CGD-O2D
4	A	614	HEC	CAA-CBA-CGA-O2A
4	A	614	HEC	CAD-CBD-CGD-O2D
4	A	614	HEC	CAA-CBA-CGA-O1A
4	A	614	HEC	CAD-CBD-CGD-O1D
4	A	616	HEC	CAD-CBD-CGD-O2D
4	A	611	HEC	CAD-CBD-CGD-O1D
4	A	616	HEC	CAD-CBD-CGD-O1D
4	A	609	HEC	C3D-CAD-CBD-CGD
4	A	612	HEC	CAD-CBD-CGD-O2D
3	A	606	HG1	CAS-CAW-CAY-CAX
4	A	611	HEC	CAA-CBA-CGA-O2A

There are no ring outliers.

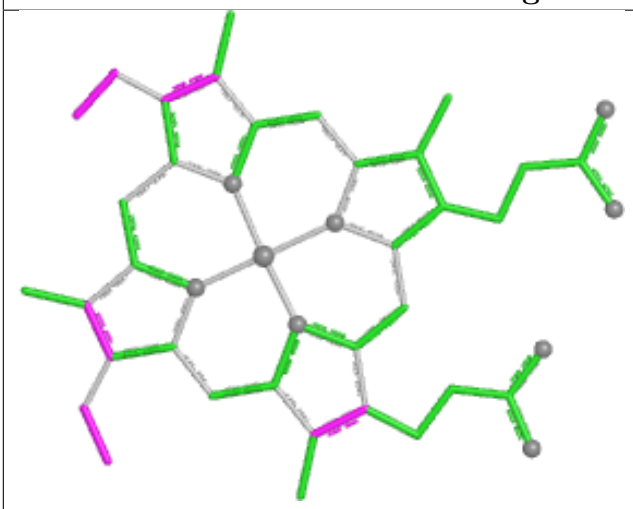
9 monomers are involved in 23 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	611	HEC	1	0
4	A	613	HEC	2	0
4	A	615	HEC	3	0
4	A	609	HEC	3	0
4	A	614	HEC	4	0
4	A	610	HEC	1	0
4	A	612	HEC	4	2
4	A	616	HEC	3	0
3	A	606	HG1	1	0

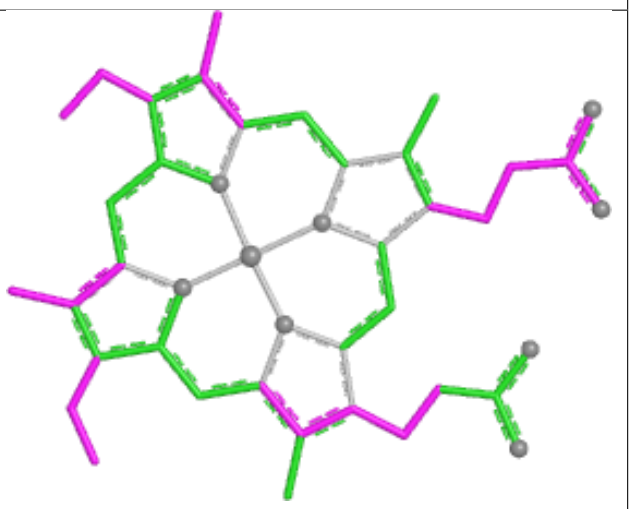
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths,

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 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.

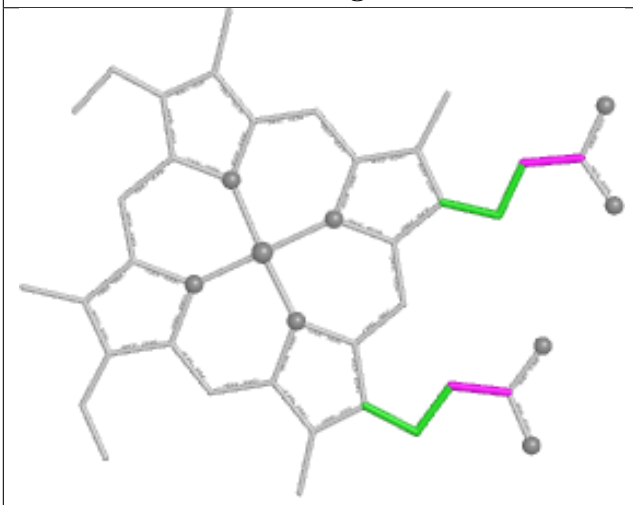
Ligand HEC A 611



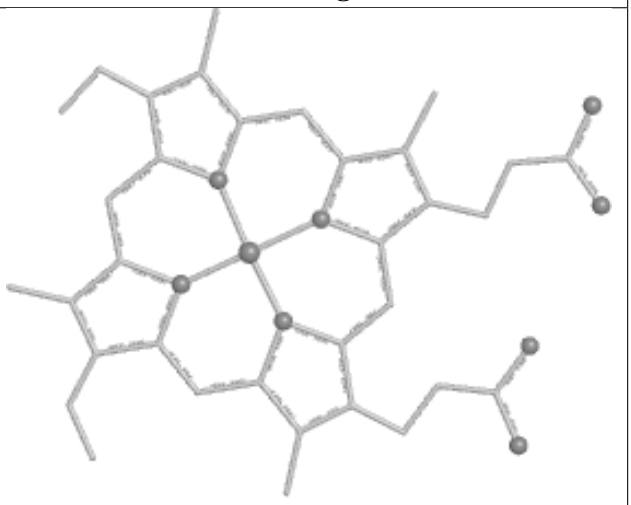
Bond lengths



Bond angles

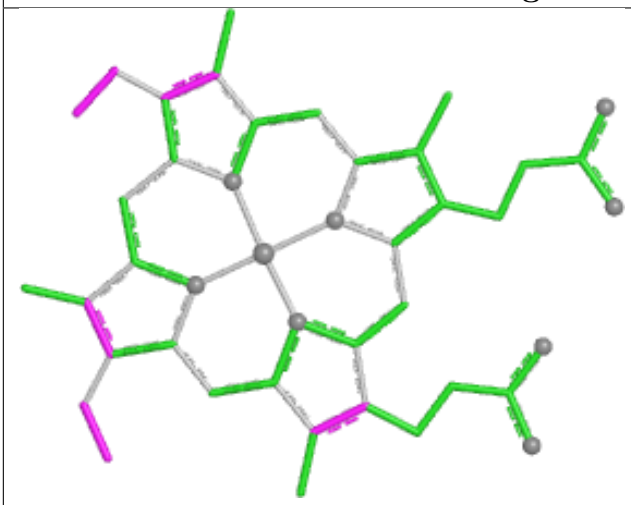


Torsions

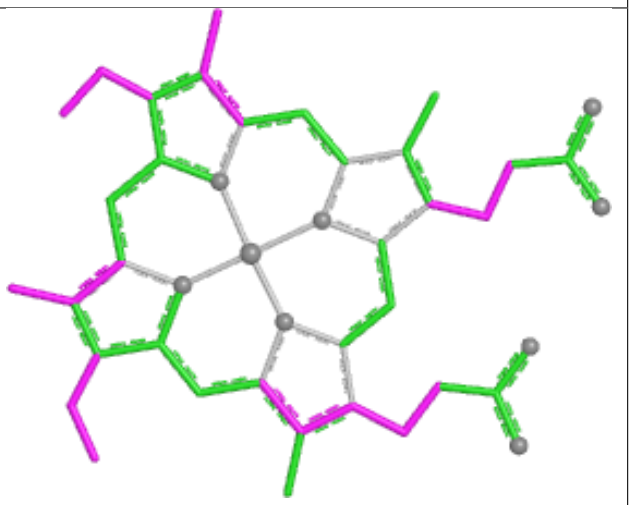


Rings

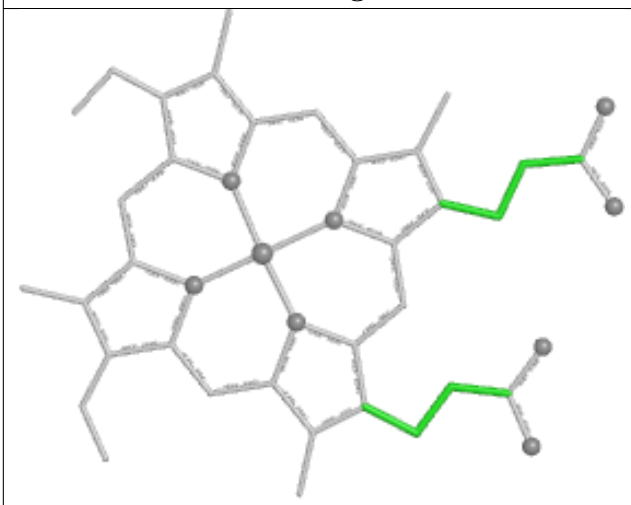
Ligand HEC A 613



Bond lengths



Bond angles



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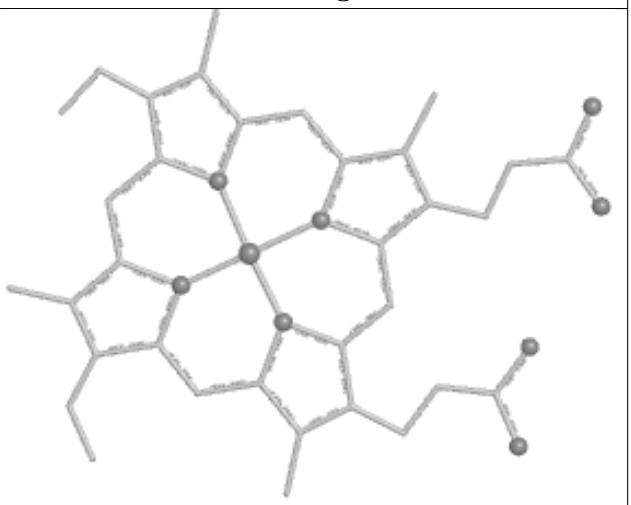

Rings

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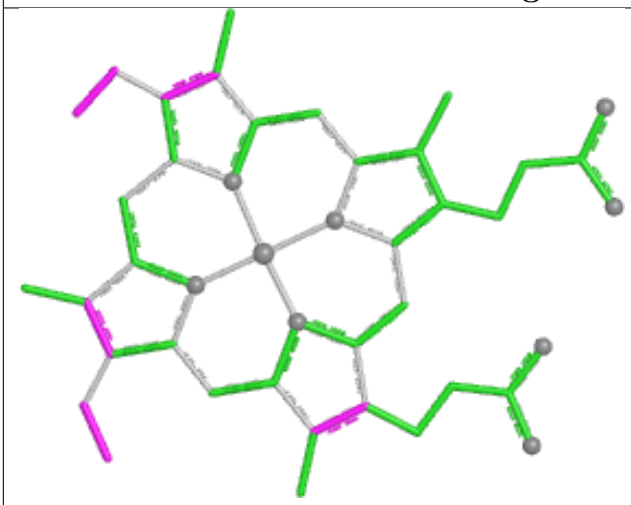
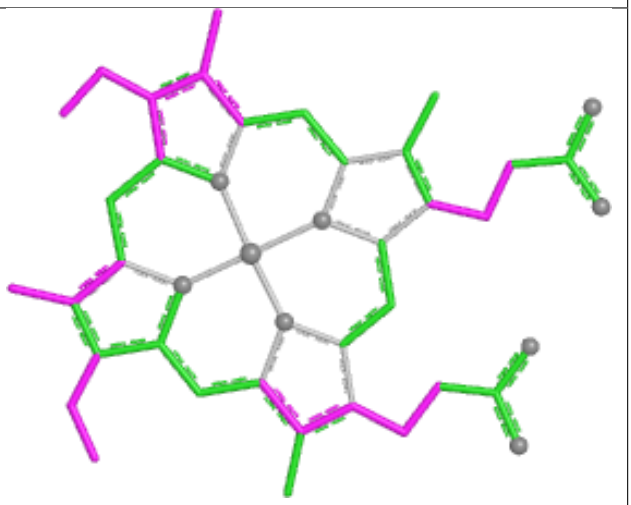
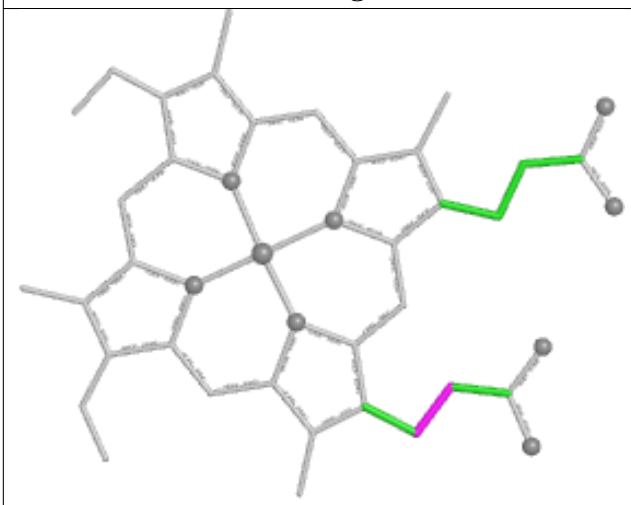


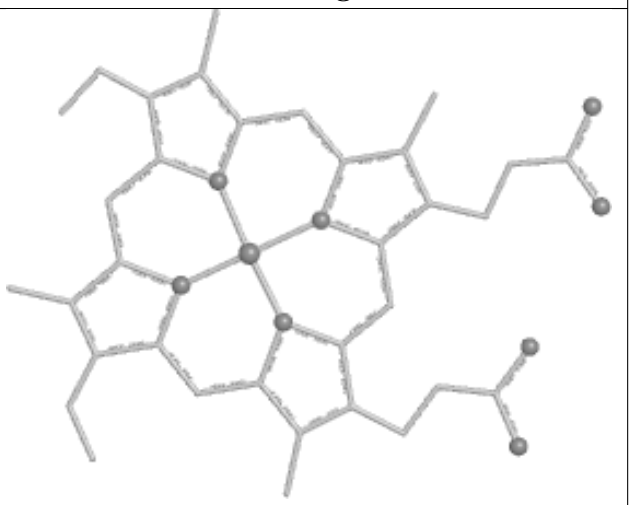
Bond lengths



Bond angles

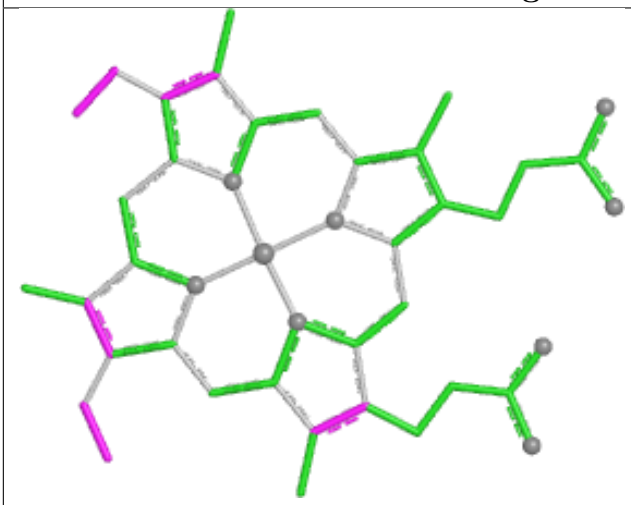


Torsions



Rings

Ligand HEC A 609



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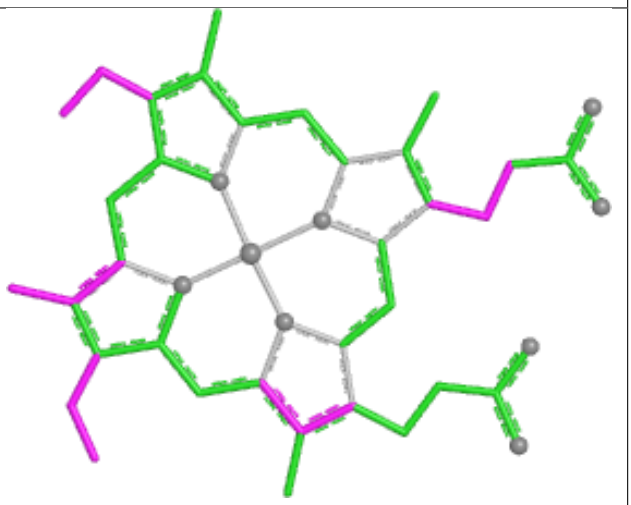

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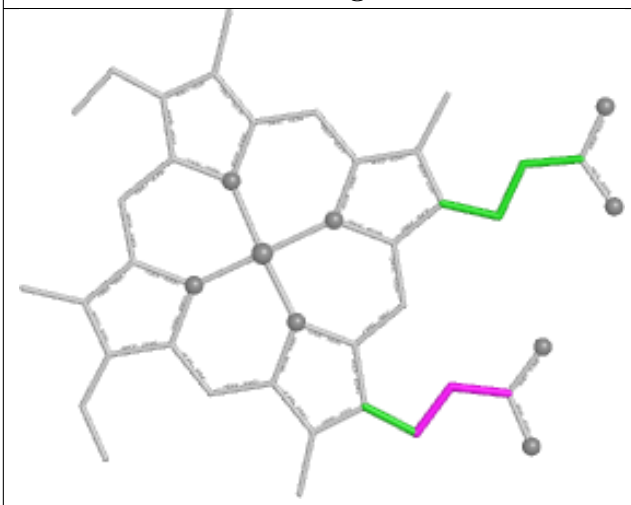

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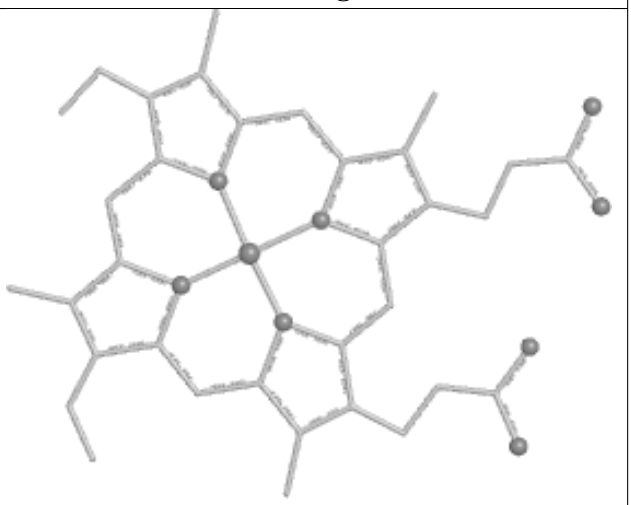
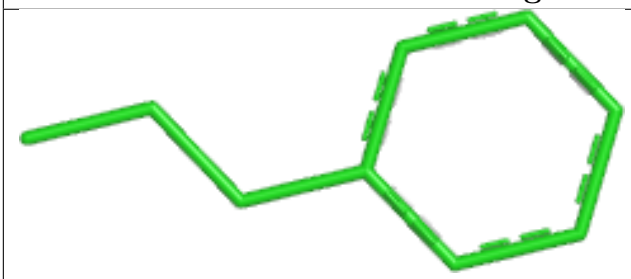


Rings

Ligand HG1 A 608



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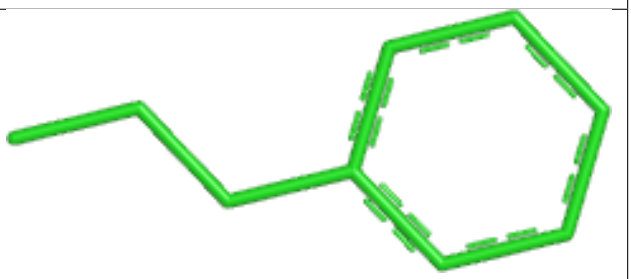
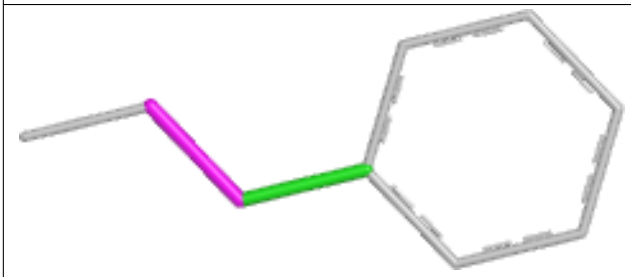


Bond angles



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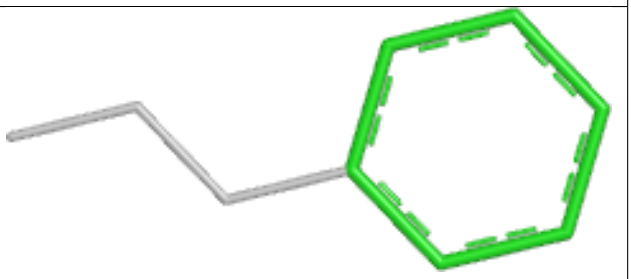

Rings

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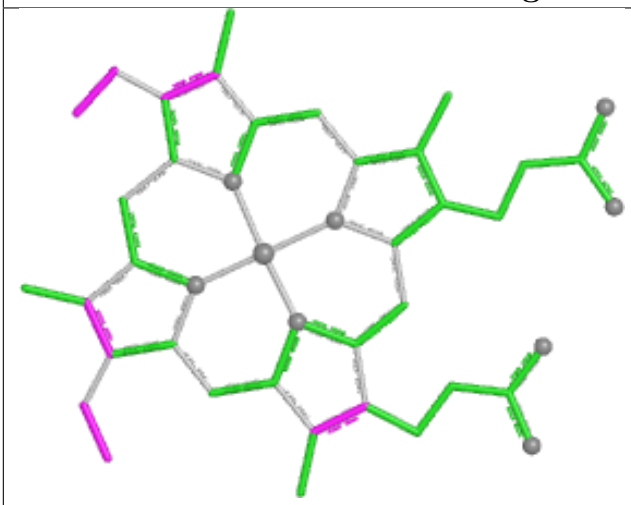

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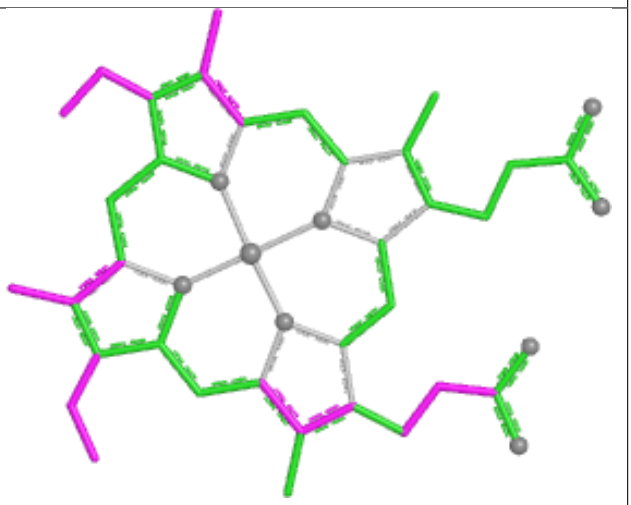
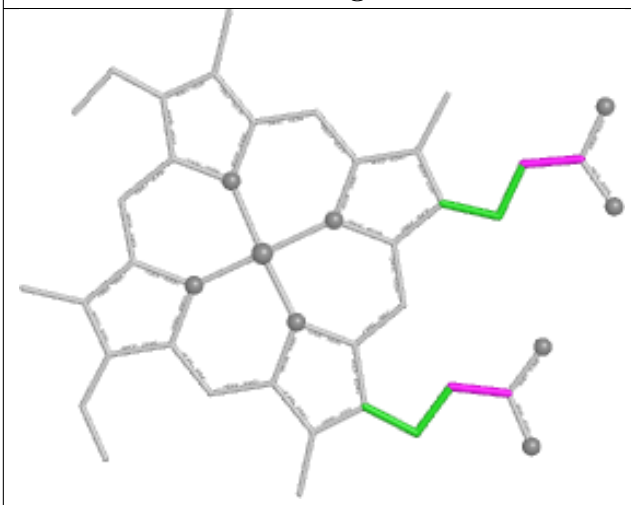


Bond angles



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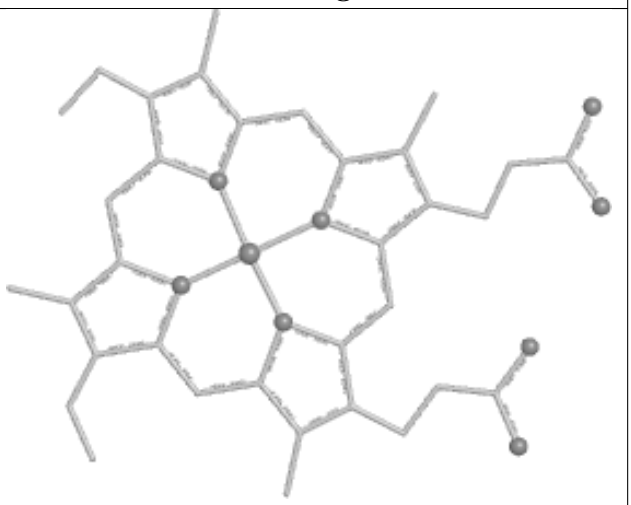
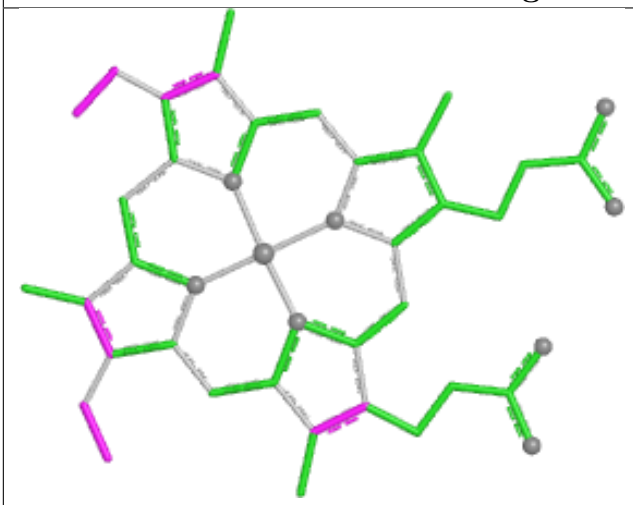


Rings

Ligand HEC A 610



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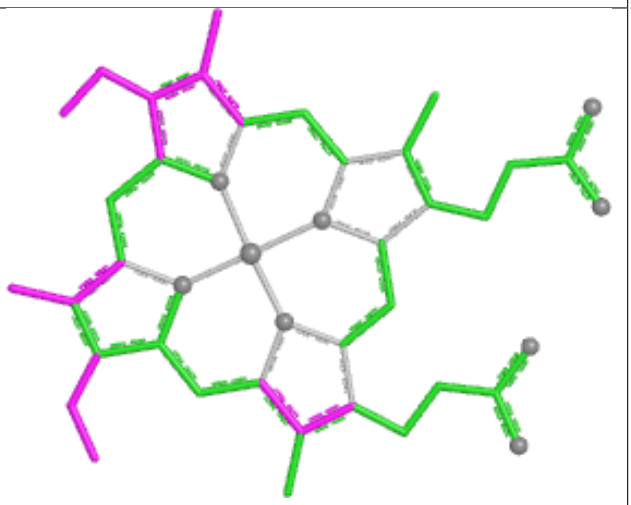
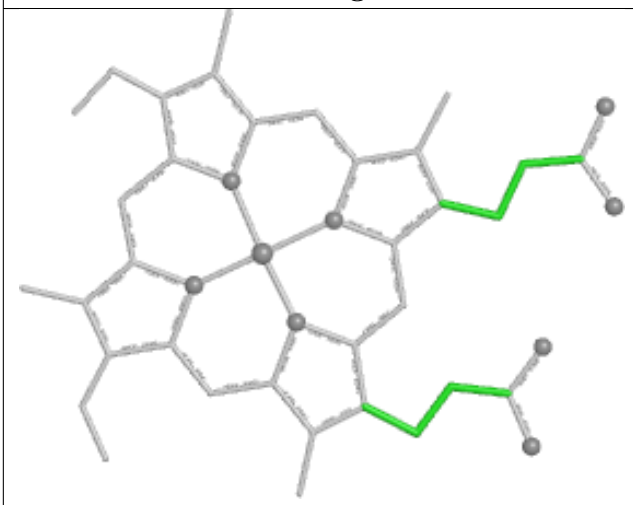


Bond angles



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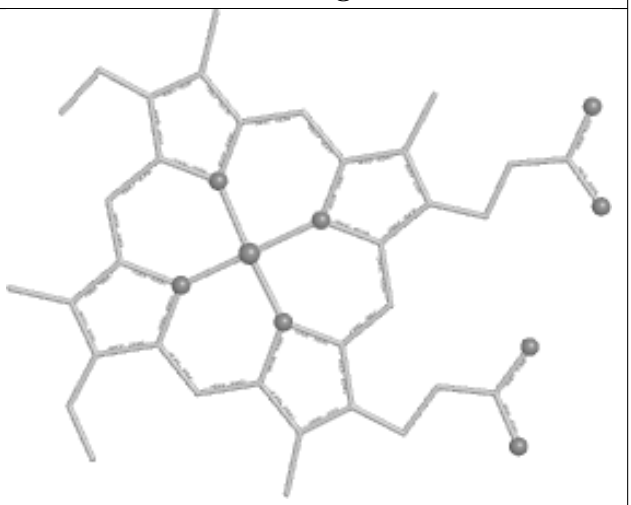
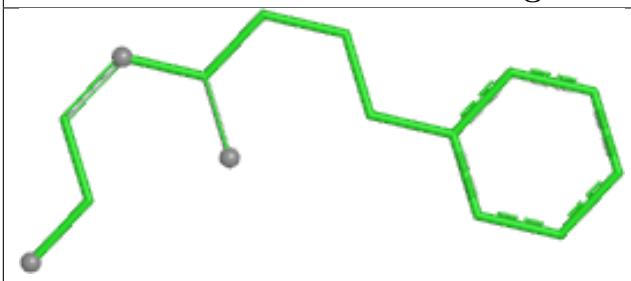
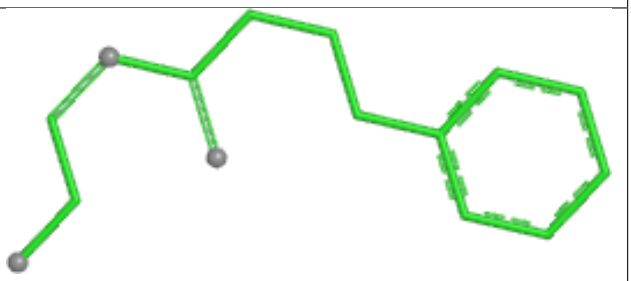


Rings

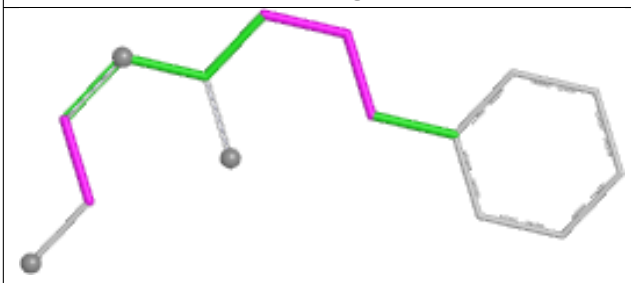
Ligand HG1 A 607



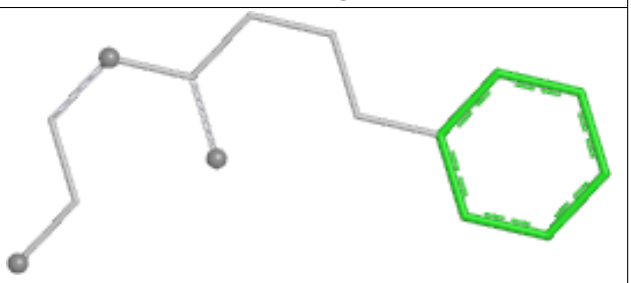
Bond lengths



Bond angles

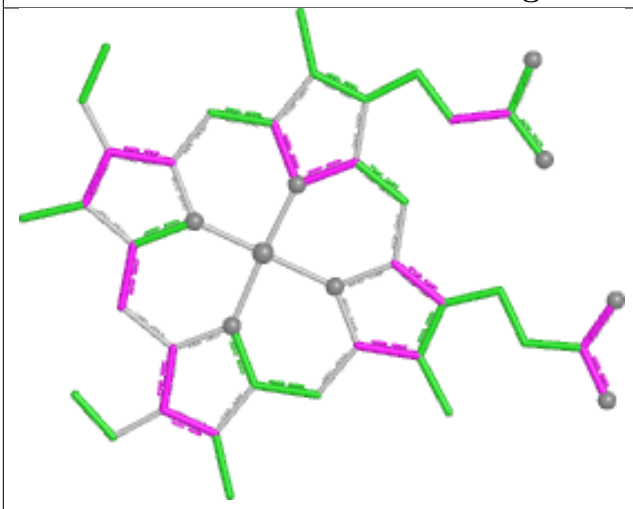


Torsions

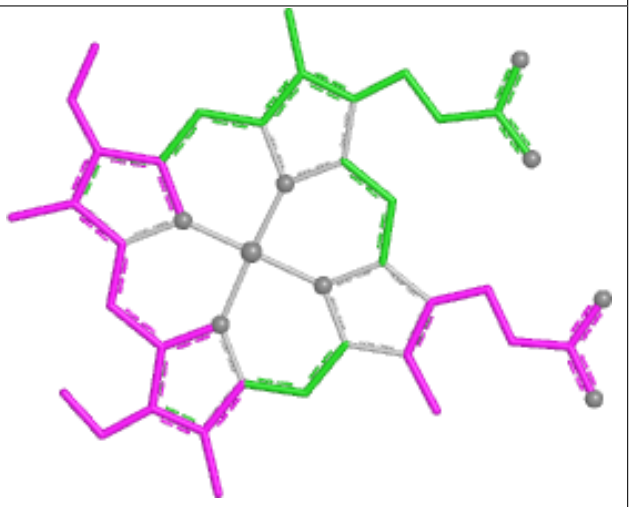


Rings

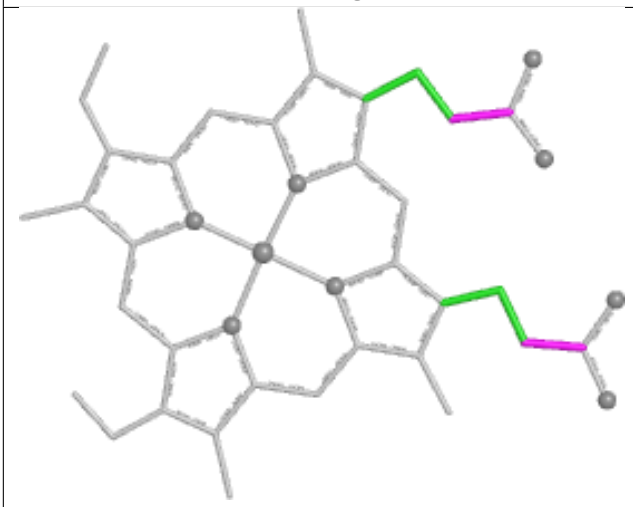
Ligand HEC A 612



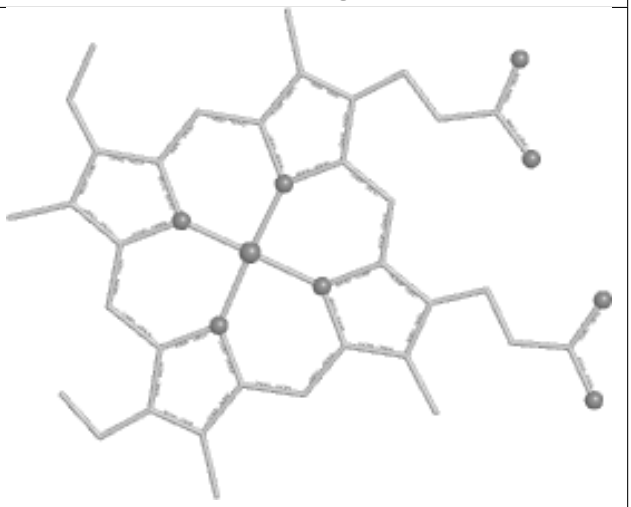
Bond lengths



Bond angles

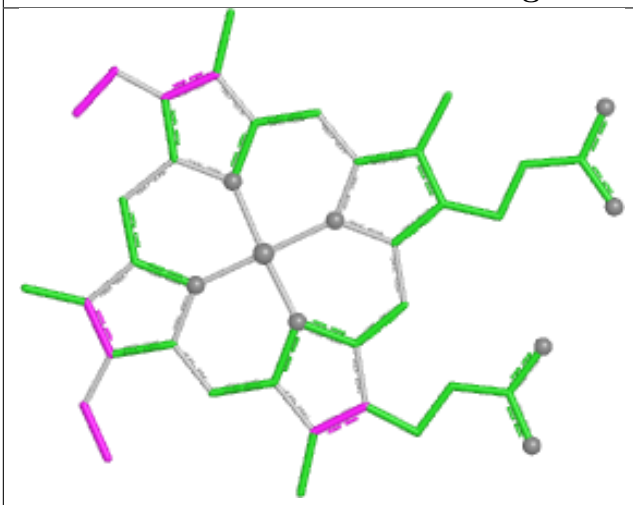


Torsions

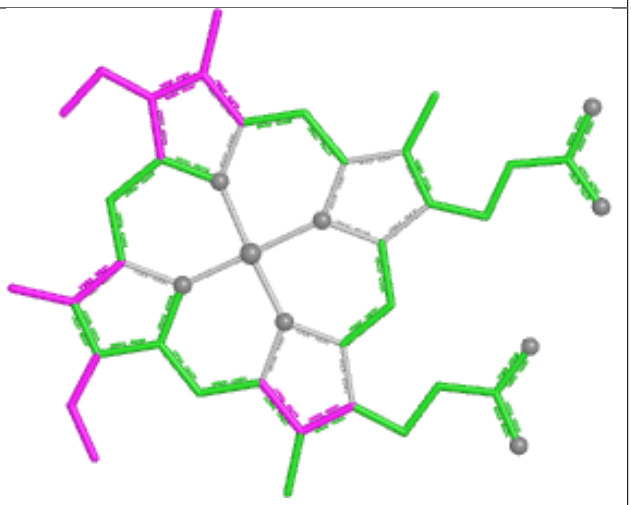


Rings

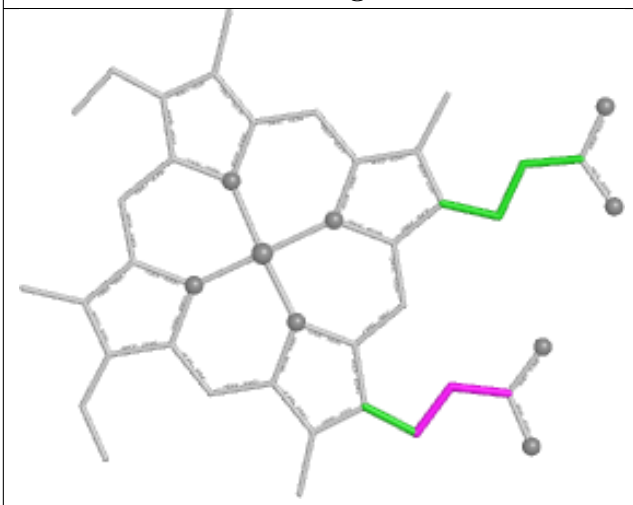
Ligand HEC A 616



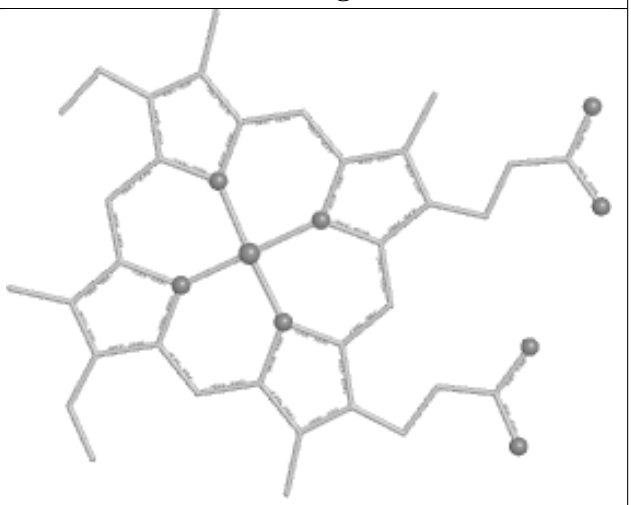
Bond lengths



Bond angles

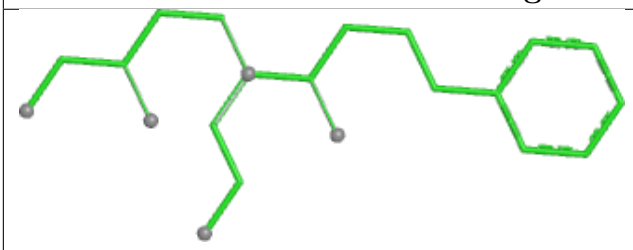


Torsions

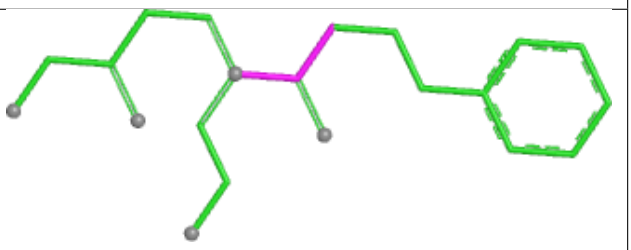


Rings

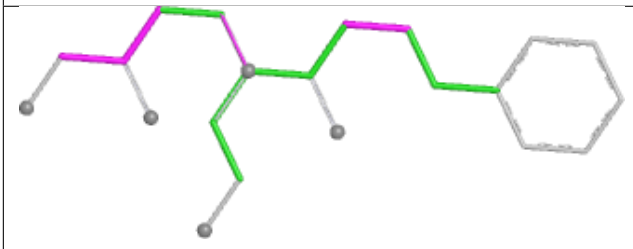
Ligand HG1 A 606



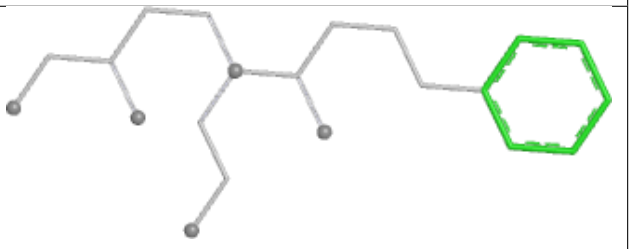
Bond lengths



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)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2	OWAB(Å^2)	Q<0.9
1	A	497/500 (99%)	-0.35	7 (1%) 75 72	9, 14, 24, 33	0

All (7) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	38	GLY	6.3
1	A	534	ILE	3.8
1	A	486	GLU	3.2
1	A	39	PRO	2.9
1	A	140	ASP	2.5
1	A	488	LYS	2.4
1	A	152	LYS	2.2

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	HG1	A	606	21/26	0.67	0.26	47,50,52,53	0

Continued on next page...

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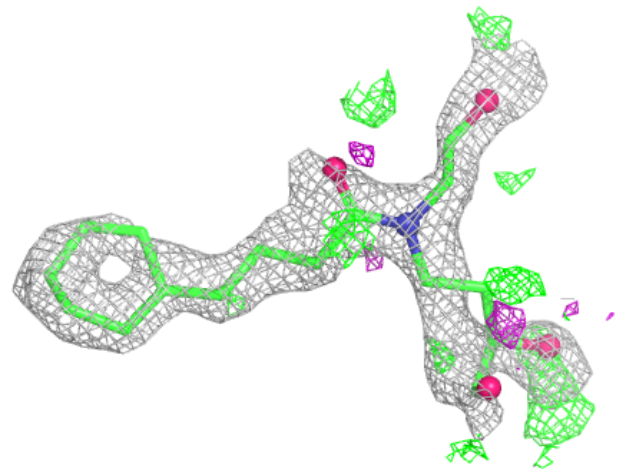
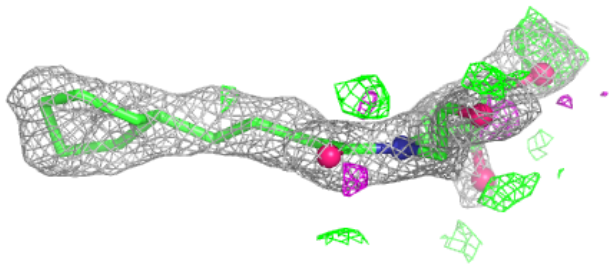
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	HG1	A	607	15/26	0.74	0.19	47,49,53,53	0
3	HG1	A	605	12/26	0.84	0.14	35,35,37,38	0
2	PO4	A	604	5/5	0.86	0.30	31,33,34,34	0
3	HG1	A	608	9/26	0.87	0.13	47,48,48,48	0
2	PO4	A	602	5/5	0.93	0.20	25,27,28,28	0
2	PO4	A	601	5/5	0.97	0.16	24,25,26,27	0
4	HEC	A	612	43/43	0.97	0.10	11,13,17,21	0
4	HEC	A	610	43/43	0.98	0.08	9,11,12,12	0
4	HEC	A	611	43/43	0.98	0.08	8,11,13,15	0
4	HEC	A	609	43/43	0.98	0.07	12,13,16,17	0
4	HEC	A	614	43/43	0.98	0.09	7,9,14,16	0
4	HEC	A	615	43/43	0.98	0.09	7,9,12,14	0
2	PO4	A	603	5/5	0.99	0.08	14,14,15,15	5
4	HEC	A	613	43/43	0.99	0.07	8,10,13,16	0
4	HEC	A	616	43/43	0.99	0.09	8,9,11,11	0

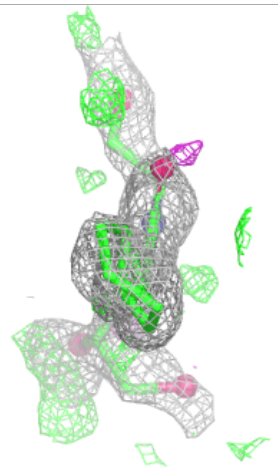
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around HG1 A 606:

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

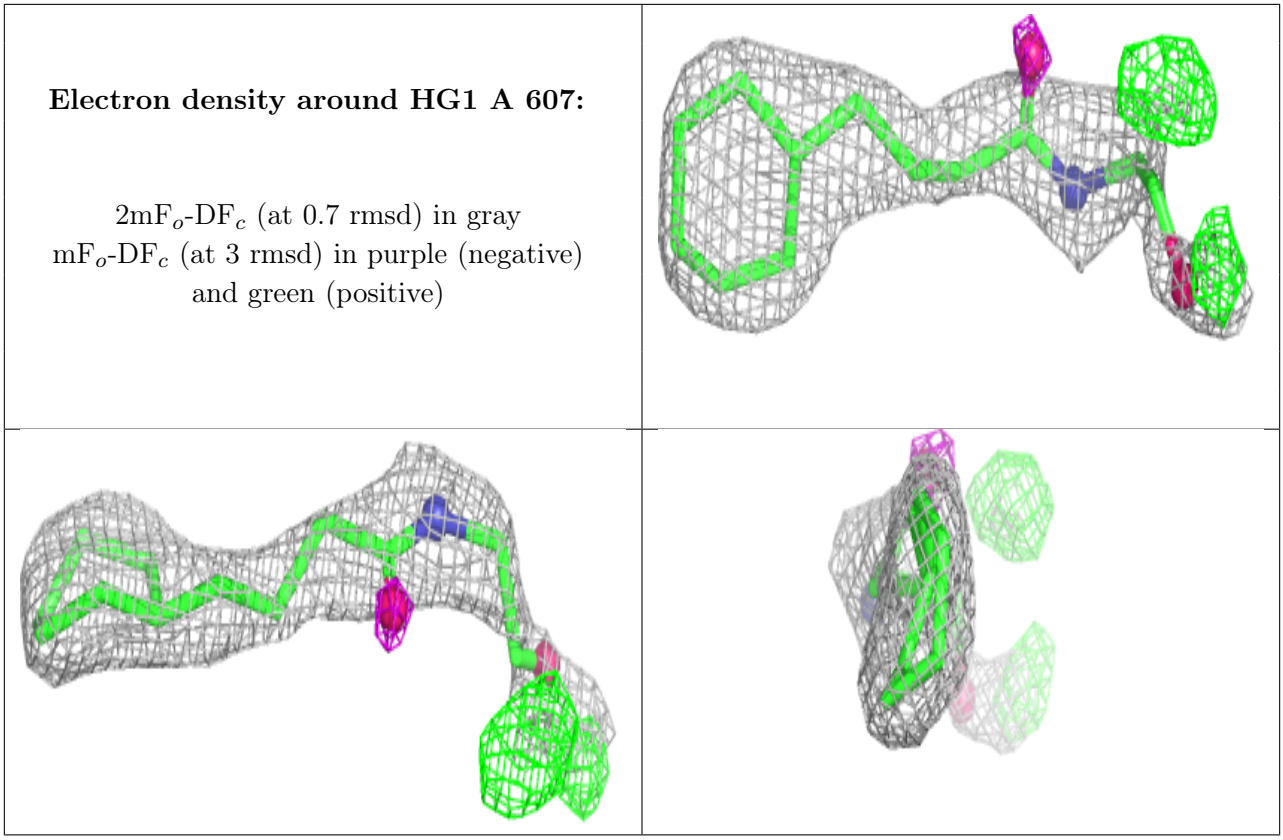






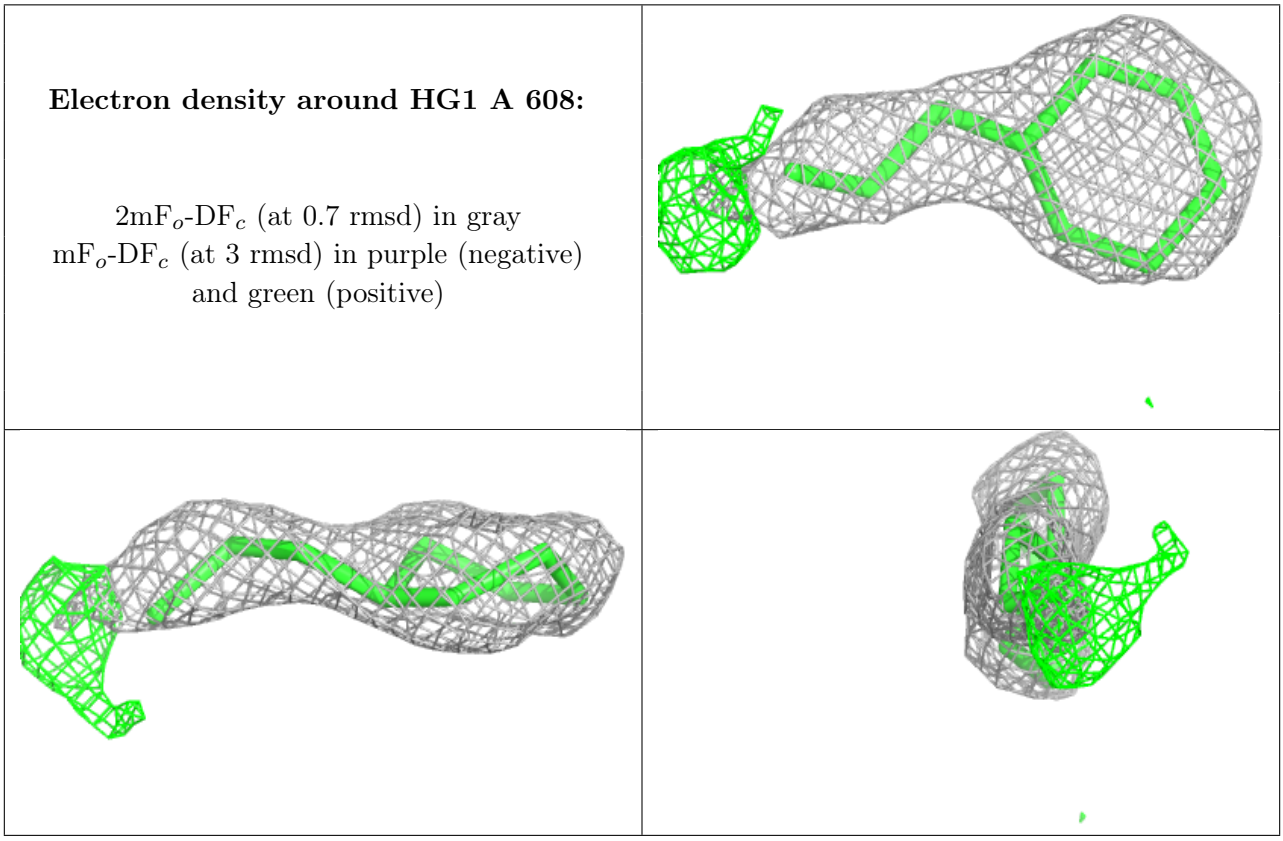
Electron density around HG1 A 607:

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



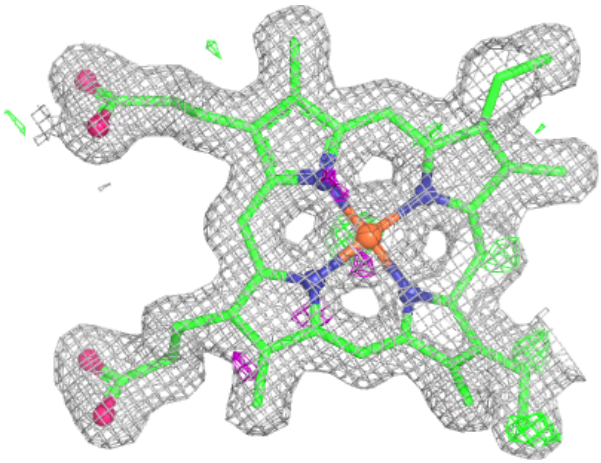
Electron density around HG1 A 608:

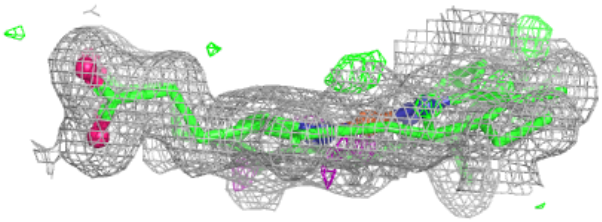
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

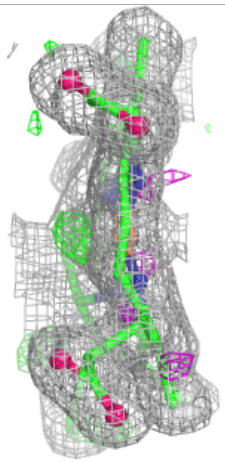


Electron density around HEC A 612:

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

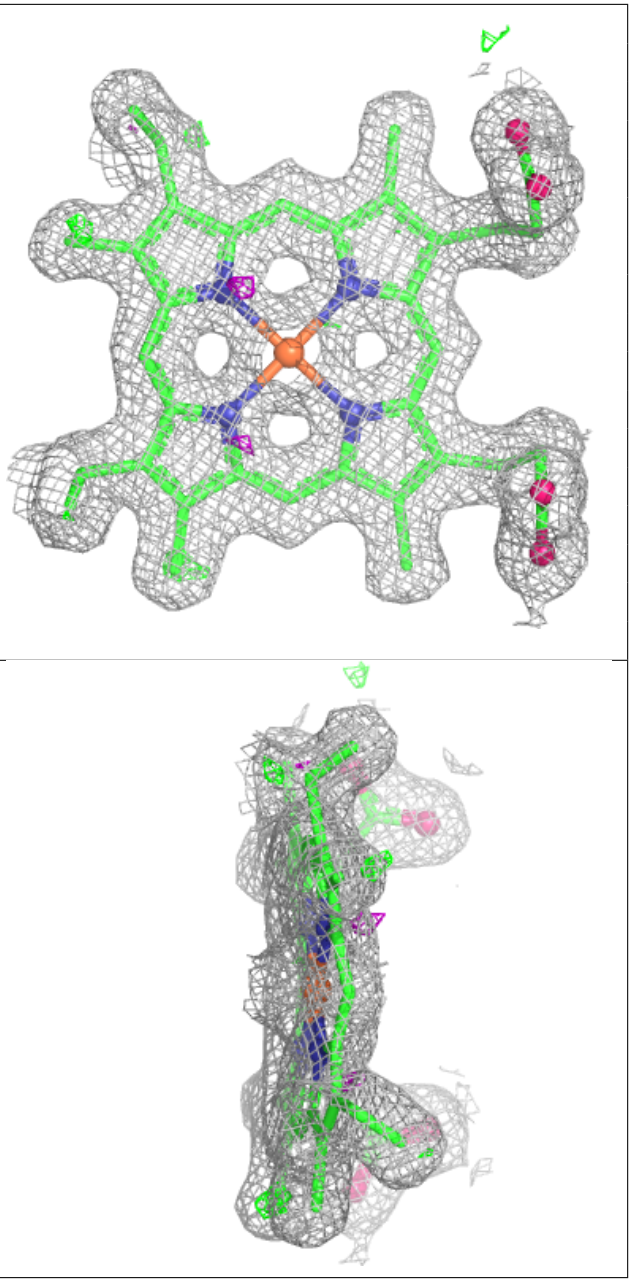






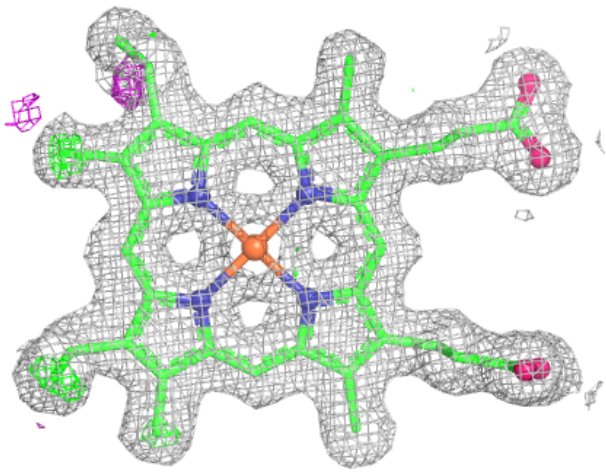
Electron density around HEC A 610:

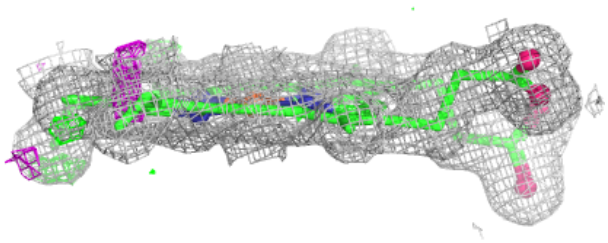
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

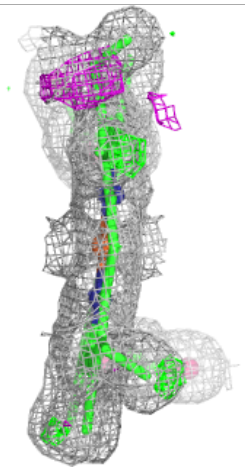


Electron density around HEC A 611:

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

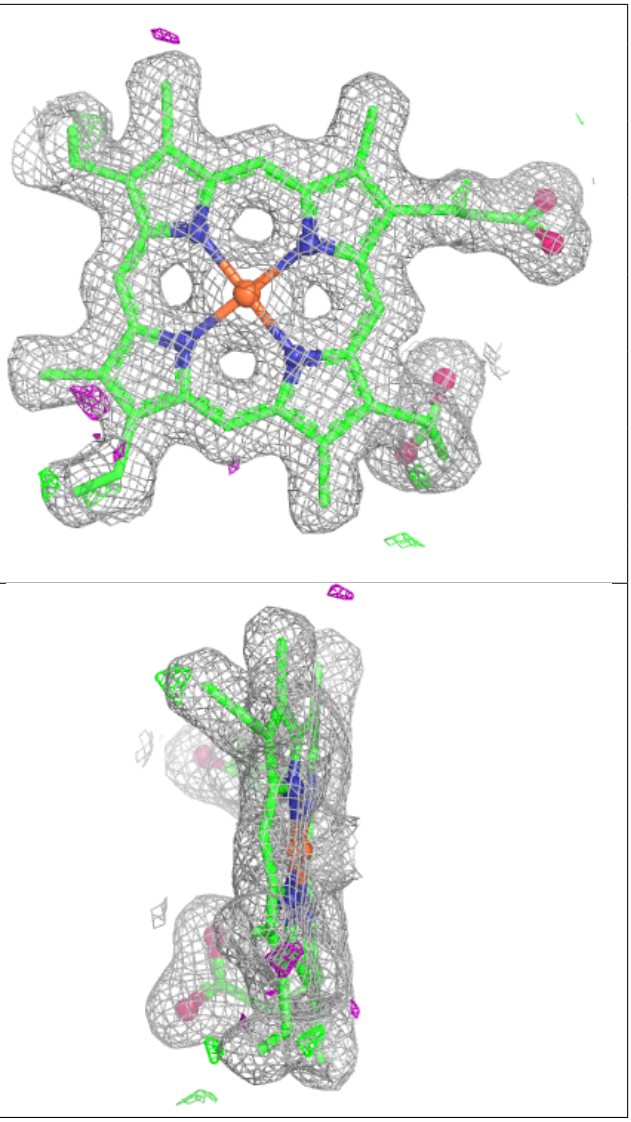






Electron density around HEC A 609:

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



Electron density around HEC A 614:

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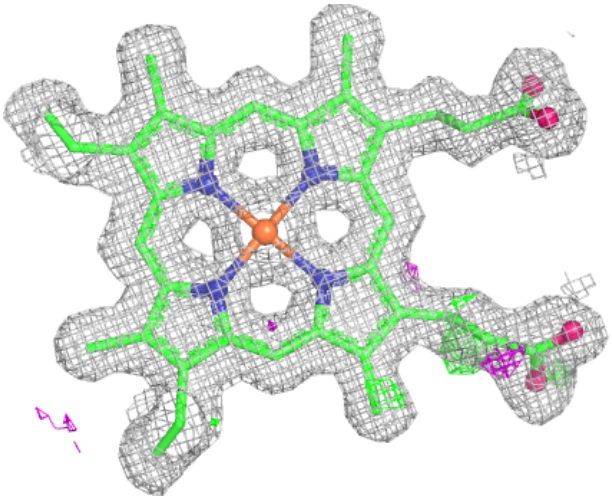
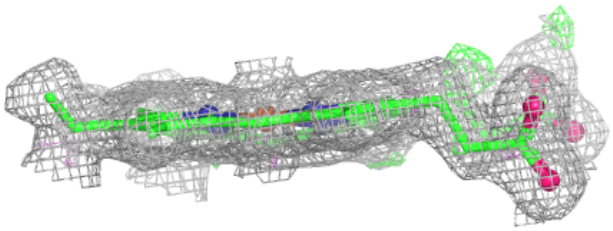
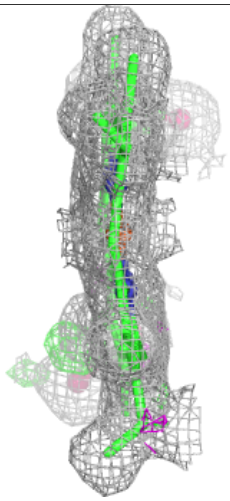
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)







Electron density around HEC A 615:

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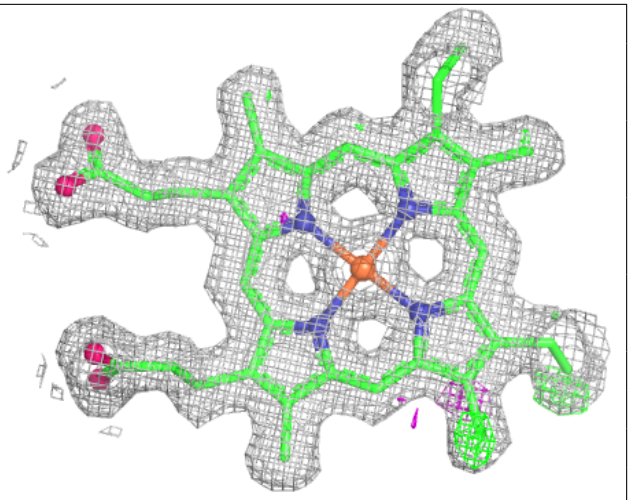
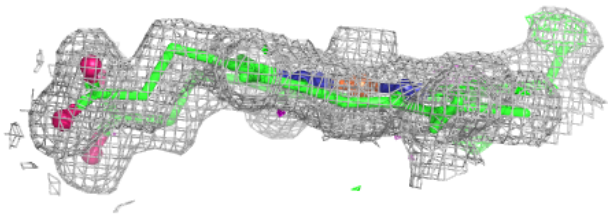
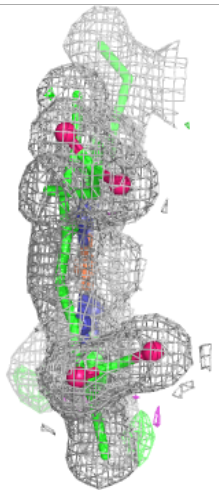
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

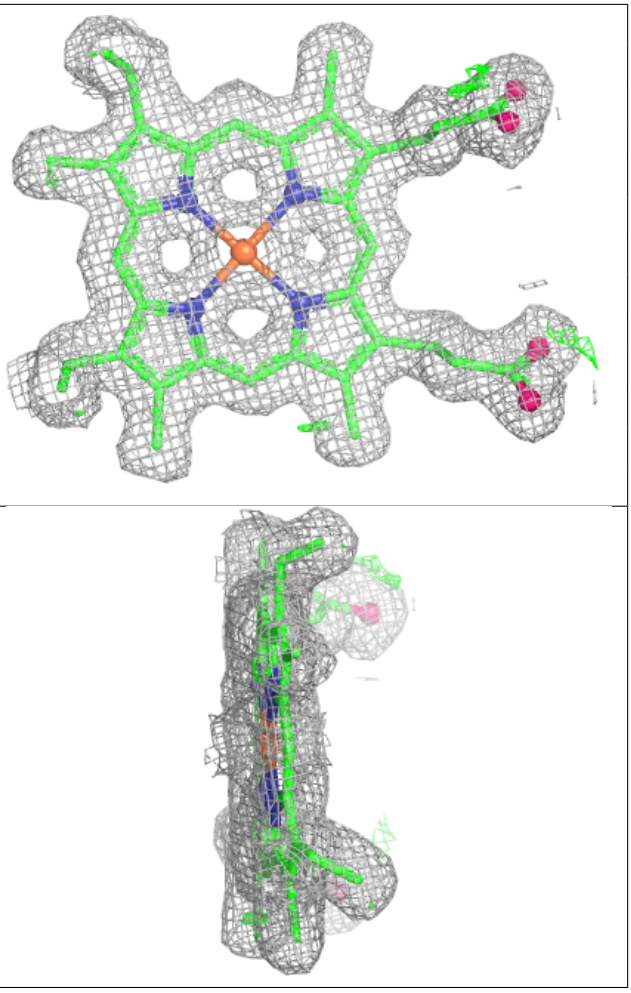






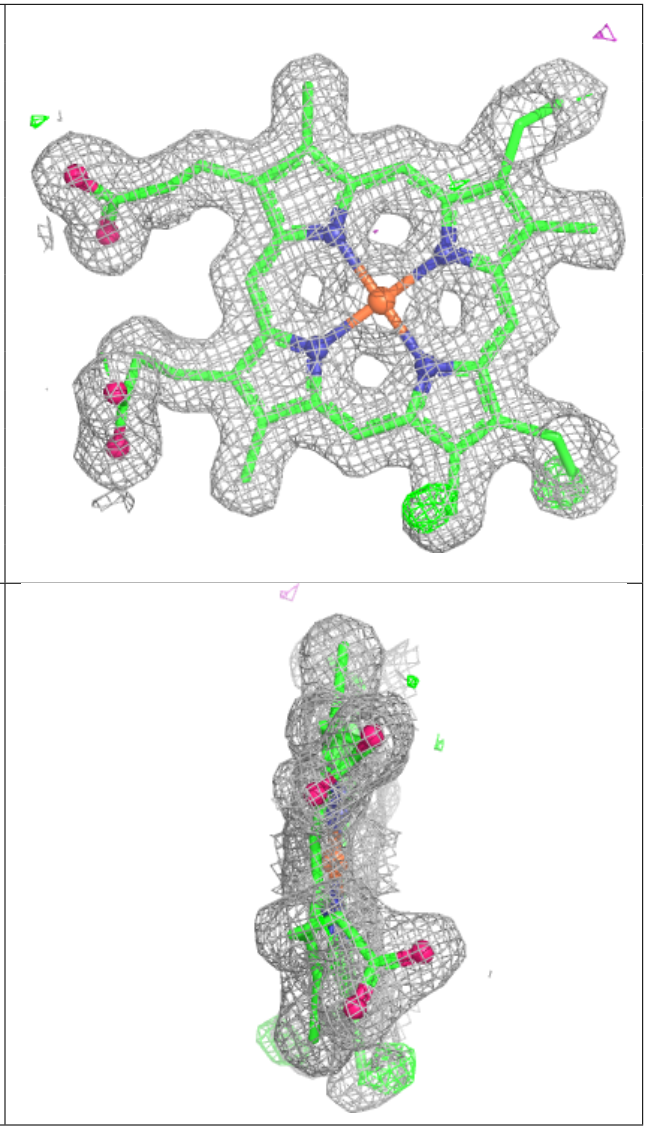
Electron density around HEC A 613:

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



Electron density around HEC A 616:

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



6.5 Other polymers (i)

There are no such residues in this entry.