Full wwPDB X-ray Structure Validation Report (i)

May 14, 2020 – 03:24 pm BST

PDB ID	:	2VXI
Title	:	The binding of heme and zinc in Escherichia coli Bacterioferritin
Authors	:	Willies, S.C.; Isupov, M.N.; Garman, E.F.; Littlechild, J.A.
Deposited on	:	2008-07-04
Resolution	:	1.91 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 following versions of software and data (see references (i)) were used in the production of this report:

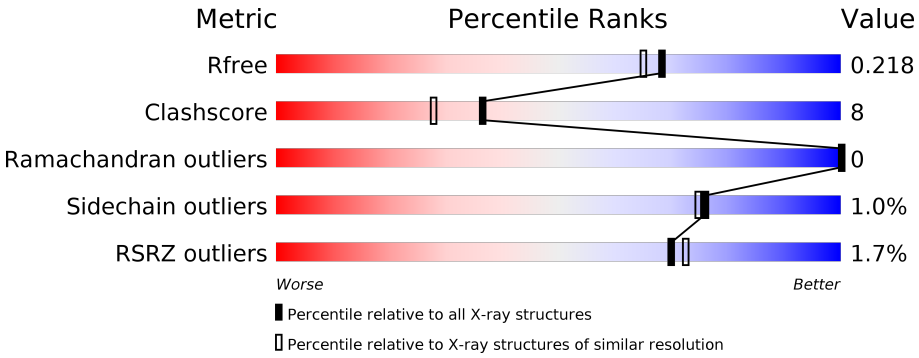
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.11

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 1.91 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	7937 (1.94-1.90)
Clashscore	141614	8644 (1.94-1.90)
Ramachandran outliers	138981	8530 (1.94-1.90)
Sidechain outliers	138945	8530 (1.94-1.90)
RSRZ outliers	127900	7793 (1.94-1.90)

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 on 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	158	3% 86% 13%
1	B	158	3% 89% 10%
1	C	158	% 92% 6%
1	D	158	2% 89% 9%
1	E	158	% 94% 5%
1	F	158	2% 93% 6%

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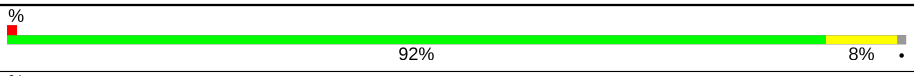
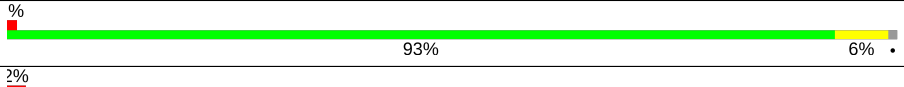
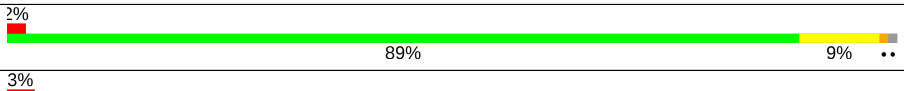
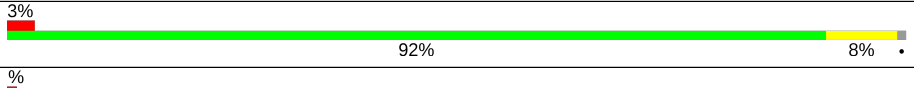
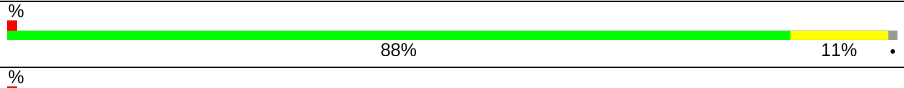
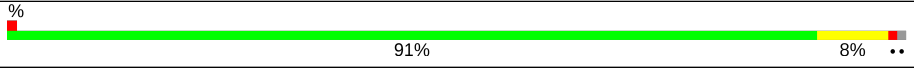
Mol	Chain	Length	Quality of chain
1	G	158	
1	H	158	
1	I	158	
1	J	158	
1	K	158	
1	L	158	

2 Entry composition

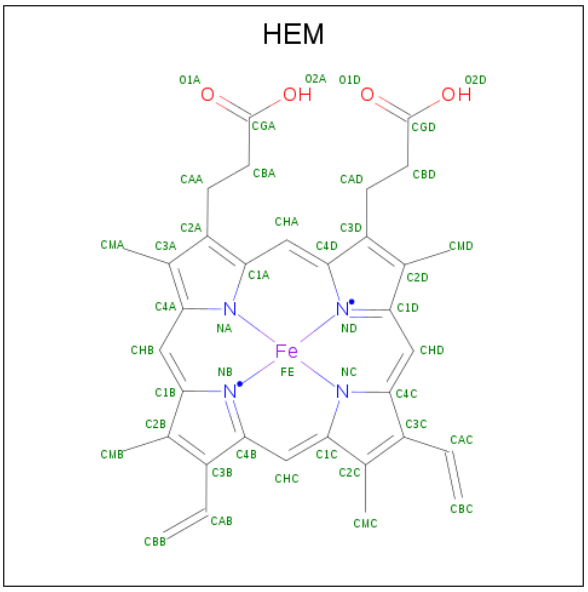
There are 5 unique types of molecules in this entry. The entry contains 18818 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 BACTERIOFERRITIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	157	Total	C	N	O	S	0	10	0
			1346	850	229	260	7			
1	B	157	Total	C	N	O	S	0	5	0
			1325	836	229	253	7			
1	C	157	Total	C	N	O	S	0	5	0
			1325	836	229	253	7			
1	D	157	Total	C	N	O	S	0	5	0
			1323	835	228	253	7			
1	E	157	Total	C	N	O	S	0	6	0
			1328	838	229	254	7			
1	F	157	Total	C	N	O	S	0	6	0
			1326	837	228	254	7			
1	G	157	Total	C	N	O	S	0	5	0
			1325	837	227	254	7			
1	H	157	Total	C	N	O	S	0	5	0
			1323	835	228	253	7			
1	I	157	Total	C	N	O	S	0	4	0
			1318	832	227	252	7			
1	J	157	Total	C	N	O	S	0	5	0
			1326	837	228	254	7			
1	K	157	Total	C	N	O	S	0	4	0
			1318	832	225	254	7			
1	L	157	Total	C	N	O	S	0	4	0
			1318	832	227	252	7			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).

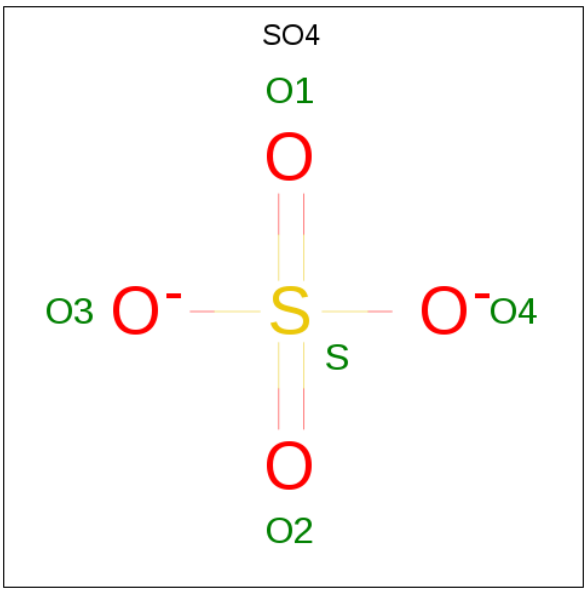


Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	B	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	C	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	D	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	E	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	F	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	G	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	H	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	I	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	J	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	K	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		
2	L	1	Total	C	Fe	N	O	0	1
			45	36	1	4	4		

- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	G	2	Total Zn 2 2	0	0
3	J	2	Total Zn 2 2	0	0
3	D	2	Total Zn 2 2	0	0
3	K	2	Total Zn 2 2	0	0
3	E	2	Total Zn 2 2	0	0
3	H	2	Total Zn 2 2	0	0
3	B	2	Total Zn 2 2	0	0
3	I	2	Total Zn 2 2	0	0
3	C	2	Total Zn 2 2	0	0
3	A	2	Total Zn 2 2	0	0
3	L	2	Total Zn 2 2	0	0
3	F	2	Total Zn 2 2	0	0

- Molecule 4 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	A	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	B	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0
4	C	1	Total O S 5 4 1	0	0
4	D	1	Total O S 5 4 1	0	0
4	D	1	Total O S 5 4 1	0	0
4	E	1	Total O S 5 4 1	0	0
4	E	1	Total O S 5 4 1	0	0
4	E	1	Total O S 5 4 1	0	0
4	F	1	Total O S 5 4 1	0	0
4	G	1	Total O S 5 4 1	0	0
4	G	1	Total O S 5 4 1	0	0
4	H	1	Total O S 5 4 1	0	0
4	H	1	Total O S 5 4 1	0	0
4	H	1	Total O S 5 4 1	0	0
4	I	1	Total O S 5 4 1	0	0

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Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	I	1	Total O S 5 4 1	0	0
4	J	1	Total O S 5 4 1	0	0
4	J	1	Total O S 5 4 1	0	0
4	K	1	Total O S 5 4 1	0	0
4	K	1	Total O S 5 4 1	0	0
4	L	1	Total O S 5 4 1	0	0
4	L	1	Total O S 5 4 1	0	0

- Molecule 5 is water.

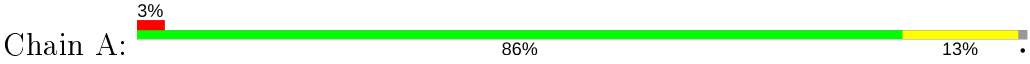
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	190	Total O 190 190	0	0
5	B	184	Total O 184 184	0	0
5	C	189	Total O 189 189	0	0
5	D	181	Total O 181 181	0	0
5	E	191	Total O 191 191	0	0
5	F	179	Total O 179 179	0	0
5	G	195	Total O 195 195	0	0
5	H	188	Total O 188 188	0	0
5	I	194	Total O 194 194	0	0
5	J	179	Total O 179 179	0	0
5	K	173	Total O 173 173	0	0
5	L	165	Total O 165 165	0	0

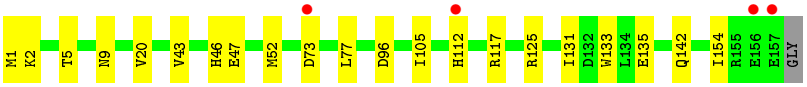
3 Residue-property plots (i)

These plots are drawn for all protein, RNA and DNA 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: BACTERIOFERRITIN

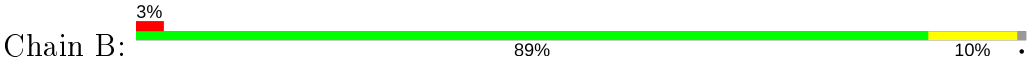
Chain A:

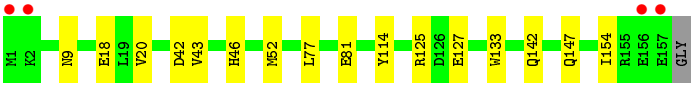




- Molecule 1: BACTERIOFERRITIN

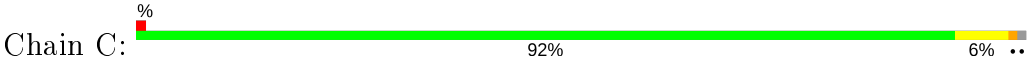
Chain B:

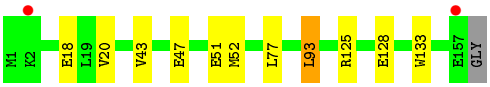




- Molecule 1: BACTERIOFERRITIN

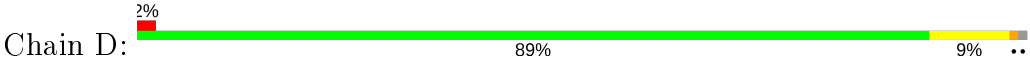
Chain C:

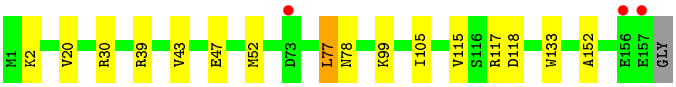




- Molecule 1: BACTERIOFERRITIN

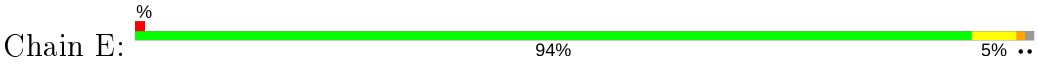
Chain D:

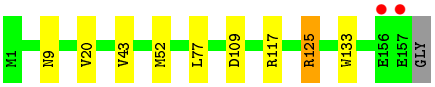




- Molecule 1: BACTERIOFERRITIN

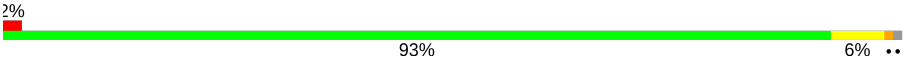
Chain E:





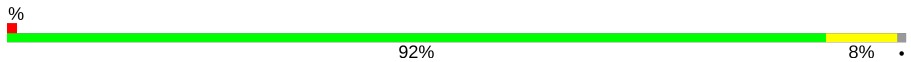
- Molecule 1: BACTERIOFERRITIN

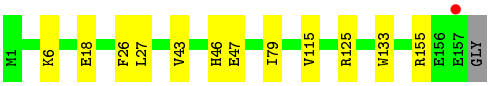
Chain F:



• Molecule 1: BACTERIOFERRITIN

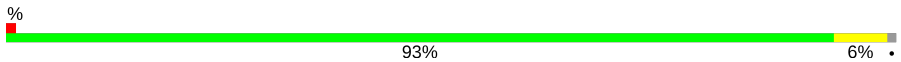
Chain G:

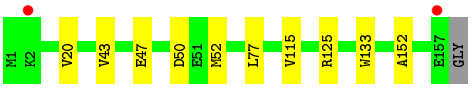




• Molecule 1: BACTERIOFERRITIN

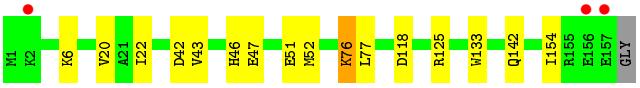
Chain H:





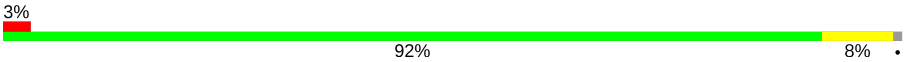
• Molecule 1: BACTERIOFERRITIN

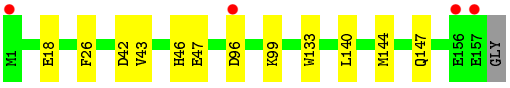
Chain I:



• Molecule 1: BACTERIOFERRITIN

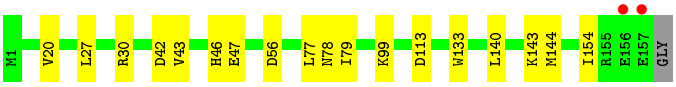
Chain J:





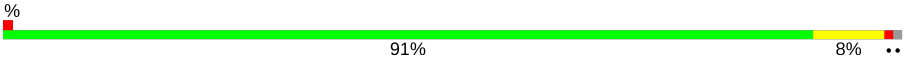
• Molecule 1: BACTERIOFERRITIN

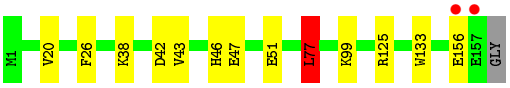
Chain K:



• Molecule 1: BACTERIOFERRITIN

Chain L:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 42 21 2	Depositor
Cell constants a, b, c, α, β, γ	208.10Å 208.10Å 142.77Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	147.44 – 1.91 23.27 – 1.91	Depositor EDS
% Data completeness (in resolution range)	98.7 (147.44-1.91) 97.9 (23.27-1.91)	Depositor EDS
R_{merge}	0.10	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.36 (at 1.92Å)	Xtriage
Refinement program	REFMAC 5.4.0057	Depositor
R, R_{free}	0.179 , 0.216 0.183 , 0.218	Depositor DCC
R_{free} test set	9100 reflections (2.01%)	wwPDB-VP
Wilson B-factor (Å^2)	22.6	Xtriage
Anisotropy	0.018	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.38 , 59.7	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.98	EDS
Total number of atoms	18818	wwPDB-VP
Average B, all atoms (Å^2)	24.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The analyses of the Patterson function reveals a significant off-origin peak that is 49.24 % of the origin peak, indicating pseudo-translational symmetry. The chance of finding a peak of this or larger height randomly in a structure without pseudo-translational symmetry is equal to 7.6691e-05. The detected translational NCS is most likely also responsible for the elevated intensity ratio.*

[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

5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: HEM, ZN, SO4

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	0.88	1/1399 (0.1%)	0.76	0/1881
1	B	0.92	2/1360 (0.1%)	0.79	1/1828 (0.1%)
1	C	0.89	1/1360 (0.1%)	0.83	1/1828 (0.1%)
1	D	0.89	0/1358	0.79	2/1825 (0.1%)
1	E	0.84	0/1366	0.75	0/1836
1	F	0.87	0/1364	0.75	1/1833 (0.1%)
1	G	0.89	1/1363 (0.1%)	0.80	1/1832 (0.1%)
1	H	0.84	0/1358	0.79	0/1825
1	I	0.90	0/1350	0.80	0/1814
1	J	0.85	0/1361	0.79	1/1829 (0.1%)
1	K	0.85	0/1350	0.79	3/1815 (0.2%)
1	L	0.87	0/1350	0.77	1/1814 (0.1%)
All	All	0.87	5/16339 (0.0%)	0.78	11/21960 (0.1%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	1	MET	N-CA	8.05	1.62	1.46
1	B	127	GLU	CD-OE2	-6.02	1.19	1.25
1	B	114	TYR	CD2-CE2	5.32	1.47	1.39
1	G	115	VAL	CB-CG1	5.11	1.63	1.52
1	C	128	GLU	CB-CG	5.06	1.61	1.52

All (11) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	18	GLU	OE1-CD-OE2	-7.09	114.79	123.30
1	J	18	GLU	OE1-CD-OE2	-6.39	115.63	123.30
1	F	18	GLU	OE1-CD-OE2	-6.30	115.74	123.30
1	K	30	ARG	NE-CZ-NH2	-6.15	117.22	120.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	G	18	GLU	OE1-CD-OE2	-6.06	116.03	123.30
1	B	18	GLU	OE1-CD-OE2	-5.94	116.17	123.30
1	K	113	ASP	CB-CG-OD1	5.44	123.19	118.30
1	L	77	LEU	CA-CB-CG	5.20	127.27	115.30
1	D	39	ARG	NE-CZ-NH1	5.19	122.89	120.30
1	D	30	ARG	NE-CZ-NH2	-5.09	117.75	120.30
1	K	56	ASP	CB-CG-OD1	5.07	122.86	118.30

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	1346	0	1348	18	0
1	B	1325	0	1317	9	0
1	C	1325	0	1317	9	0
1	D	1323	0	1316	12	0
1	E	1328	0	1322	7	0
1	F	1326	0	1321	19	0
1	G	1325	0	1322	8	0
1	H	1323	0	1316	9	0
1	I	1318	0	1310	14	0
1	J	1326	0	1317	9	0
1	K	1318	0	1304	10	0
1	L	1318	0	1310	15	0
2	A	45	0	6	9	0
2	B	45	0	6	7	0
2	C	45	0	6	10	0
2	D	45	0	6	11	0
2	E	45	0	6	16	0
2	F	45	0	6	11	0
2	G	45	0	6	12	0
2	H	45	0	6	12	0
2	I	45	0	6	13	0
2	J	45	0	6	10	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	K	45	0	6	10	0
2	L	45	0	6	11	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
3	C	2	0	0	0	0
3	D	2	0	0	0	0
3	E	2	0	0	0	0
3	F	2	0	0	0	0
3	G	2	0	0	0	0
3	H	2	0	0	0	0
3	I	2	0	0	0	0
3	J	2	0	0	0	0
3	K	2	0	0	0	0
3	L	2	0	0	0	0
4	A	15	0	0	0	0
4	B	20	0	0	1	0
4	C	15	0	0	0	0
4	D	10	0	0	0	0
4	E	15	0	0	0	0
4	F	5	0	0	0	0
4	G	10	0	0	0	0
4	H	15	0	0	0	0
4	I	10	0	0	0	0
4	J	10	0	0	1	0
4	K	10	0	0	0	0
4	L	10	0	0	0	0
5	A	190	0	0	9	0
5	B	184	0	0	3	0
5	C	189	0	0	1	0
5	D	181	0	0	2	0
5	E	191	0	0	1	0
5	F	179	0	0	0	0
5	G	195	0	0	3	0
5	H	188	0	0	1	0
5	I	194	0	0	2	0
5	J	179	0	0	1	0
5	K	173	0	0	1	0
5	L	165	0	0	0	0
All	All	18818	0	15892	251	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 8.

All (251) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:L:125[B]:ARG:NH1	1:L:125[B]:ARG:HB3	1.36	1.40
1:L:125[B]:ARG:CB	1:L:125[B]:ARG:HH11	1.35	1.35
1:A:46[A]:HIS:CD2	5:A:2072:HOH:O	1.68	1.32
1:A:112:HIS:ND1	5:A:2150:HOH:O	1.82	1.10
1:I:125[B]:ARG:NH1	1:I:125[B]:ARG:HG2	1.64	1.04
1:I:125[B]:ARG:HG2	1:I:125[B]:ARG:HH11	0.85	1.02
1:I:125[B]:ARG:HH11	1:I:125[B]:ARG:CG	1.71	1.02
1:J:96:ASP:HB2	5:J:2127:HOH:O	1.65	0.97
1:D:115:VAL:HG13	1:F:125[A]:ARG:HD2	1.48	0.94
1:L:125[B]:ARG:HH11	1:L:125[B]:ARG:CG	1.81	0.93
1:H:50:ASP:OD2	5:H:2066:HOH:O	1.88	0.89
1:L:125[B]:ARG:HH11	1:L:125[B]:ARG:HB3	0.73	0.88
1:F:125[B]:ARG:HH11	1:F:125[B]:ARG:HG2	1.38	0.86
1:A:46[A]:HIS:NE2	5:A:2072:HOH:O	1.73	0.85
1:F:125[B]:ARG:HH11	1:F:125[B]:ARG:CG	1.88	0.85
1:A:112:HIS:CE1	5:A:2150:HOH:O	2.21	0.85
1:A:9[A]:ASN:ND2	5:A:2012:HOH:O	2.19	0.74
1:C:20:VAL:HG13	1:C:77:LEU:HD12	1.72	0.72
1:A:96[A]:ASP:OD2	5:A:2134:HOH:O	2.09	0.69
1:K:42:ASP:O	1:K:46[A]:HIS:HD2	1.76	0.68
1:L:125[B]:ARG:CB	1:L:125[B]:ARG:NH1	2.17	0.66
1:D:20:VAL:HG13	1:D:77:LEU:HD12	1.78	0.66
1:B:125[A]:ARG:NH2	4:B:1160:SO4:O4	2.29	0.65
1:E:20:VAL:HG13	1:E:77:LEU:HD23	1.79	0.65
1:L:42:ASP:O	1:L:46[A]:HIS:HD2	1.81	0.64
2:E:200[A]:HEM:HBC1	1:F:26:PHE:CE1	2.32	0.63
1:A:9[B]:ASN:ND2	5:A:2013:HOH:O	2.20	0.63
1:C:125[B]:ARG:HH11	1:C:125[B]:ARG:CG	2.12	0.62
1:L:42:ASP:O	1:L:46[A]:HIS:CD2	2.53	0.61
1:B:147:GLN:NE2	5:B:2160:HOH:O	2.19	0.60
1:D:43:VAL:HG11	1:D:133:TRP:CE2	2.37	0.60
1:L:125[B]:ARG:NH1	1:L:125[B]:ARG:CG	2.49	0.60
1:E:9[A]:ASN:ND2	5:E:2015:HOH:O	2.36	0.59
1:K:43:VAL:O	1:K:47[A]:GLU:HG2	2.02	0.58
1:B:9[A]:ASN:ND2	5:B:2014:HOH:O	2.36	0.58
1:A:43:VAL:HG11	1:A:133:TRP:CE2	2.38	0.58
1:B:20:VAL:HG13	1:B:77:LEU:HD12	1.86	0.57
1:F:125[B]:ARG:HG2	1:F:125[B]:ARG:NH1	2.14	0.57
1:I:42:ASP:O	1:I:46[A]:HIS:HD2	1.87	0.57
1:D:118:ASP:OD2	1:F:125[A]:ARG:CZ	2.54	0.56

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:G:125[B]:ARG:NH2	5:G:2158:HOH:O	2.38	0.56
1:I:76:LYS:HG3	5:I:2112:HOH:O	2.06	0.55
1:F:20:VAL:HG13	1:F:77:LEU:HD23	1.89	0.55
1:H:20:VAL:HG13	1:H:77:LEU:HD23	1.89	0.55
1:K:46[B]:HIS:CD2	5:K:2065:HOH:O	2.60	0.54
1:D:115:VAL:HG13	1:F:125[A]:ARG:HH11	1.73	0.54
2:E:200[A]:HEM:CBC	1:F:26:PHE:CE1	2.91	0.54
1:C:47:GLU:O	1:C:51:GLU:HG2	2.09	0.53
1:G:43:VAL:HG11	1:G:133:TRP:CE2	2.44	0.53
1:C:43:VAL:O	1:C:47:GLU:HG2	2.09	0.53
1:H:125[B]:ARG:HG2	1:H:125[B]:ARG:HH21	1.74	0.52
1:G:26:PHE:CE1	2:H:200[A]:HEM:HBC1	2.44	0.52
1:C:125[B]:ARG:HH11	1:C:125[B]:ARG:HG3	1.73	0.52
1:C:43:VAL:HG11	1:C:133:TRP:CE2	2.45	0.52
1:D:43:VAL:O	1:D:47:GLU:HG2	2.10	0.52
1:F:125[B]:ARG:NH1	1:F:125[B]:ARG:CG	2.59	0.52
1:B:43:VAL:HG11	1:B:133:TRP:CE2	2.45	0.51
1:I:42:ASP:O	1:I:46[A]:HIS:CD2	2.63	0.51
1:A:154:ILE:HD11	1:D:152:ALA:HB1	1.93	0.51
1:F:43:VAL:HG11	1:F:133:TRP:CE2	2.45	0.51
1:J:42:ASP:O	1:J:46[A]:HIS:HD2	1.94	0.51
1:L:47:GLU:O	1:L:51:GLU:HG2	2.11	0.50
1:K:42:ASP:O	1:K:46[A]:HIS:CD2	2.60	0.50
2:I:200[A]:HEM:HBC1	1:J:26:PHE:CE1	2.47	0.49
1:A:43:VAL:O	1:A:47:GLU:HG2	2.12	0.49
1:G:27:LEU:HD23	1:G:79:ILE:HD12	1.93	0.49
1:G:43:VAL:O	1:G:47:GLU:HG2	2.13	0.49
1:B:42:ASP:O	1:B:46[A]:HIS:HD2	1.95	0.49
1:A:131:ILE:O	1:A:135:GLU:HG3	2.14	0.48
2:E:200[A]:HEM:HBC2	1:F:26:PHE:CE2	2.49	0.48
1:I:76:LYS:HE2	5:I:2040:HOH:O	2.13	0.47
1:J:140:LEU:O	1:J:144:MET:HG2	2.14	0.47
1:G:26:PHE:CE1	2:H:200[A]:HEM:CBC	2.97	0.47
1:L:38:LYS:HB3	1:L:156:GLU:HG2	1.95	0.47
1:A:20:VAL:HG13	1:A:77:LEU:HD23	1.94	0.47
1:I:125[B]:ARG:NH1	1:I:125[B]:ARG:CG	2.43	0.47
1:H:125[B]:ARG:HG2	1:H:125[B]:ARG:NH2	2.29	0.47
1:J:43:VAL:O	1:J:47[A]:GLU:HG2	2.15	0.47
1:D:99:LYS:HE3	1:D:99:LYS:HB2	1.38	0.47
1:G:46[B]:HIS:CD2	5:G:2075:HOH:O	2.67	0.47
1:I:43:VAL:HG11	1:I:133:TRP:CE2	2.49	0.47

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:152:ALA:HB1	1:I:154:ILE:HD11	1.97	0.47
1:A:73:ASP:C	5:A:2106:HOH:O	2.54	0.46
1:D:2:LYS:HG2	5:D:2085:HOH:O	2.15	0.46
1:A:9[B]:ASN:OD1	5:A:2013:HOH:O	2.20	0.46
1:G:6:LYS:HE2	5:G:2007:HOH:O	2.15	0.46
1:H:43:VAL:O	1:H:47:GLU:HG2	2.16	0.46
1:J:43:VAL:HG11	1:J:133:TRP:CE2	2.51	0.46
1:D:105:ILE:HG23	1:D:117:ARG:HG3	1.97	0.45
1:B:81:GLU:CD	5:B:2111:HOH:O	2.53	0.45
1:F:43:VAL:O	1:F:47:GLU:HG2	2.16	0.45
1:L:43:VAL:HG11	1:L:133:TRP:CE2	2.51	0.45
2:I:200[A]:HEM:CBC	1:J:26:PHE:CE1	2.99	0.45
1:K:43:VAL:HG11	1:K:133:TRP:CE2	2.51	0.45
1:K:99:LYS:HB2	1:K:99:LYS:HE3	1.49	0.45
1:H:115:VAL:HG13	1:L:125[B]:ARG:HG2	1.98	0.45
5:D:2039:HOH:O	1:K:143:LYS:HE3	2.17	0.45
1:A:105:ILE:HG23	1:A:117:ARG:HG3	1.99	0.45
2:E:200[A]:HEM:CBC	1:F:26:PHE:CD1	3.00	0.45
1:B:142[B]:GLN:HA	1:B:142[B]:GLN:HE21	1.83	0.44
1:L:26:PHE:CE1	2:L:200[A]:HEM:HBC1	2.53	0.44
1:L:20:VAL:HG13	1:L:77:LEU:HD12	2.00	0.44
1:H:43:VAL:HG11	1:H:133:TRP:CE2	2.53	0.44
2:E:200[A]:HEM:HBC1	1:F:26:PHE:CD1	2.52	0.43
1:K:27:LEU:HD23	1:K:79:ILE:HD12	2.00	0.43
1:J:147:GLN:HG3	4:J:1159:SO4:O1	2.19	0.43
1:K:140:LEU:O	1:K:144:MET:HG2	2.18	0.43
1:A:43:VAL:HG11	1:A:133:TRP:CZ2	2.54	0.43
1:C:125[B]:ARG:CG	1:C:125[B]:ARG:NH1	2.78	0.42
1:E:43:VAL:HG11	1:E:133:TRP:CE2	2.54	0.42
1:J:99:LYS:HE3	1:J:99:LYS:HB2	1.85	0.42
1:D:115:VAL:CG1	1:F:125[A]:ARG:HH11	2.33	0.42
1:C:93:LEU:HD22	5:C:2125:HOH:O	2.20	0.41
1:F:109:ASP:HB2	1:F:117:ARG:HH11	1.86	0.41
1:E:125[B]:ARG:HB3	1:E:125[B]:ARG:HE	1.64	0.41
1:L:43:VAL:O	1:L:47:GLU:HG2	2.20	0.41
1:I:20:VAL:HG13	1:I:77:LEU:HD12	2.03	0.41
2:E:200[A]:HEM:HBC2	1:F:26:PHE:CZ	2.55	0.41
1:I:22:ILE:HD11	1:I:52:MET:HA	2.02	0.41
2:E:200[A]:HEM:CBC	1:F:26:PHE:CZ	3.04	0.40
1:A:125[A]:ARG:CZ	1:I:118:ASP:OD2	2.70	0.40
1:E:109:ASP:HB2	1:E:117:ARG:HH11	1.86	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:I:47:GLU:O	1:I:51:GLU:HG2	2.21	0.40
1:K:20:VAL:HG13	1:K:77:LEU:HD23	2.04	0.40

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	166/158 (105%)	166 (100%)	0	0	100	100
1	B	160/158 (101%)	159 (99%)	1 (1%)	0	100	100
1	C	160/158 (101%)	160 (100%)	0	0	100	100
1	D	160/158 (101%)	160 (100%)	0	0	100	100
1	E	161/158 (102%)	161 (100%)	0	0	100	100
1	F	161/158 (102%)	161 (100%)	0	0	100	100
1	G	161/158 (102%)	161 (100%)	0	0	100	100
1	H	160/158 (101%)	159 (99%)	1 (1%)	0	100	100
1	I	159/158 (101%)	159 (100%)	0	0	100	100
1	J	160/158 (101%)	158 (99%)	2 (1%)	0	100	100
1	K	159/158 (101%)	157 (99%)	2 (1%)	0	100	100
1	L	159/158 (101%)	159 (100%)	0	0	100	100
All	All	1926/1896 (102%)	1920 (100%)	6 (0%)	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	150/139 (108%)	145 (97%)	5 (3%)	38 28
1	B	144/139 (104%)	143 (99%)	1 (1%)	84 83
1	C	144/139 (104%)	143 (99%)	1 (1%)	84 83
1	D	144/139 (104%)	141 (98%)	3 (2%)	53 46
1	E	145/139 (104%)	143 (99%)	2 (1%)	67 63
1	F	145/139 (104%)	143 (99%)	2 (1%)	67 63
1	G	145/139 (104%)	144 (99%)	1 (1%)	84 83
1	H	144/139 (104%)	144 (100%)	0	100 100
1	I	143/139 (103%)	139 (97%)	4 (3%)	43 34
1	J	144/139 (104%)	144 (100%)	0	100 100
1	K	143/139 (103%)	140 (98%)	3 (2%)	53 46
1	L	143/139 (103%)	141 (99%)	2 (1%)	67 63
All	All	1734/1668 (104%)	1710 (99%)	24 (1%)	76 63

All (24) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	2	LYS
1	A	5[A]	THR
1	A	5[B]	THR
1	A	142[A]	GLN
1	A	142[B]	GLN
1	B	154	ILE
1	C	93	LEU
1	D	77	LEU
1	D	78[A]	ASN
1	D	78[B]	ASN
1	E	125[A]	ARG
1	E	125[B]	ARG
1	F	125[A]	ARG
1	F	125[B]	ARG

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Mol	Chain	Res	Type
1	G	155	ARG
1	I	6	LYS
1	I	76	LYS
1	I	142[A]	GLN
1	I	142[B]	GLN
1	K	78[A]	ASN
1	K	78[B]	ASN
1	K	154	ILE
1	L	77	LEU
1	L	99	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (34) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	151	GLN
1	B	17	ASN
1	B	100	ASN
1	B	151	GLN
1	C	17	ASN
1	C	100	ASN
1	C	151	GLN
1	D	112	HIS
1	D	151	GLN
1	E	17	ASN
1	E	100	ASN
1	E	151	GLN
1	F	9	ASN
1	F	151	GLN
1	G	9	ASN
1	G	34	ASN
1	G	151	GLN
1	H	17	ASN
1	H	100	ASN
1	I	9	ASN
1	I	17	ASN
1	I	100	ASN
1	I	151	GLN
1	J	9	ASN
1	J	17	ASN
1	J	100	ASN
1	J	151	GLN
1	K	17	ASN

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Mol	Chain	Res	Type
1	K	100	ASN
1	K	112	HIS
1	K	151	GLN
1	L	17	ASN
1	L	100	ASN
1	L	151	GLN

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 carbohydrates in this entry.

5.6 Ligand geometry

Of 77 ligands modelled in this entry, 24 are monoatomic - leaving 53 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	$\#\|Z\| > 2$
2	HEM	G	200[B]	-	27,50,50	2.74	5 (18%)	17,82,82	2.38	7 (41%)
4	SO4	C	1158	-	4,4,4	0.26	0	6,6,6	0.64	0
2	HEM	F	200[A]	-	27,50,50	2.57	4 (14%)	17,82,82	1.94	7 (41%)
2	HEM	F	200[B]	2	27,50,50	2.59	5 (18%)	17,82,82	1.94	7 (41%)
4	SO4	J	1158	-	4,4,4	0.21	0	6,6,6	0.27	0
4	SO4	D	1158	-	4,4,4	0.20	0	6,6,6	0.39	0
4	SO4	B	1159	-	4,4,4	0.18	0	6,6,6	0.41	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
4	SO4	B	1161	-	4,4,4	0.19	0	6,6,6	0.27	0
4	SO4	C	1160	-	4,4,4	0.11	0	6,6,6	0.28	0
4	SO4	E	1160	-	4,4,4	0.41	0	6,6,6	0.48	0
2	HEM	H	200[A]	-	27,50,50	2.54	7 (25%)	17,82,82	1.97	5 (29%)
4	SO4	L	1158	-	4,4,4	0.34	0	6,6,6	0.31	0
4	SO4	G	1158	-	4,4,4	0.31	0	6,6,6	0.33	0
4	SO4	A	1158	-	4,4,4	0.29	0	6,6,6	0.36	0
4	SO4	L	1159	-	4,4,4	0.15	0	6,6,6	0.30	0
2	HEM	E	200[B]	-	27,50,50	2.72	7 (25%)	17,82,82	1.64	4 (23%)
2	HEM	H	200[B]	-	27,50,50	2.57	8 (29%)	17,82,82	1.97	5 (29%)
2	HEM	E	200[A]	-	27,50,50	2.69	6 (22%)	17,82,82	1.64	4 (23%)
4	SO4	B	1158	-	4,4,4	0.19	0	6,6,6	0.26	0
4	SO4	A	1159	-	4,4,4	0.24	0	6,6,6	0.21	0
4	SO4	A	1160	-	4,4,4	0.17	0	6,6,6	0.20	0
4	SO4	G	1159	-	4,4,4	0.28	0	6,6,6	0.42	0
2	HEM	K	200[B]	-	27,50,50	2.45	7 (25%)	17,82,82	2.34	8 (47%)
4	SO4	K	1159	-	4,4,4	0.26	0	6,6,6	0.41	0
4	SO4	B	1160	-	4,4,4	0.10	0	6,6,6	0.27	0
2	HEM	K	200[A]	-	27,50,50	2.45	6 (22%)	17,82,82	2.34	8 (47%)
2	HEM	D	200[B]	-	27,50,50	2.45	6 (22%)	17,82,82	1.50	4 (23%)
4	SO4	H	1160	-	4,4,4	0.28	0	6,6,6	0.55	0
4	SO4	J	1159	-	4,4,4	0.08	0	6,6,6	0.30	0
2	HEM	L	200[B]	-	27,50,50	2.46	6 (22%)	17,82,82	1.71	4 (23%)
2	HEM	D	200[A]	-	27,50,50	2.45	6 (22%)	17,82,82	1.50	4 (23%)
2	HEM	B	200[A]	-	27,50,50	2.33	6 (22%)	17,82,82	1.66	3 (17%)
2	HEM	B	200[B]	-	27,50,50	2.37	7 (25%)	17,82,82	1.66	3 (17%)
2	HEM	L	200[A]	-	27,50,50	2.45	6 (22%)	17,82,82	1.71	4 (23%)
4	SO4	C	1159	-	4,4,4	0.32	0	6,6,6	0.56	0
2	HEM	J	200[A]	-	27,50,50	2.61	6 (22%)	17,82,82	1.86	3 (17%)
4	SO4	E	1159	-	4,4,4	0.31	0	6,6,6	0.46	0
2	HEM	J	200[B]	-	27,50,50	2.60	6 (22%)	17,82,82	1.86	3 (17%)
4	SO4	I	1159	-	4,4,4	0.16	0	6,6,6	0.27	0
4	SO4	H	1159	-	4,4,4	0.10	0	6,6,6	0.49	0
4	SO4	E	1158	-	4,4,4	0.23	0	6,6,6	0.37	0
4	SO4	I	1158	-	4,4,4	0.24	0	6,6,6	0.38	0
2	HEM	A	200[B]	-	27,50,50	2.51	7 (25%)	17,82,82	1.58	4 (23%)
2	HEM	C	200[A]	-	27,50,50	2.37	10 (37%)	17,82,82	2.07	8 (47%)
2	HEM	A	200[A]	-	27,50,50	2.51	7 (25%)	17,82,82	1.58	4 (23%)
4	SO4	K	1158	-	4,4,4	0.21	0	6,6,6	0.23	0

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
2	HEM	I	200[B]	2	27,50,50	2.54	7 (25%)	17,82,82	1.95	5 (29%)
4	SO4	H	1158	-	4,4,4	0.30	0	6,6,6	0.34	0
2	HEM	C	200[B]	-	27,50,50	2.35	9 (33%)	17,82,82	2.07	8 (47%)
2	HEM	I	200[A]	-	27,50,50	2.52	7 (25%)	17,82,82	1.95	5 (29%)
4	SO4	F	1158	-	4,4,4	0.16	0	6,6,6	0.49	0
2	HEM	G	200[A]	-	27,50,50	2.64	4 (14%)	17,82,82	2.38	7 (41%)
4	SO4	D	1159	-	4,4,4	0.17	0	6,6,6	0.22	0

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	HEM	G	200[B]	-	-	0/6/54/54	-
2	HEM	F	200[A]	-	-	0/6/54/54	-
2	HEM	F	200[B]	2	-	0/6/54/54	-
2	HEM	H	200[B]	-	-	0/6/54/54	-
2	HEM	H	200[A]	-	-	0/6/54/54	-
2	HEM	E	200[B]	-	-	0/6/54/54	-
2	HEM	E	200[A]	-	-	0/6/54/54	-
2	HEM	K	200[A]	-	-	0/6/54/54	-
2	HEM	K	200[B]	-	-	0/6/54/54	-
2	HEM	D	200[B]	-	-	0/6/54/54	-
2	HEM	L	200[B]	-	-	0/6/54/54	-
2	HEM	D	200[A]	-	-	0/6/54/54	-
2	HEM	B	200[A]	-	-	0/6/54/54	-
2	HEM	B	200[B]	-	-	0/6/54/54	-
2	HEM	L	200[A]	-	-	0/6/54/54	-
2	HEM	J	200[A]	-	-	0/6/54/54	-
2	HEM	J	200[B]	-	-	0/6/54/54	-
2	HEM	A	200[B]	-	-	0/6/54/54	-
2	HEM	C	200[A]	-	-	0/6/54/54	-
2	HEM	A	200[A]	-	-	0/6/54/54	-
2	HEM	I	200[B]	2	-	0/6/54/54	-
2	HEM	C	200[B]	-	-	0/6/54/54	-
2	HEM	I	200[A]	-	-	0/6/54/54	-
2	HEM	G	200[A]	-	-	0/6/54/54	-

All (155) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	I	200[B]	HEM	C3C-C2C	-8.61	1.28	1.40
2	I	200[A]	HEM	C3C-C2C	-8.61	1.28	1.40
2	G	200[B]	HEM	C3C-C2C	-8.38	1.28	1.40
2	G	200[A]	HEM	C3C-C2C	-8.38	1.28	1.40
2	J	200[A]	HEM	C3B-C2B	-8.31	1.28	1.40
2	J	200[B]	HEM	C3B-C2B	-8.31	1.28	1.40
2	F	200[A]	HEM	C3C-C2C	-8.05	1.29	1.40
2	F	200[B]	HEM	C3C-C2C	-8.05	1.29	1.40
2	D	200[B]	HEM	C3C-C2C	-8.03	1.29	1.40
2	D	200[A]	HEM	C3C-C2C	-8.03	1.29	1.40
2	J	200[A]	HEM	C3C-C2C	-8.02	1.29	1.40
2	J	200[B]	HEM	C3C-C2C	-8.02	1.29	1.40
2	E	200[B]	HEM	C3C-C2C	-7.78	1.29	1.40
2	E	200[A]	HEM	C3C-C2C	-7.78	1.29	1.40
2	L	200[B]	HEM	C3C-C2C	-7.66	1.29	1.40
2	L	200[A]	HEM	C3C-C2C	-7.66	1.29	1.40
2	G	200[B]	HEM	C3B-C2B	-7.65	1.29	1.40
2	G	200[A]	HEM	C3B-C2B	-7.65	1.29	1.40
2	A	200[B]	HEM	C3B-C2B	-7.64	1.29	1.40
2	A	200[A]	HEM	C3B-C2B	-7.64	1.29	1.40
2	E	200[B]	HEM	C3B-C2B	-7.60	1.29	1.40
2	E	200[A]	HEM	C3B-C2B	-7.60	1.29	1.40
2	H	200[B]	HEM	C3C-C2C	-7.48	1.30	1.40
2	H	200[A]	HEM	C3C-C2C	-7.48	1.30	1.40
2	B	200[A]	HEM	C3C-C2C	-7.46	1.30	1.40
2	B	200[B]	HEM	C3C-C2C	-7.46	1.30	1.40
2	C	200[A]	HEM	C3C-C2C	-7.46	1.30	1.40
2	C	200[B]	HEM	C3C-C2C	-7.46	1.30	1.40
2	A	200[B]	HEM	C3C-C2C	-7.44	1.30	1.40
2	A	200[A]	HEM	C3C-C2C	-7.44	1.30	1.40
2	K	200[A]	HEM	C3C-C2C	-7.42	1.30	1.40
2	K	200[B]	HEM	C3C-C2C	-7.42	1.30	1.40
2	F	200[A]	HEM	C3B-C2B	-7.28	1.30	1.40
2	F	200[B]	HEM	C3B-C2B	-7.28	1.30	1.40
2	H	200[B]	HEM	C3B-C2B	-7.21	1.30	1.40
2	H	200[A]	HEM	C3B-C2B	-7.21	1.30	1.40
2	K	200[A]	HEM	C3B-C2B	-7.15	1.30	1.40
2	K	200[B]	HEM	C3B-C2B	-7.15	1.30	1.40
2	L	200[B]	HEM	C3B-C2B	-7.11	1.30	1.40
2	L	200[A]	HEM	C3B-C2B	-7.11	1.30	1.40
2	B	200[A]	HEM	C3B-C2B	-6.69	1.31	1.40
2	B	200[B]	HEM	C3B-C2B	-6.69	1.31	1.40
2	D	200[B]	HEM	C3B-C2B	-6.66	1.31	1.40

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	200[A]	HEM	C3B-C2B	-6.66	1.31	1.40
2	I	200[B]	HEM	C3B-C2B	-6.39	1.31	1.40
2	I	200[A]	HEM	C3B-C2B	-6.39	1.31	1.40
2	C	200[A]	HEM	C3B-C2B	-5.22	1.33	1.40
2	C	200[B]	HEM	C3B-C2B	-5.22	1.33	1.40
2	L	200[B]	HEM	C3D-C2D	4.52	1.51	1.37
2	L	200[A]	HEM	C3D-C2D	4.52	1.51	1.37
2	I	200[B]	HEM	C3D-C2D	4.48	1.50	1.37
2	I	200[A]	HEM	C3D-C2D	4.48	1.50	1.37
2	E	200[B]	HEM	C3D-C2D	4.48	1.50	1.37
2	E	200[A]	HEM	C3D-C2D	4.48	1.50	1.37
2	F	200[A]	HEM	C3D-C2D	4.39	1.50	1.37
2	F	200[B]	HEM	C3D-C2D	4.39	1.50	1.37
2	H	200[B]	HEM	C3D-C2D	4.37	1.50	1.37
2	H	200[A]	HEM	C3D-C2D	4.37	1.50	1.37
2	A	200[B]	HEM	C3D-C2D	4.35	1.50	1.37
2	A	200[A]	HEM	C3D-C2D	4.35	1.50	1.37
2	C	200[A]	HEM	C3D-C2D	4.29	1.50	1.37
2	C	200[B]	HEM	C3D-C2D	4.29	1.50	1.37
2	K	200[A]	HEM	C3D-C2D	4.23	1.50	1.37
2	K	200[B]	HEM	C3D-C2D	4.23	1.50	1.37
2	D	200[B]	HEM	C3D-C2D	4.23	1.50	1.37
2	D	200[A]	HEM	C3D-C2D	4.23	1.50	1.37
2	E	200[B]	HEM	CBB-CAB	-4.20	1.01	1.29
2	E	200[A]	HEM	CBB-CAB	-4.20	1.01	1.29
2	G	200[B]	HEM	C3D-C2D	4.19	1.50	1.37
2	G	200[A]	HEM	C3D-C2D	4.19	1.50	1.37
2	F	200[A]	HEM	CBB-CAB	-4.11	1.01	1.29
2	F	200[B]	HEM	CBB-CAB	-4.11	1.01	1.29
2	G	200[B]	HEM	C3C-CAC	4.08	1.56	1.47
2	J	200[A]	HEM	C3D-C2D	3.95	1.49	1.37
2	J	200[B]	HEM	C3D-C2D	3.95	1.49	1.37
2	B	200[A]	HEM	C3D-C2D	3.92	1.49	1.37
2	B	200[B]	HEM	C3D-C2D	3.92	1.49	1.37
2	H	200[B]	HEM	CBB-CAB	-3.89	1.03	1.29
2	H	200[A]	HEM	CBB-CAB	-3.89	1.03	1.29
2	G	200[B]	HEM	CBB-CAB	-3.85	1.03	1.29
2	G	200[A]	HEM	CBB-CAB	-3.85	1.03	1.29
2	E	200[B]	HEM	C3B-CAB	3.84	1.55	1.47
2	E	200[A]	HEM	C3B-CAB	3.84	1.55	1.47
2	H	200[B]	HEM	C3B-CAB	2.85	1.53	1.47
2	H	200[A]	HEM	C3B-CAB	2.85	1.53	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	200[A]	HEM	C3B-CAB	2.70	1.53	1.47
2	C	200[B]	HEM	C3B-CAB	2.70	1.53	1.47
2	B	200[A]	HEM	C1D-CHD	-2.70	1.33	1.41
2	B	200[B]	HEM	C1D-CHD	-2.70	1.33	1.41
2	J	200[A]	HEM	C3B-CAB	2.55	1.53	1.47
2	J	200[B]	HEM	C3B-CAB	2.55	1.53	1.47
2	C	200[A]	HEM	C4B-CHC	-2.46	1.34	1.41
2	C	200[B]	HEM	C4B-CHC	-2.46	1.34	1.41
2	A	200[B]	HEM	C4A-CHB	-2.41	1.34	1.41
2	A	200[A]	HEM	C4A-CHB	-2.41	1.34	1.41
2	E	200[B]	HEM	C4B-CHC	-2.35	1.34	1.41
2	E	200[A]	HEM	C4B-CHC	-2.35	1.34	1.41
2	I	200[B]	HEM	C2A-C3A	-2.35	1.30	1.37
2	I	200[A]	HEM	C2A-C3A	-2.35	1.30	1.37
2	I	200[B]	HEM	C1D-CHD	-2.33	1.34	1.41
2	I	200[A]	HEM	C1D-CHD	-2.33	1.34	1.41
2	A	200[B]	HEM	C2A-C3A	-2.30	1.30	1.37
2	A	200[A]	HEM	C2A-C3A	-2.30	1.30	1.37
2	I	200[B]	HEM	C4A-CHB	-2.29	1.34	1.41
2	I	200[A]	HEM	C4A-CHB	-2.29	1.34	1.41
2	I	200[B]	HEM	C4B-CHC	-2.28	1.34	1.41
2	I	200[A]	HEM	C4B-CHC	-2.28	1.34	1.41
2	L	200[B]	HEM	C3B-CAB	2.27	1.52	1.47
2	L	200[A]	HEM	C3B-CAB	2.27	1.52	1.47
2	B	200[B]	HEM	C3C-CAC	2.26	1.52	1.47
2	A	200[B]	HEM	C1D-CHD	-2.18	1.34	1.41
2	A	200[A]	HEM	C1D-CHD	-2.18	1.34	1.41
2	A	200[B]	HEM	C4B-CHC	-2.17	1.35	1.41
2	A	200[A]	HEM	C4B-CHC	-2.17	1.35	1.41
2	J	200[A]	HEM	C4A-CHB	-2.16	1.35	1.41
2	J	200[B]	HEM	C4A-CHB	-2.16	1.35	1.41
2	C	200[A]	HEM	C4A-CHB	-2.15	1.35	1.41
2	C	200[B]	HEM	C4A-CHB	-2.15	1.35	1.41
2	B	200[A]	HEM	C4A-CHB	-2.13	1.35	1.41
2	B	200[B]	HEM	C4A-CHB	-2.13	1.35	1.41
2	D	200[B]	HEM	C4A-NA	-2.12	1.31	1.36
2	D	200[A]	HEM	C4A-NA	-2.12	1.31	1.36
2	D	200[B]	HEM	C4A-CHB	-2.12	1.35	1.41
2	D	200[A]	HEM	C4A-CHB	-2.12	1.35	1.41
2	C	200[A]	HEM	C3C-CAC	-2.11	1.43	1.47
2	C	200[A]	HEM	C1B-C2B	2.11	1.47	1.42
2	C	200[B]	HEM	C1B-C2B	2.11	1.47	1.42

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	L	200[B]	HEM	C1D-CHD	-2.10	1.35	1.41
2	L	200[A]	HEM	C1D-CHD	-2.10	1.35	1.41
2	K	200[A]	HEM	C4B-CHC	-2.09	1.35	1.41
2	K	200[B]	HEM	C4B-CHC	-2.09	1.35	1.41
2	C	200[A]	HEM	C2A-C3A	-2.09	1.31	1.37
2	C	200[B]	HEM	C2A-C3A	-2.09	1.31	1.37
2	C	200[A]	HEM	CBB-CAB	-2.08	1.15	1.29
2	C	200[B]	HEM	CBB-CAB	-2.08	1.15	1.29
2	H	200[B]	HEM	C1A-CHA	-2.08	1.35	1.41
2	H	200[A]	HEM	C1A-CHA	-2.08	1.35	1.41
2	L	200[B]	HEM	C4B-CHC	-2.06	1.35	1.41
2	L	200[A]	HEM	C4B-CHC	-2.06	1.35	1.41
2	B	200[A]	HEM	C3B-CAB	2.06	1.52	1.47
2	B	200[B]	HEM	C3B-CAB	2.06	1.52	1.47
2	J	200[A]	HEM	C1D-CHD	-2.06	1.35	1.41
2	J	200[B]	HEM	C1D-CHD	-2.06	1.35	1.41
2	K	200[B]	HEM	C3C-CAC	2.05	1.52	1.47
2	E	200[B]	HEM	C3C-CAC	2.04	1.52	1.47
2	K	200[A]	HEM	C1A-CHA	-2.04	1.35	1.41
2	K	200[B]	HEM	C1A-CHA	-2.04	1.35	1.41
2	H	200[B]	HEM	C3C-CAC	2.03	1.52	1.47
2	F	200[B]	HEM	C3C-CAC	2.03	1.52	1.47
2	K	200[A]	HEM	C1D-CHD	-2.02	1.35	1.41
2	K	200[B]	HEM	C1D-CHD	-2.02	1.35	1.41
2	H	200[B]	HEM	C2A-C3A	-2.02	1.31	1.37
2	H	200[A]	HEM	C2A-C3A	-2.02	1.31	1.37
2	D	200[B]	HEM	C1D-CHD	-2.00	1.35	1.41
2	D	200[A]	HEM	C1D-CHD	-2.00	1.35	1.41

All (124) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	G	200[B]	HEM	C4C-C3C-C2C	6.32	111.31	106.90
2	G	200[A]	HEM	C4C-C3C-C2C	6.32	111.31	106.90
2	J	200[A]	HEM	CBD-CAD-C3D	-4.79	103.66	112.48
2	J	200[B]	HEM	CBD-CAD-C3D	-4.79	103.66	112.48
2	K	200[A]	HEM	C1D-C2D-C3D	-4.62	103.78	107.00
2	K	200[B]	HEM	C1D-C2D-C3D	-4.62	103.78	107.00
2	J	200[A]	HEM	C4C-C3C-C2C	4.30	109.90	106.90
2	J	200[B]	HEM	C4C-C3C-C2C	4.30	109.90	106.90
2	F	200[A]	HEM	C4C-C3C-C2C	4.26	109.87	106.90
2	F	200[B]	HEM	C4C-C3C-C2C	4.26	109.87	106.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	H	200[B]	HEM	C4C-C3C-C2C	4.20	109.83	106.90
2	H	200[A]	HEM	C4C-C3C-C2C	4.20	109.83	106.90
2	I	200[B]	HEM	CBA-CAA-C2A	-4.13	104.86	112.49
2	I	200[A]	HEM	CBA-CAA-C2A	-4.13	104.86	112.49
2	C	200[A]	HEM	C1D-C2D-C3D	-4.04	104.19	107.00
2	C	200[B]	HEM	C1D-C2D-C3D	-4.04	104.19	107.00
2	K	200[A]	HEM	C4C-C3C-C2C	3.92	109.63	106.90
2	K	200[B]	HEM	C4C-C3C-C2C	3.92	109.63	106.90
2	G	200[B]	HEM	CBD-CAD-C3D	-3.83	105.42	112.48
2	G	200[A]	HEM	CBD-CAD-C3D	-3.83	105.42	112.48
2	B	200[A]	HEM	CBD-CAD-C3D	-3.81	105.47	112.48
2	B	200[B]	HEM	CBD-CAD-C3D	-3.81	105.47	112.48
2	H	200[B]	HEM	CBD-CAD-C3D	-3.78	105.51	112.48
2	H	200[A]	HEM	CBD-CAD-C3D	-3.78	105.51	112.48
2	C	200[A]	HEM	C4C-C3C-C2C	3.74	109.51	106.90
2	C	200[B]	HEM	C4C-C3C-C2C	3.74	109.51	106.90
2	L	200[B]	HEM	CBD-CAD-C3D	-3.74	105.59	112.48
2	L	200[A]	HEM	CBD-CAD-C3D	-3.74	105.59	112.48
2	K	200[A]	HEM	CBD-CAD-C3D	-3.67	105.71	112.48
2	K	200[B]	HEM	CBD-CAD-C3D	-3.67	105.71	112.48
2	B	200[A]	HEM	C1D-C2D-C3D	-3.65	104.46	107.00
2	B	200[B]	HEM	C1D-C2D-C3D	-3.65	104.46	107.00
2	A	200[B]	HEM	C4C-C3C-C2C	3.61	109.42	106.90
2	A	200[A]	HEM	C4C-C3C-C2C	3.61	109.42	106.90
2	I	200[B]	HEM	C4C-C3C-C2C	3.60	109.41	106.90
2	I	200[A]	HEM	C4C-C3C-C2C	3.60	109.41	106.90
2	E	200[B]	HEM	C4C-C3C-C2C	3.39	109.27	106.90
2	E	200[A]	HEM	C4C-C3C-C2C	3.39	109.27	106.90
2	L	200[B]	HEM	C4C-C3C-C2C	3.38	109.26	106.90
2	L	200[A]	HEM	C4C-C3C-C2C	3.38	109.26	106.90
2	H	200[B]	HEM	CBA-CAA-C2A	-3.31	106.38	112.49
2	H	200[A]	HEM	CBA-CAA-C2A	-3.31	106.38	112.49
2	I	200[B]	HEM	CMB-C2B-C3B	3.30	130.85	124.68
2	I	200[A]	HEM	CMB-C2B-C3B	3.30	130.85	124.68
2	C	200[A]	HEM	C3B-C4B-NB	-3.25	105.00	109.21
2	C	200[B]	HEM	C3B-C4B-NB	-3.25	105.00	109.21
2	F	200[A]	HEM	CMA-C3A-C4A	-3.25	123.47	128.46
2	F	200[B]	HEM	CMA-C3A-C4A	-3.25	123.47	128.46
2	G	200[B]	HEM	CBA-CAA-C2A	-3.23	106.52	112.49
2	G	200[A]	HEM	CBA-CAA-C2A	-3.23	106.52	112.49
2	L	200[B]	HEM	CBA-CAA-C2A	-2.90	107.14	112.49
2	L	200[A]	HEM	CBA-CAA-C2A	-2.90	107.14	112.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	200[B]	HEM	CBA-CAA-C2A	-2.88	107.18	112.49
2	A	200[A]	HEM	CBA-CAA-C2A	-2.88	107.18	112.49
2	I	200[B]	HEM	C1D-C2D-C3D	-2.87	105.00	107.00
2	I	200[A]	HEM	C1D-C2D-C3D	-2.87	105.00	107.00
2	K	200[A]	HEM	CBA-CAA-C2A	-2.84	107.25	112.49
2	K	200[B]	HEM	CBA-CAA-C2A	-2.84	107.25	112.49
2	D	200[B]	HEM	CBD-CAD-C3D	-2.82	107.28	112.48
2	D	200[A]	HEM	CBD-CAD-C3D	-2.82	107.28	112.48
2	C	200[A]	HEM	CBD-CAD-C3D	-2.80	107.32	112.48
2	C	200[B]	HEM	CBD-CAD-C3D	-2.80	107.32	112.48
2	E	200[B]	HEM	CBD-CAD-C3D	-2.79	107.33	112.48
2	E	200[A]	HEM	CBD-CAD-C3D	-2.79	107.33	112.48
2	D	200[B]	HEM	C4C-C3C-C2C	2.75	108.82	106.90
2	D	200[A]	HEM	C4C-C3C-C2C	2.75	108.82	106.90
2	K	200[A]	HEM	CMA-C3A-C4A	-2.64	124.40	128.46
2	K	200[B]	HEM	CMA-C3A-C4A	-2.64	124.40	128.46
2	G	200[B]	HEM	C3C-C4C-NC	-2.64	105.95	110.94
2	G	200[A]	HEM	C3C-C4C-NC	-2.64	105.95	110.94
2	A	200[B]	HEM	C1D-C2D-C3D	-2.62	105.17	107.00
2	A	200[A]	HEM	C1D-C2D-C3D	-2.62	105.17	107.00
2	H	200[B]	HEM	C3C-C4C-NC	-2.59	106.06	110.94
2	H	200[A]	HEM	C3C-C4C-NC	-2.59	106.06	110.94
2	I	200[B]	HEM	CBD-CAD-C3D	-2.49	107.89	112.48
2	I	200[A]	HEM	CBD-CAD-C3D	-2.49	107.89	112.48
2	G	200[B]	HEM	CMA-C3A-C4A	-2.43	124.72	128.46
2	G	200[A]	HEM	CMA-C3A-C4A	-2.43	124.72	128.46
2	L	200[B]	HEM	C1D-C2D-C3D	-2.41	105.32	107.00
2	L	200[A]	HEM	C1D-C2D-C3D	-2.41	105.32	107.00
2	K	200[A]	HEM	CMC-C2C-C3C	2.41	129.19	124.68
2	K	200[B]	HEM	CMC-C2C-C3C	2.41	129.19	124.68
2	D	200[B]	HEM	CMB-C2B-C3B	2.37	129.12	124.68
2	D	200[A]	HEM	CMB-C2B-C3B	2.37	129.12	124.68
2	F	200[A]	HEM	C4A-C3A-C2A	2.32	108.61	107.00
2	F	200[B]	HEM	C4A-C3A-C2A	2.32	108.61	107.00
2	K	200[A]	HEM	CMD-C2D-C1D	2.27	131.95	128.46
2	K	200[B]	HEM	CMD-C2D-C1D	2.27	131.95	128.46
2	J	200[A]	HEM	C1D-C2D-C3D	-2.25	105.43	107.00
2	J	200[B]	HEM	C1D-C2D-C3D	-2.25	105.43	107.00
2	E	200[B]	HEM	C3C-C4C-NC	-2.24	106.72	110.94
2	E	200[A]	HEM	C3C-C4C-NC	-2.24	106.72	110.94
2	A	200[B]	HEM	CBD-CAD-C3D	-2.23	108.36	112.48
2	A	200[A]	HEM	CBD-CAD-C3D	-2.23	108.36	112.48

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	F	200[A]	HEM	C1D-C2D-C3D	-2.20	105.46	107.00
2	F	200[B]	HEM	C1D-C2D-C3D	-2.20	105.46	107.00
2	K	200[A]	HEM	CMB-C2B-C3B	2.15	128.70	124.68
2	K	200[B]	HEM	CMB-C2B-C3B	2.15	128.70	124.68
2	C	200[A]	HEM	CMA-C3A-C4A	-2.15	125.16	128.46
2	C	200[B]	HEM	CMA-C3A-C4A	-2.15	125.16	128.46
2	C	200[A]	HEM	CBA-CAA-C2A	-2.14	108.54	112.49
2	C	200[B]	HEM	CBA-CAA-C2A	-2.14	108.54	112.49
2	B	200[A]	HEM	CMB-C2B-C3B	2.11	128.63	124.68
2	B	200[B]	HEM	CMB-C2B-C3B	2.11	128.63	124.68
2	G	200[B]	HEM	C1D-C2D-C3D	-2.10	105.53	107.00
2	G	200[A]	HEM	C1D-C2D-C3D	-2.10	105.53	107.00
2	C	200[A]	HEM	CMC-C2C-C3C	2.08	128.57	124.68
2	C	200[B]	HEM	CMC-C2C-C3C	2.08	128.57	124.68
2	E	200[B]	HEM	C4A-C3A-C2A	2.08	108.44	107.00
2	E	200[A]	HEM	C4A-C3A-C2A	2.08	108.44	107.00
2	G	200[B]	HEM	C3B-C4B-NB	-2.07	106.53	109.21
2	G	200[A]	HEM	C3B-C4B-NB	-2.07	106.53	109.21
2	D	200[B]	HEM	CBA-CAA-C2A	-2.07	108.67	112.49
2	D	200[A]	HEM	CBA-CAA-C2A	-2.07	108.67	112.49
2	H	200[B]	HEM	C1D-C2D-C3D	-2.06	105.56	107.00
2	H	200[A]	HEM	C1D-C2D-C3D	-2.06	105.56	107.00
2	F	200[A]	HEM	CBD-CAD-C3D	-2.05	108.70	112.48
2	F	200[B]	HEM	CBD-CAD-C3D	-2.05	108.70	112.48
2	C	200[A]	HEM	C4A-C3A-C2A	2.05	108.42	107.00
2	C	200[B]	HEM	C4A-C3A-C2A	2.05	108.42	107.00
2	F	200[A]	HEM	CBA-CAA-C2A	-2.04	108.72	112.49
2	F	200[B]	HEM	CBA-CAA-C2A	-2.04	108.72	112.49
2	F	200[A]	HEM	C3B-C4B-NB	-2.04	106.57	109.21
2	F	200[B]	HEM	C3B-C4B-NB	-2.04	106.57	109.21

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

26 monomers are involved in 134 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	G	200[B]	HEM	7	0
2	F	200[A]	HEM	5	0
2	F	200[B]	HEM	6	0
2	H	200[A]	HEM	8	0

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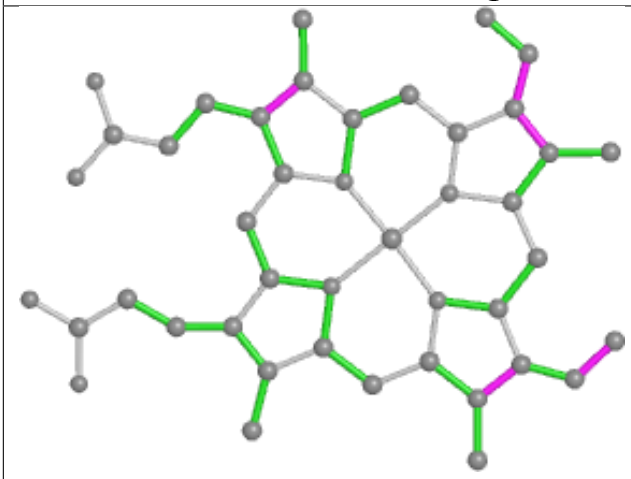
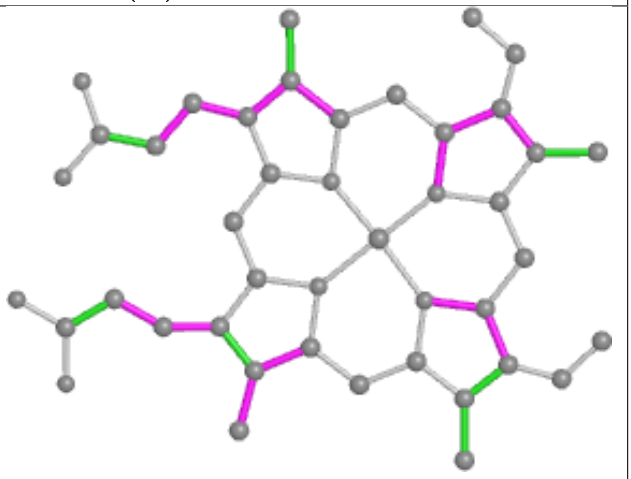
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	E	200[B]	HEM	4	0
2	H	200[B]	HEM	4	0
2	E	200[A]	HEM	12	0
2	K	200[B]	HEM	4	0
4	B	1160	SO4	1	0
2	K	200[A]	HEM	6	0
2	D	200[B]	HEM	5	0
4	J	1159	SO4	1	0
2	L	200[B]	HEM	5	0
2	D	200[A]	HEM	6	0
2	B	200[A]	HEM	3	0
2	B	200[B]	HEM	4	0
2	L	200[A]	HEM	6	0
2	J	200[A]	HEM	7	0
2	J	200[B]	HEM	3	0
2	A	200[B]	HEM	4	0
2	C	200[A]	HEM	5	0
2	A	200[A]	HEM	5	0
2	I	200[B]	HEM	5	0
2	C	200[B]	HEM	5	0
2	I	200[A]	HEM	8	0
2	G	200[A]	HEM	5	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.

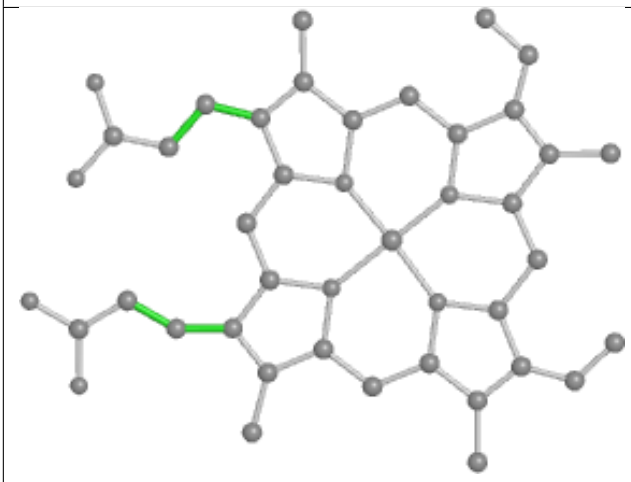
Ligand HEM G 200 (B)



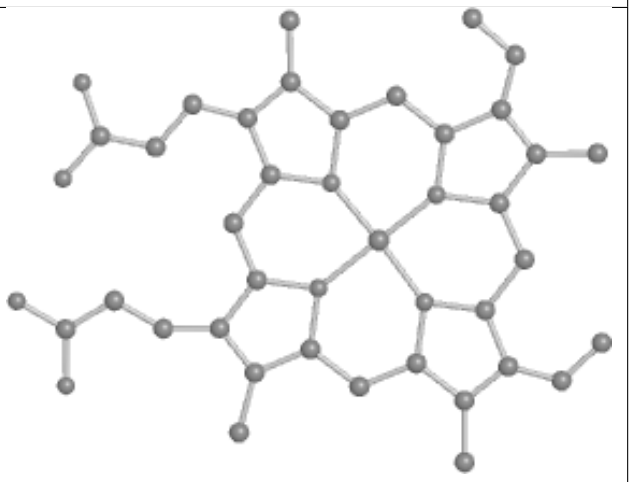
Bond lengths



Bond angles

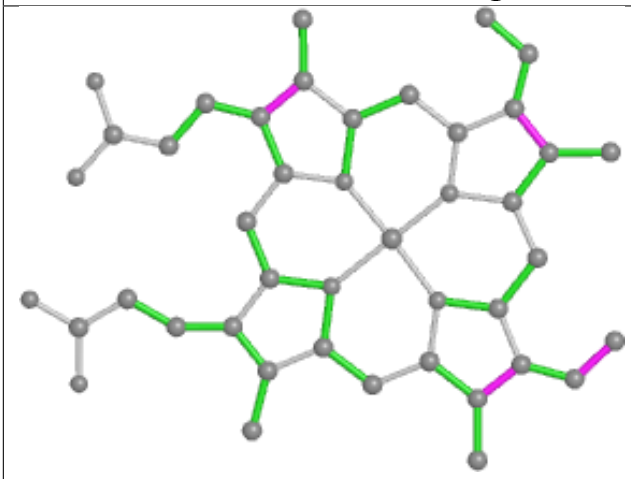


Torsions

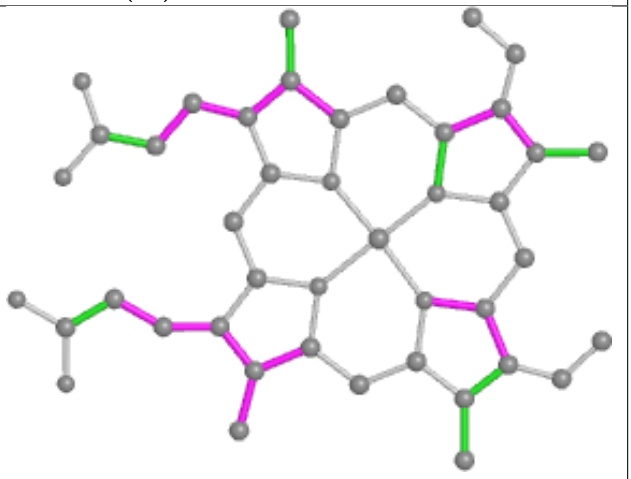


Rings

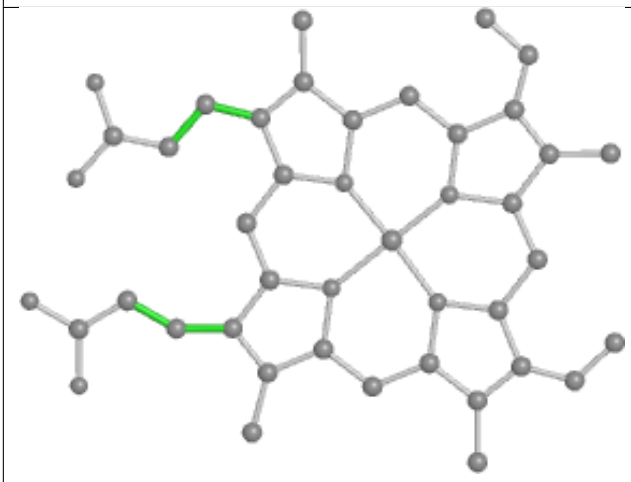
Ligand HEM F 200 (A)



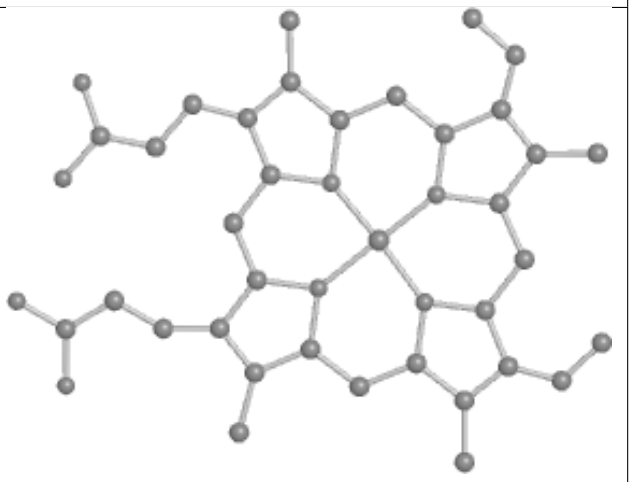
Bond lengths



Bond angles

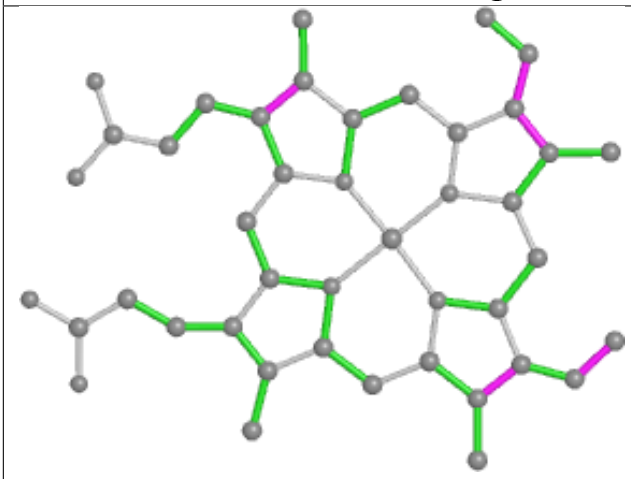


Torsions

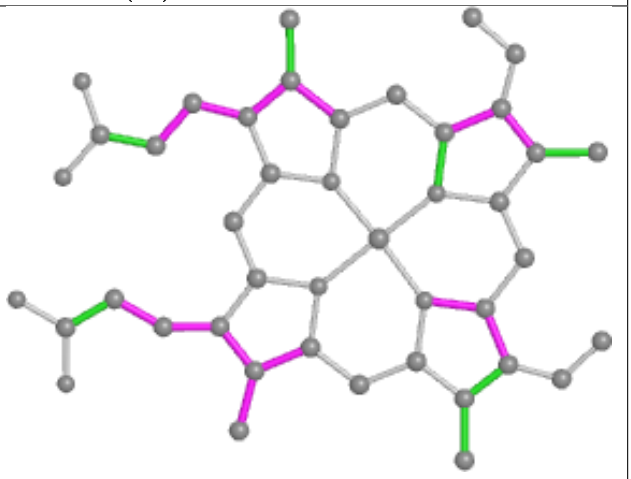


Rings

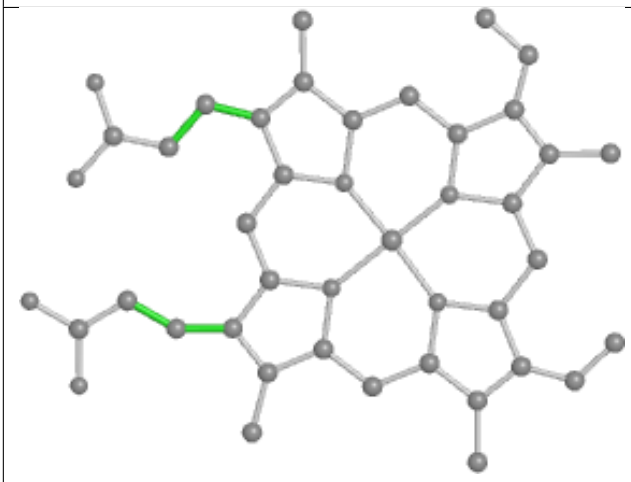
Ligand HEM F 200 (B)



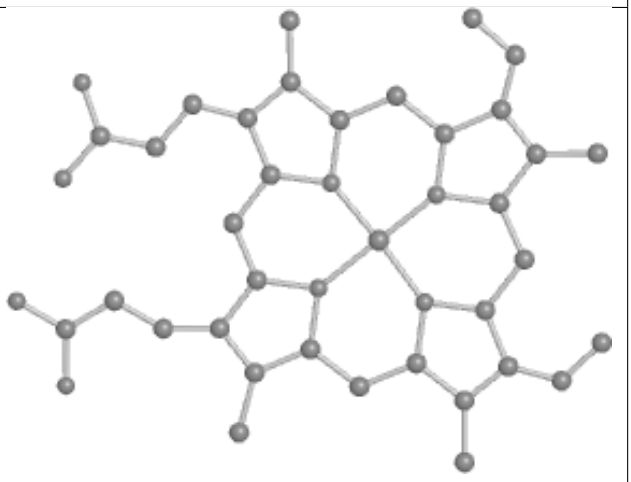
Bond lengths



Bond angles

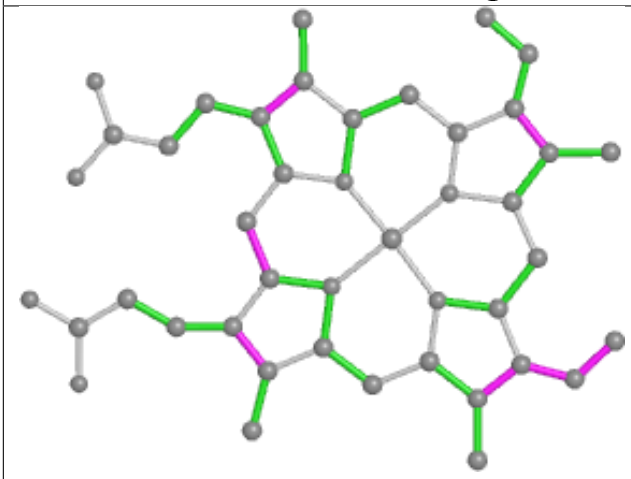


Torsions

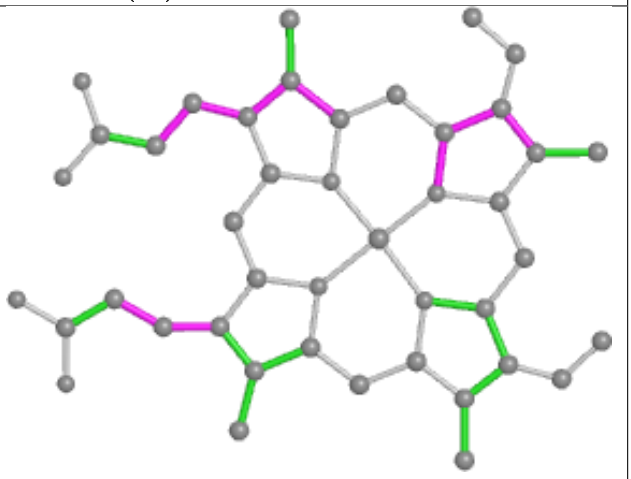


Rings

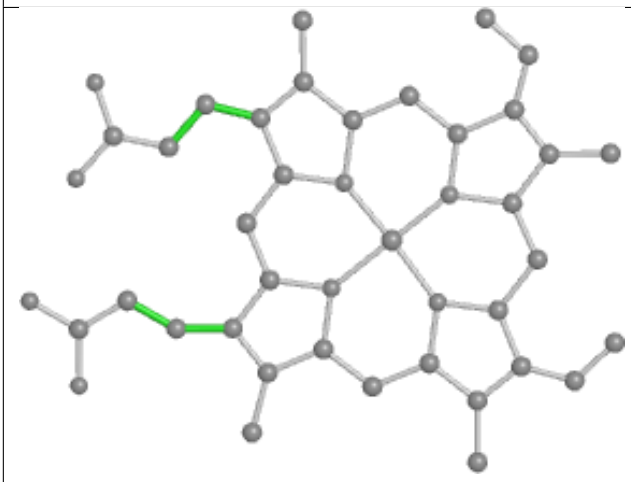
Ligand HEM H 200 (A)



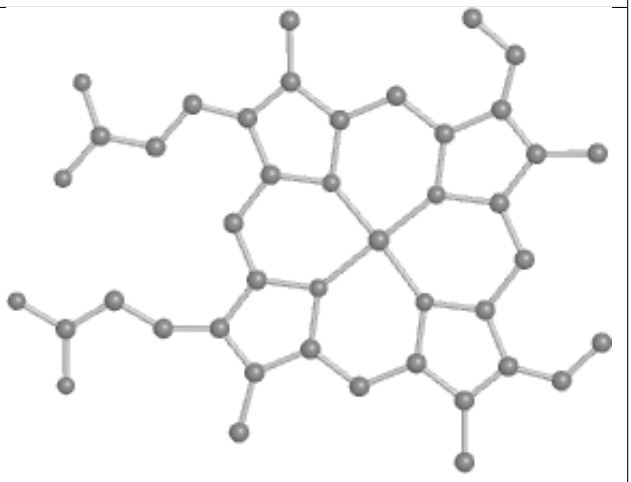
Bond lengths



Bond angles

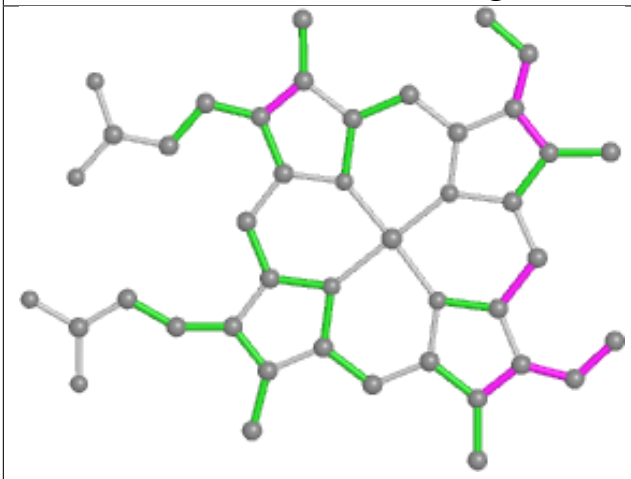


Torsions

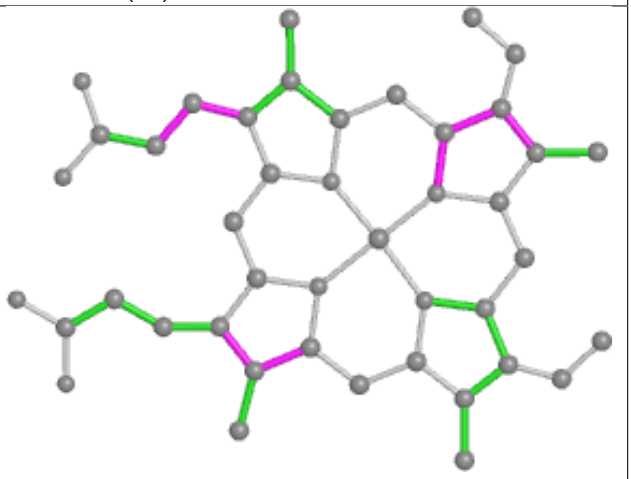


Rings

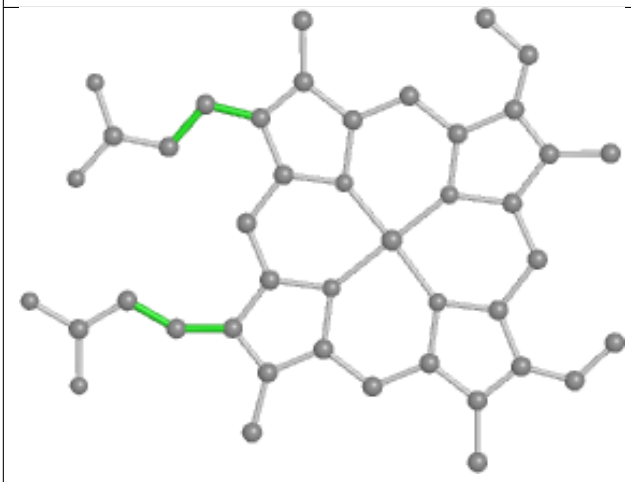
Ligand HEM E 200 (B)



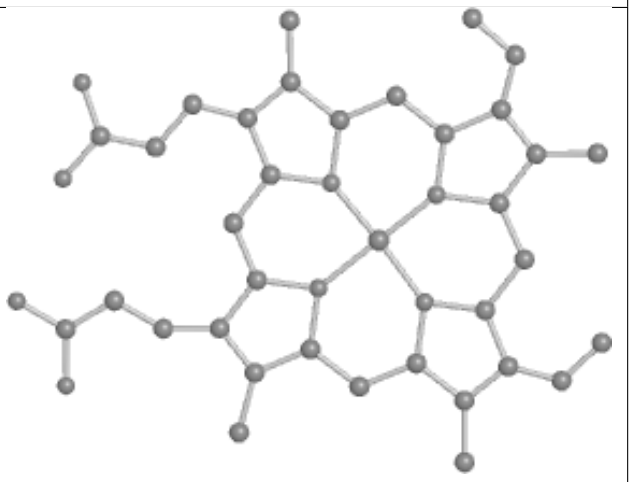
Bond lengths



Bond angles

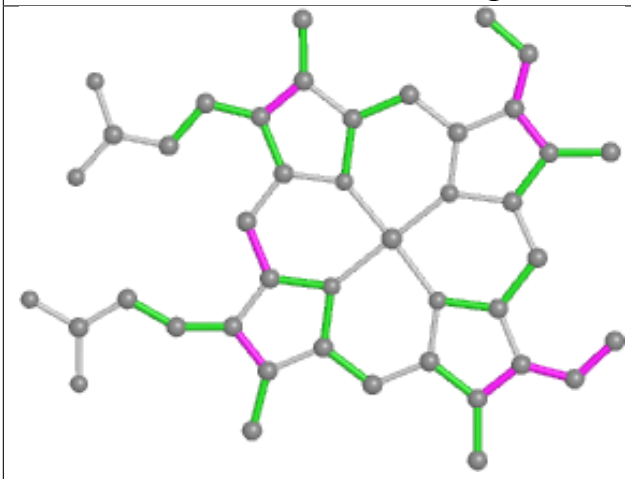


Torsions



Rings

Ligand HEM H 200 (B)



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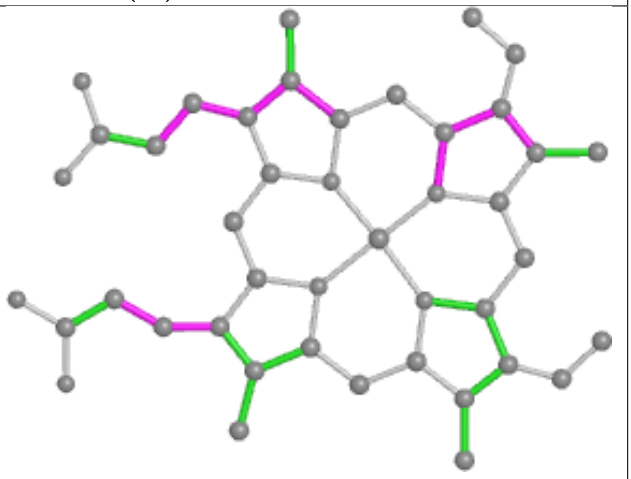
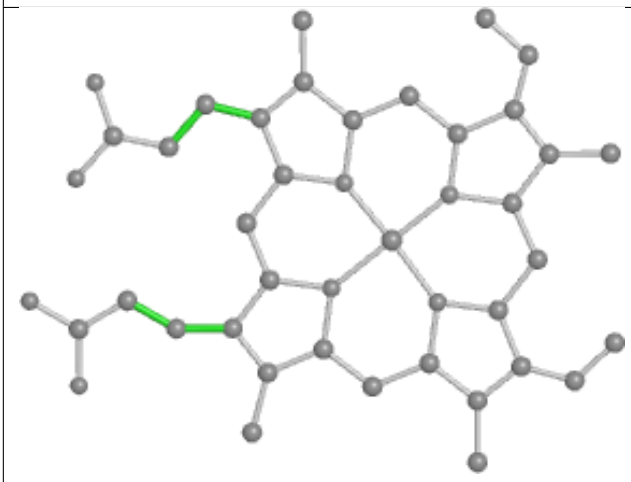


Bond angles



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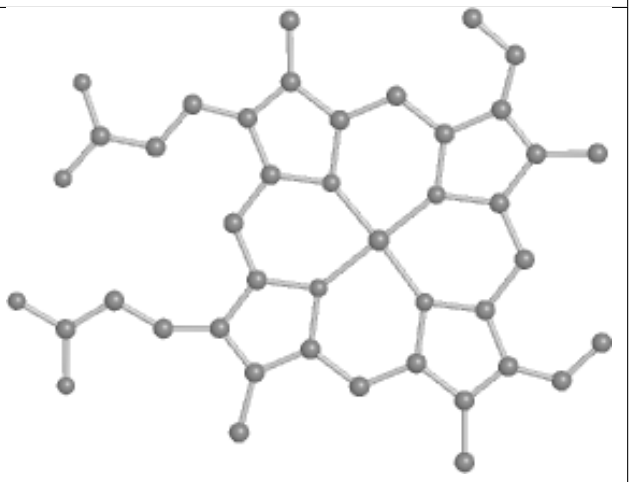
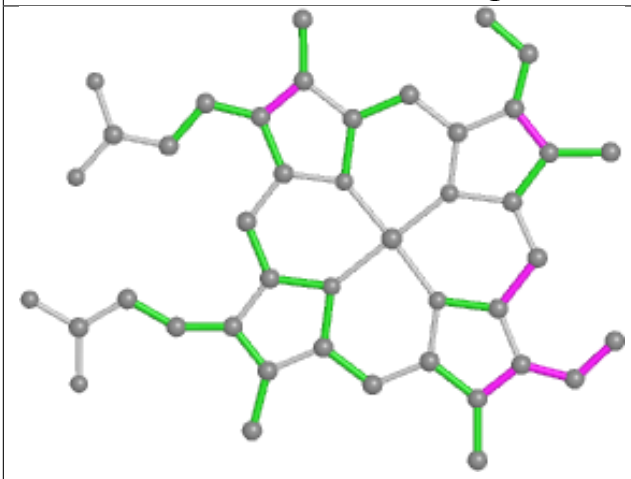


Rings

Ligand HEM E 200 (A)



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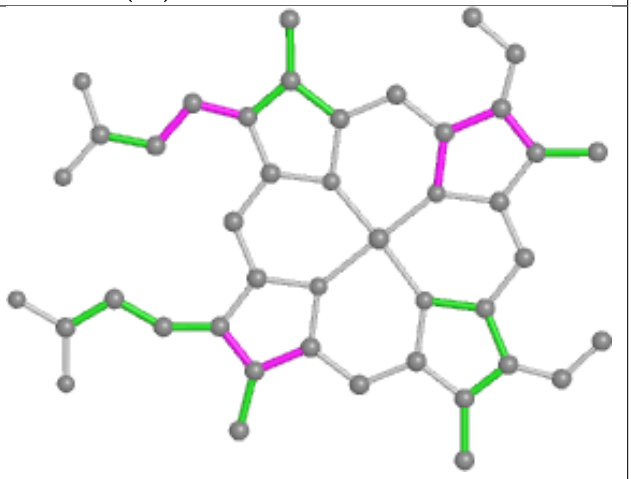
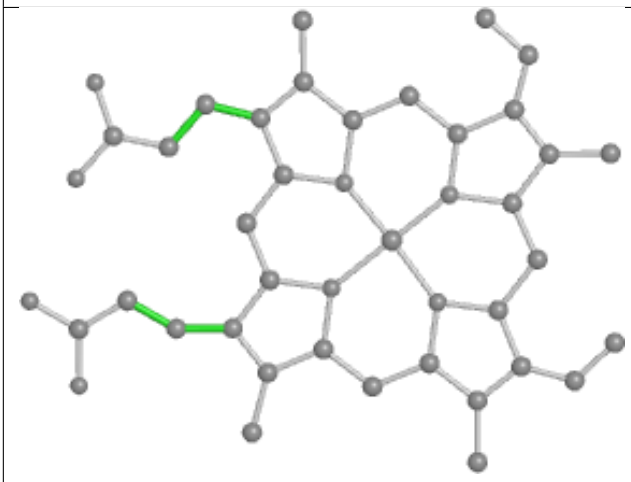


Bond angles



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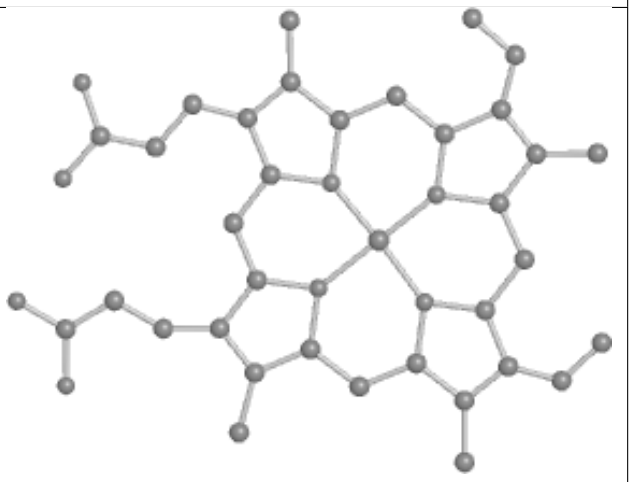
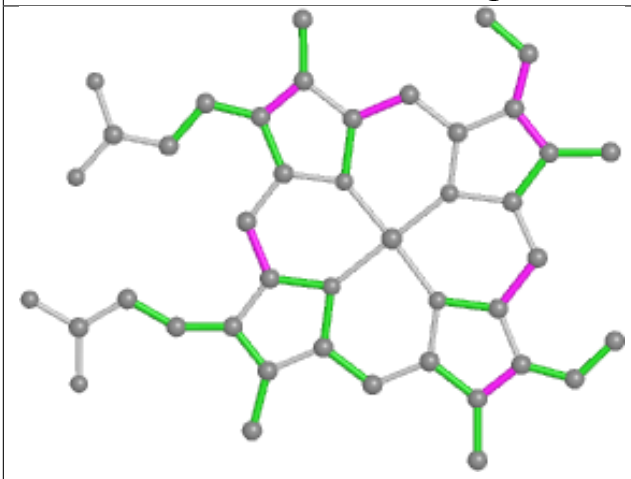


Rings

Ligand HEM K 200 (B)



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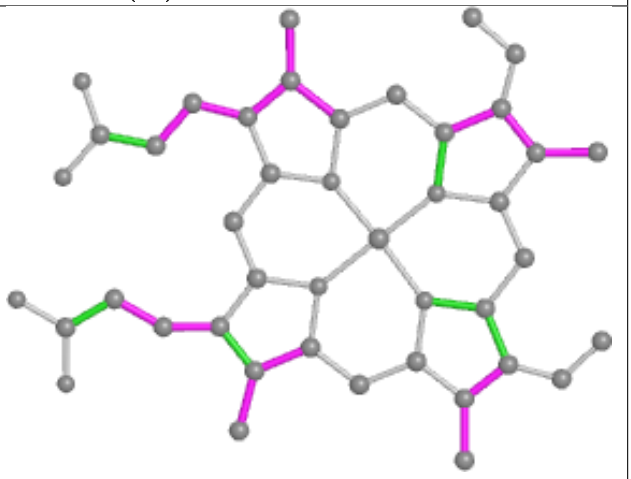
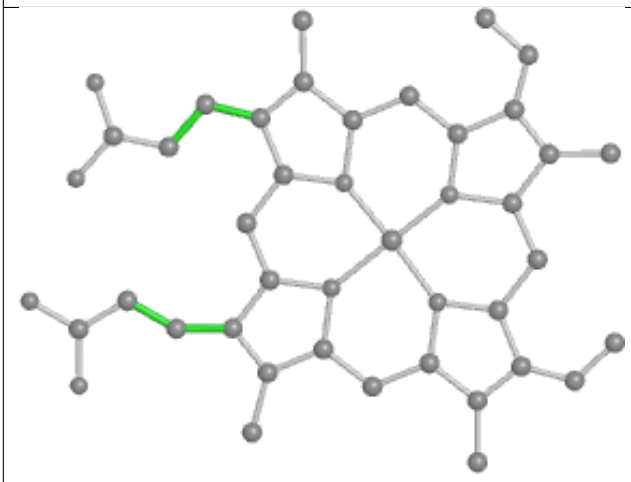
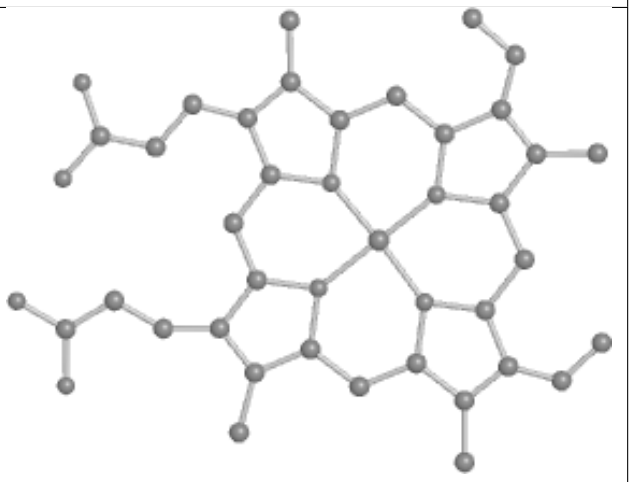


Bond angles

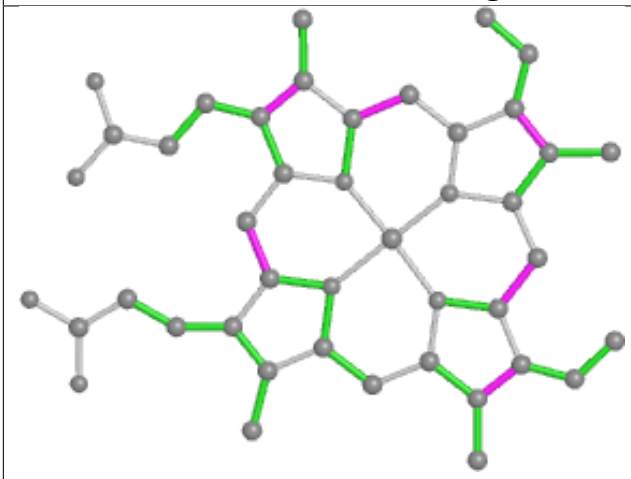


Torsions



Rings

Ligand HEM K 200 (A)



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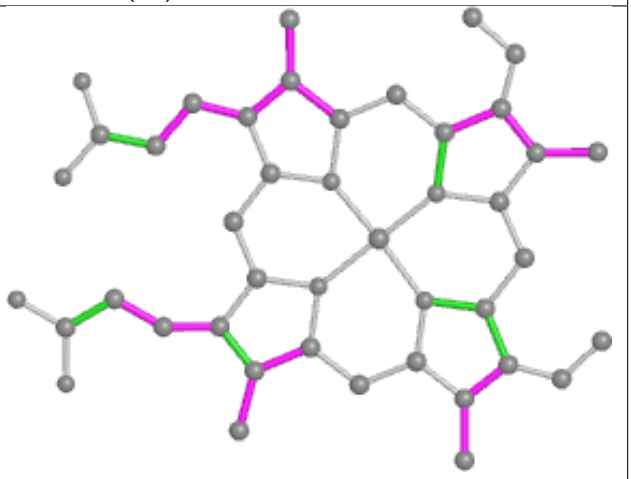
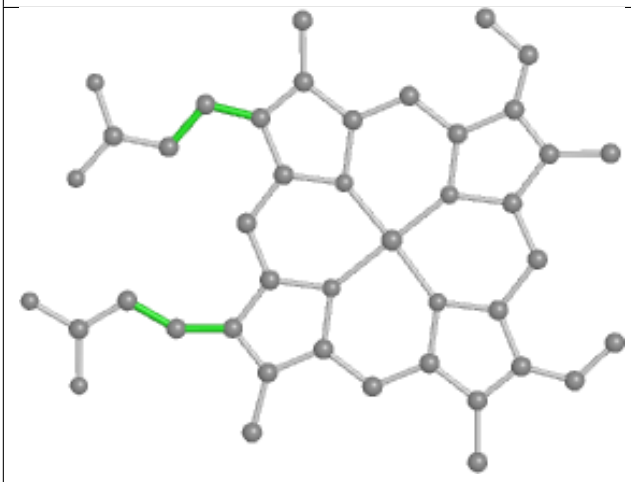


Bond angles



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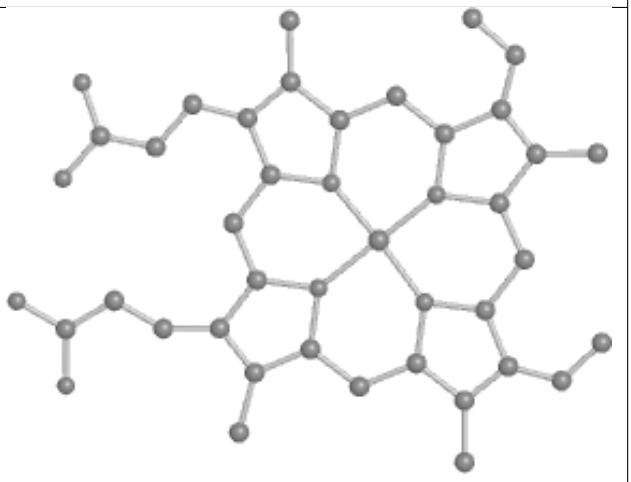
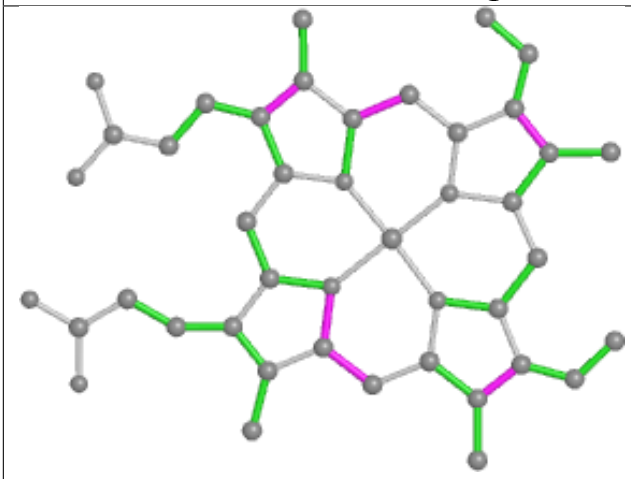


Rings

Ligand HEM D 200 (B)



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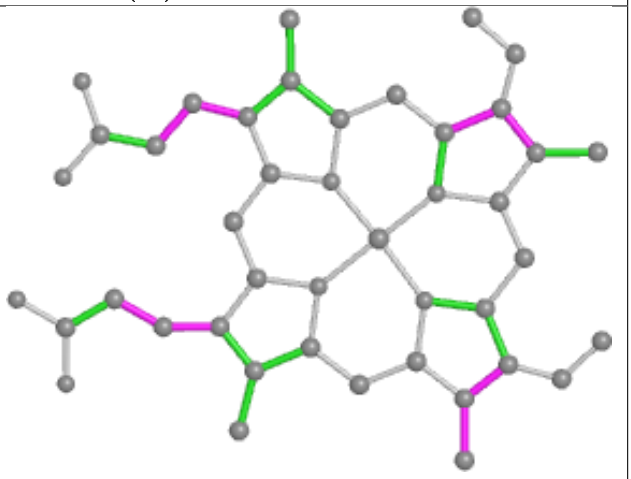
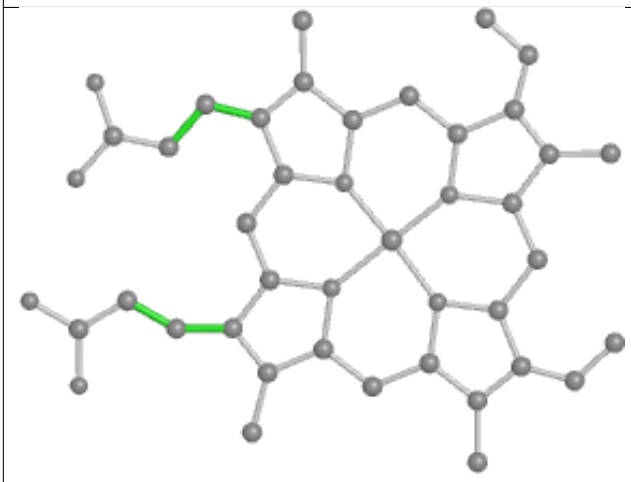


Bond angles



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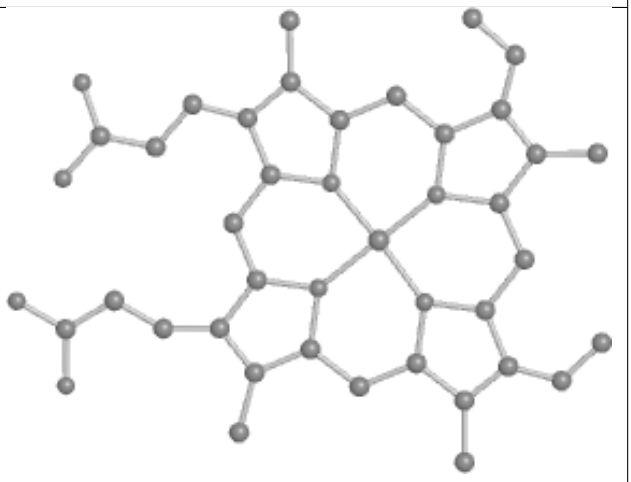
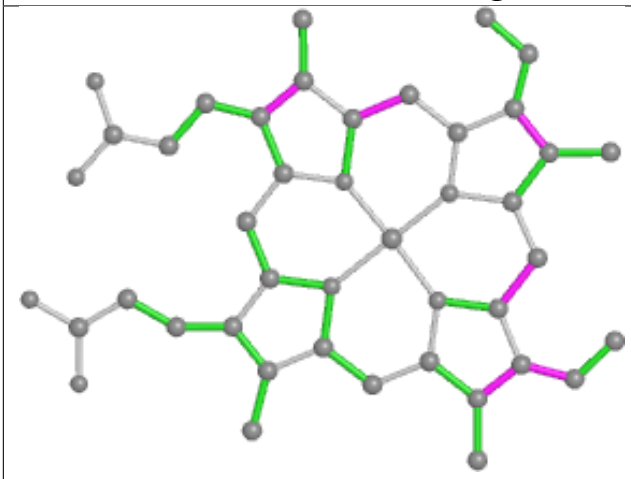


Rings

Ligand HEM L 200 (B)



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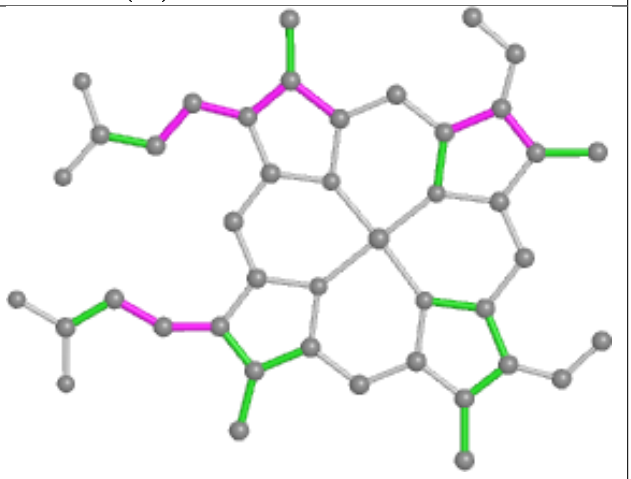
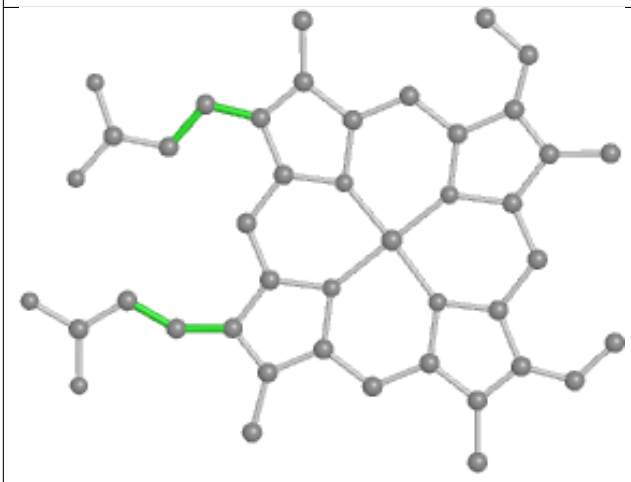
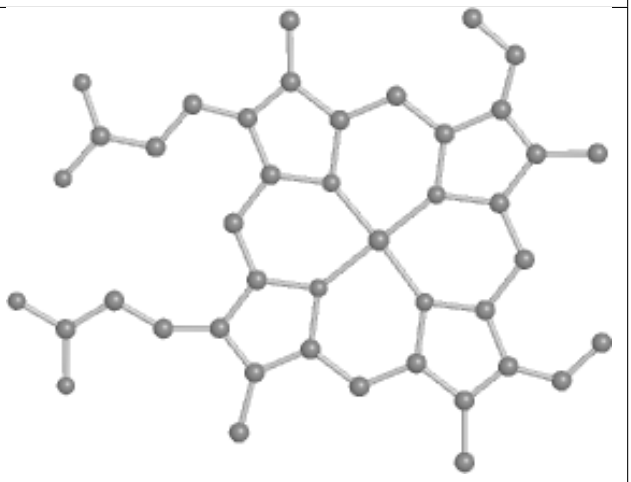


Bond angles

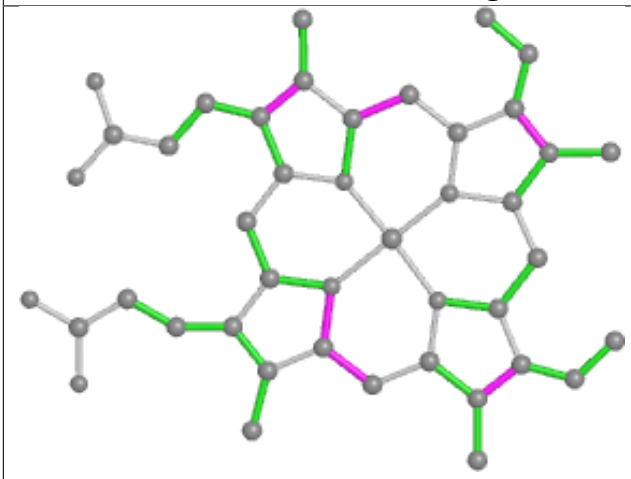


Torsions



Rings

Ligand HEM D 200 (A)



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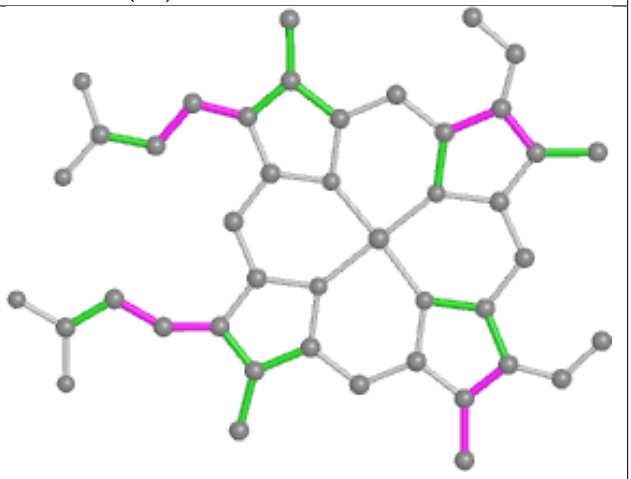
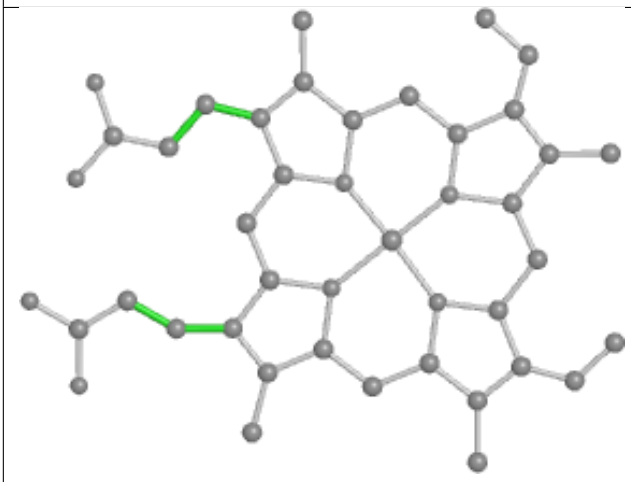


Bond angles



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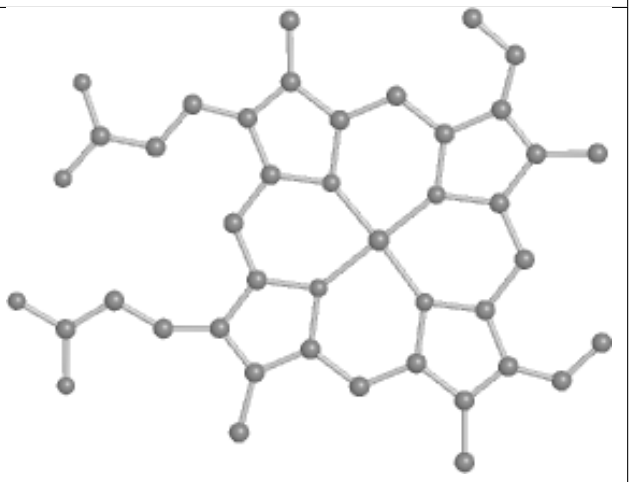
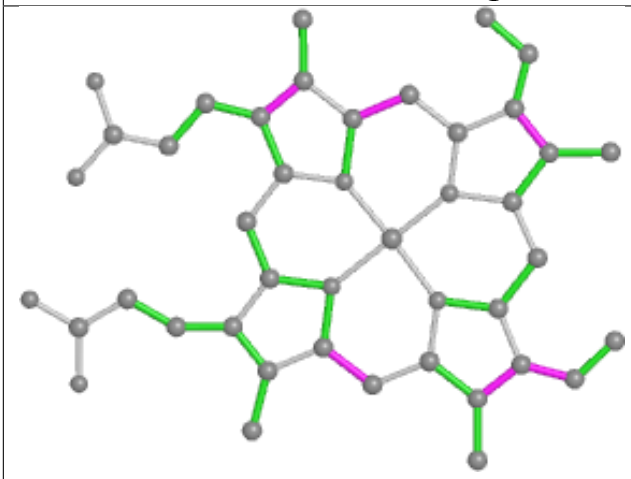


Rings

Ligand HEM B 200 (A)



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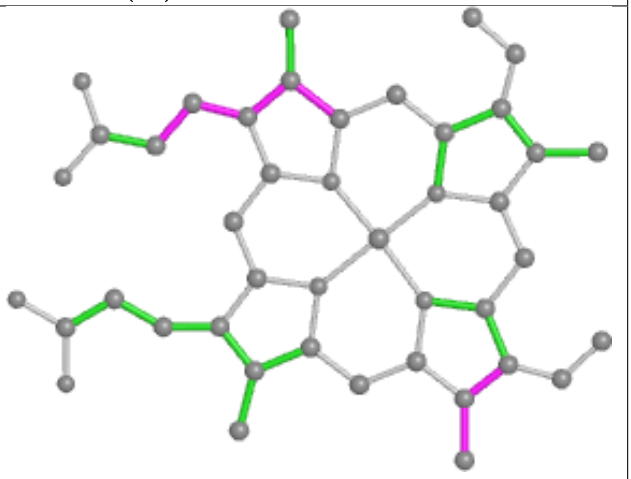
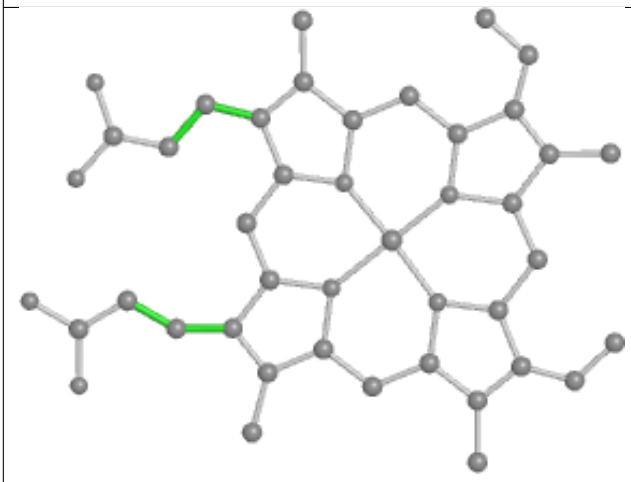
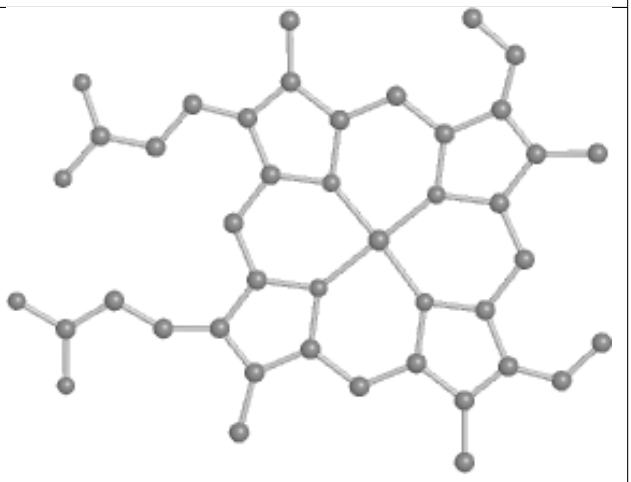


Bond angles

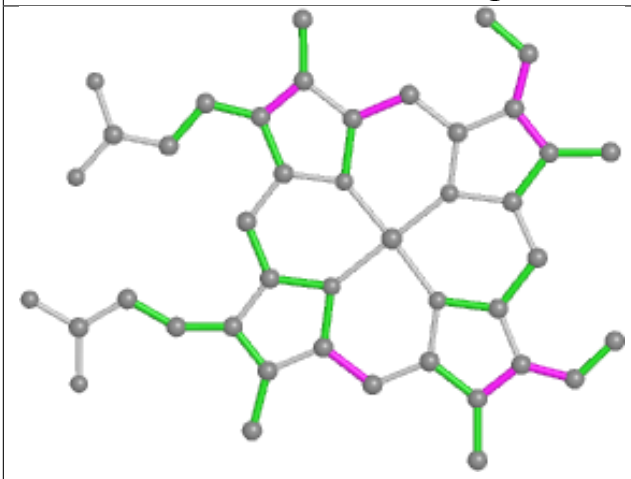


Torsions



Rings

Ligand HEM B 200 (B)



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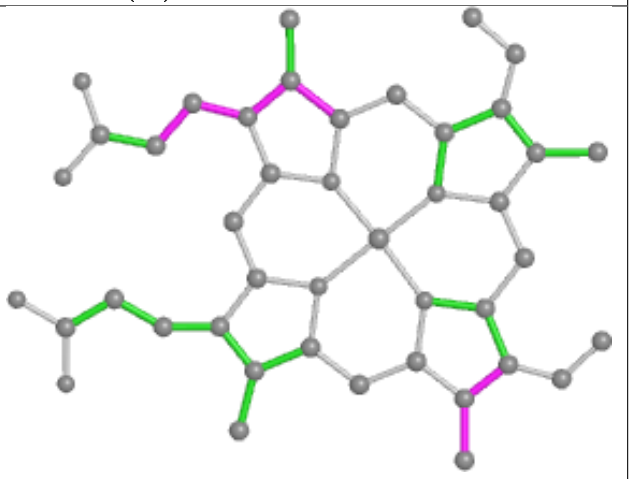
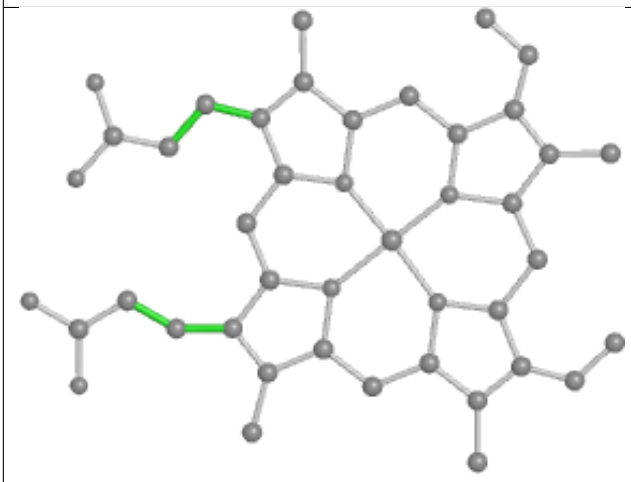
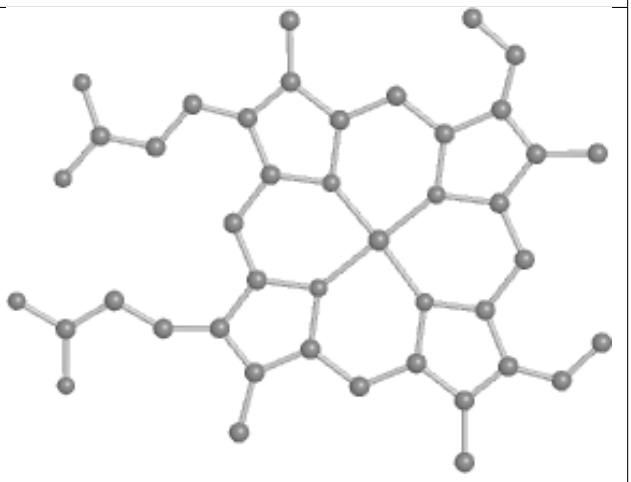


Bond angles

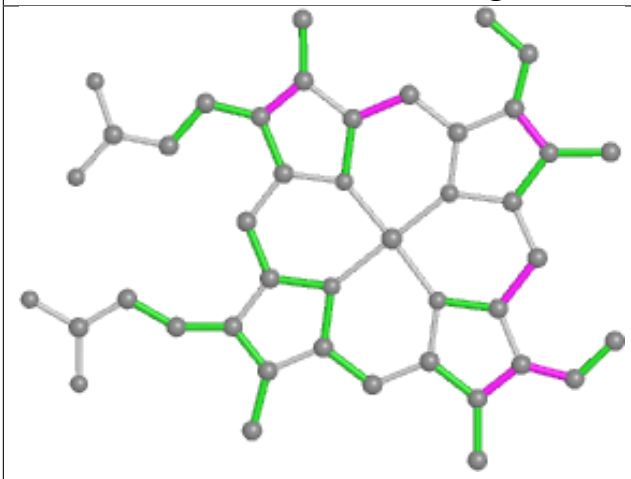


Torsions



Rings

Ligand HEM L 200 (A)



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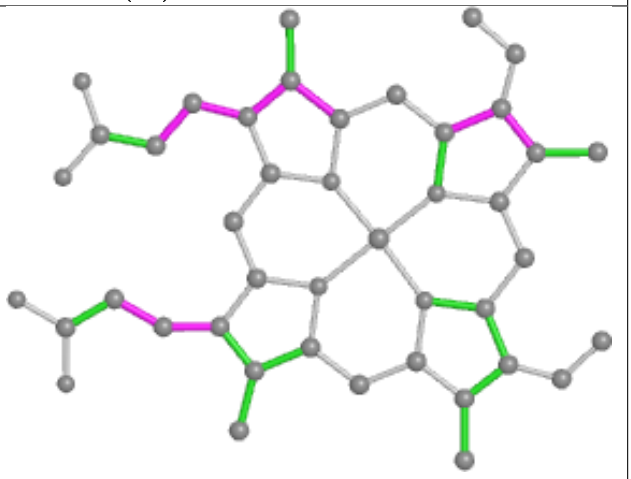
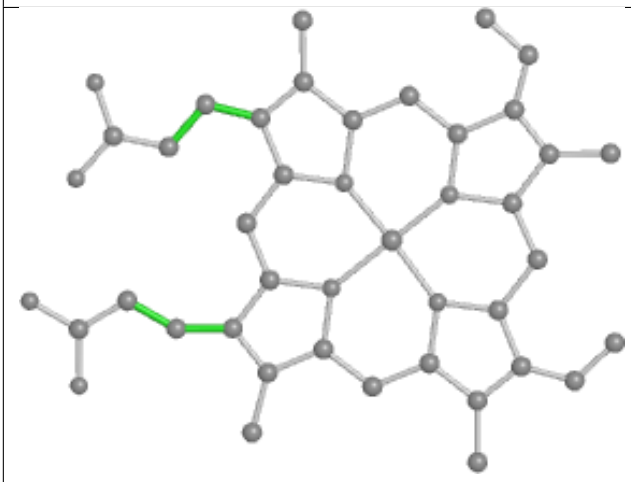
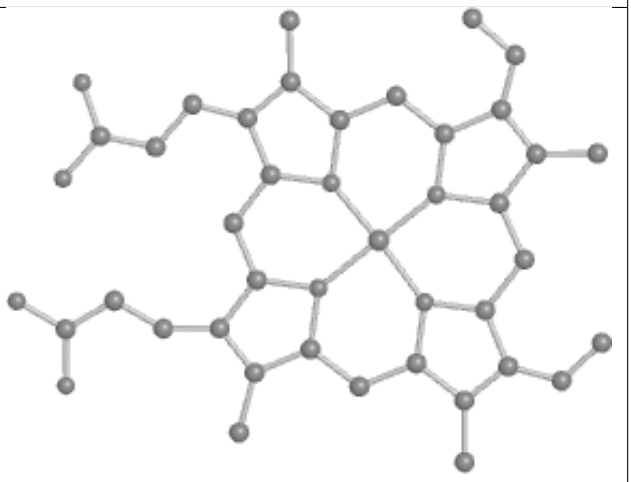


Bond angles

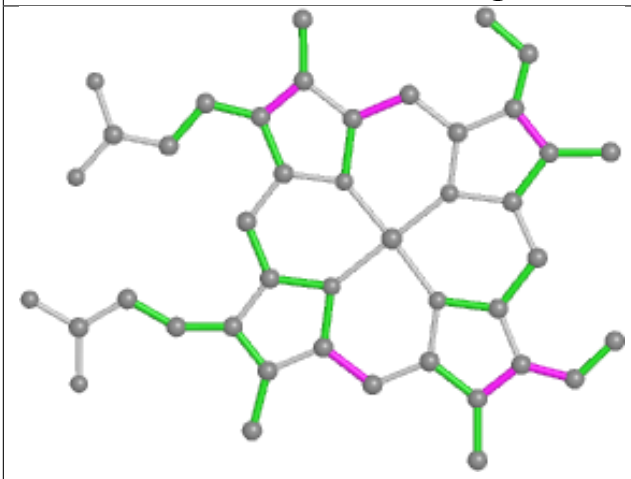


Torsions



Rings

Ligand HEM J 200 (A)



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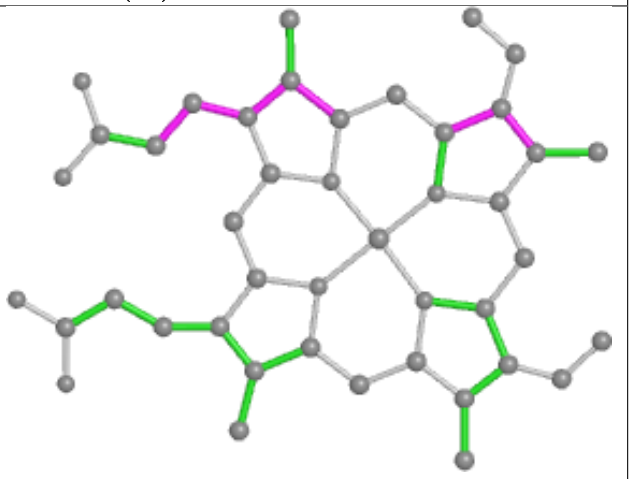
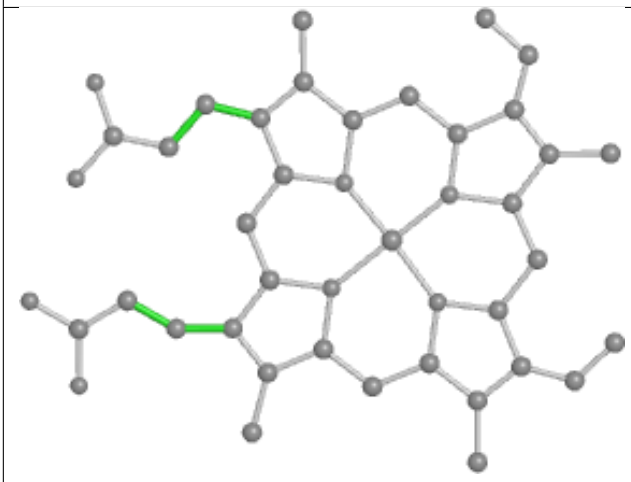
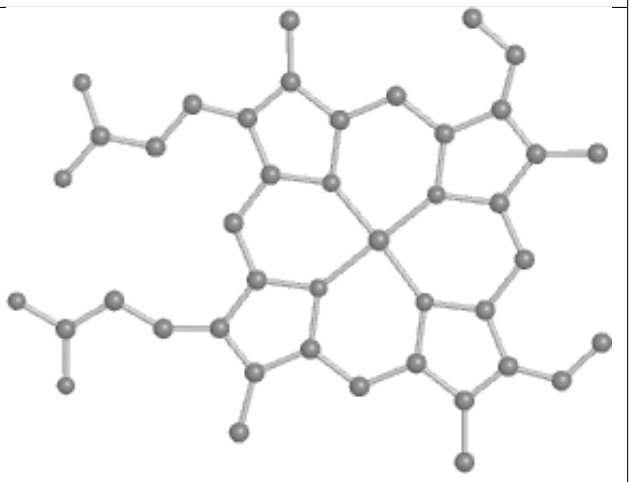


Bond angles

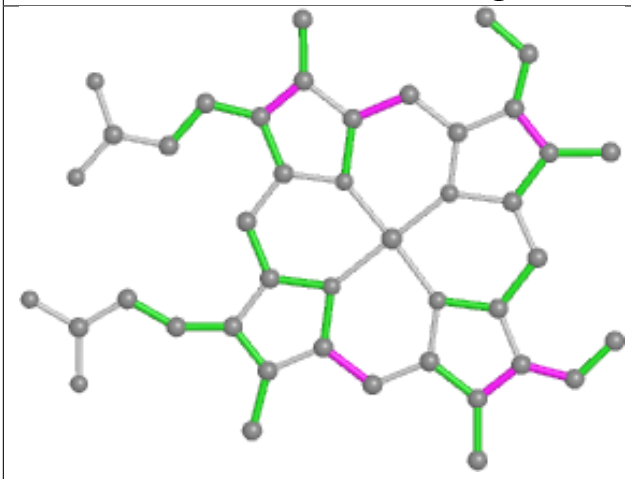


Torsions

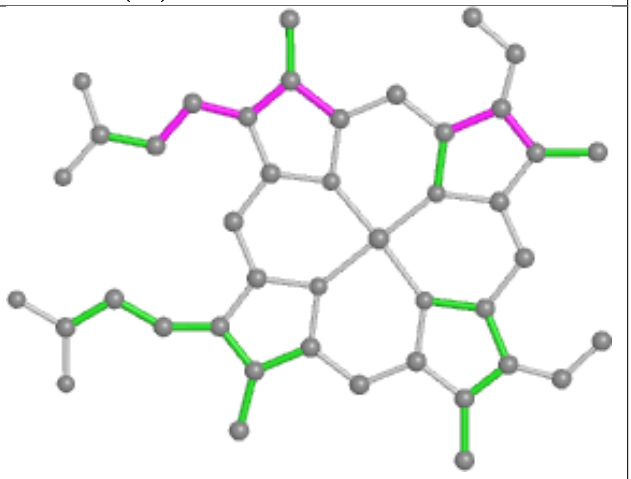


Rings

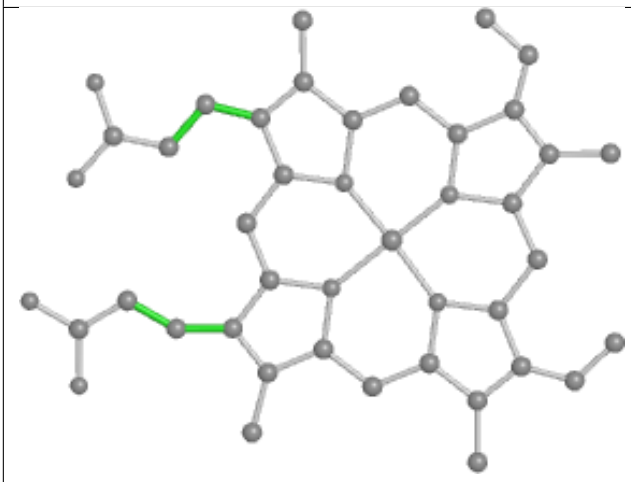
Ligand HEM J 200 (B)



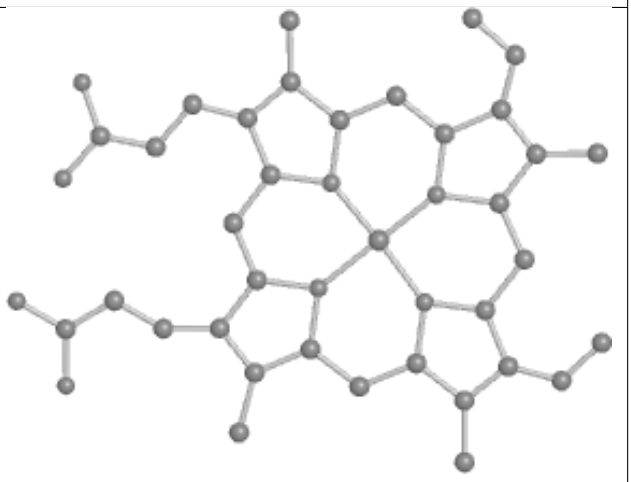
Bond lengths



Bond angles

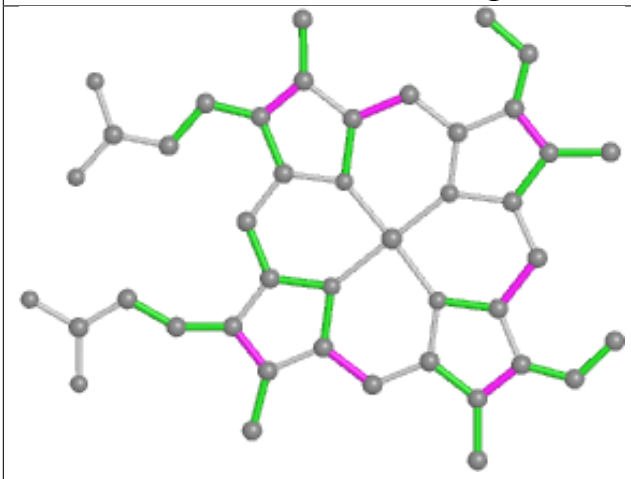


Torsions

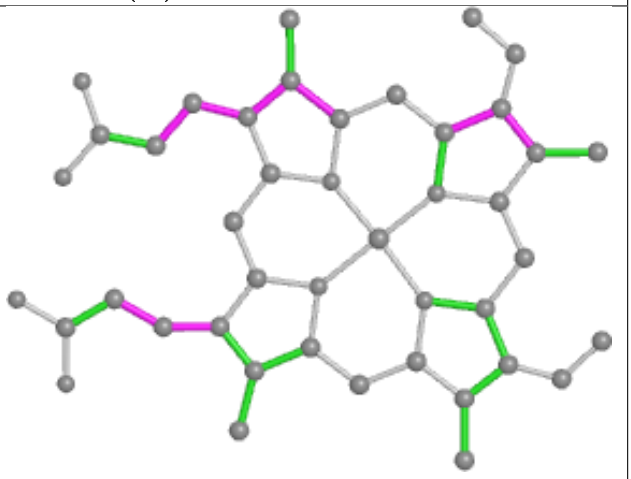


Rings

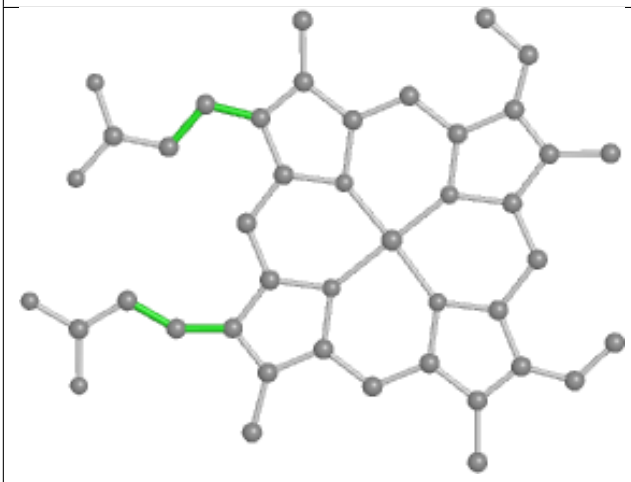
Ligand HEM A 200 (B)



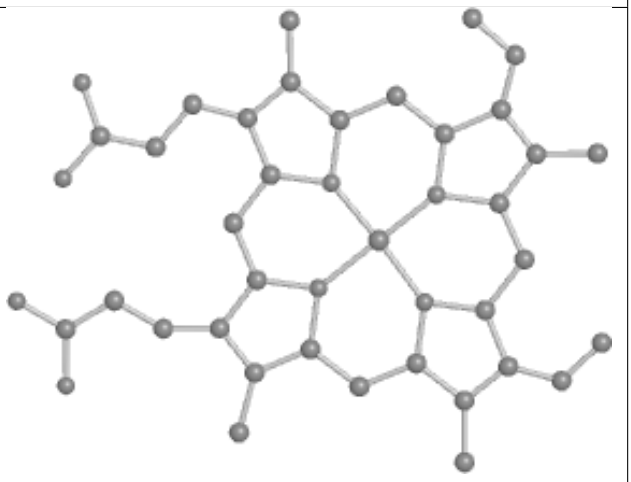
Bond lengths



Bond angles

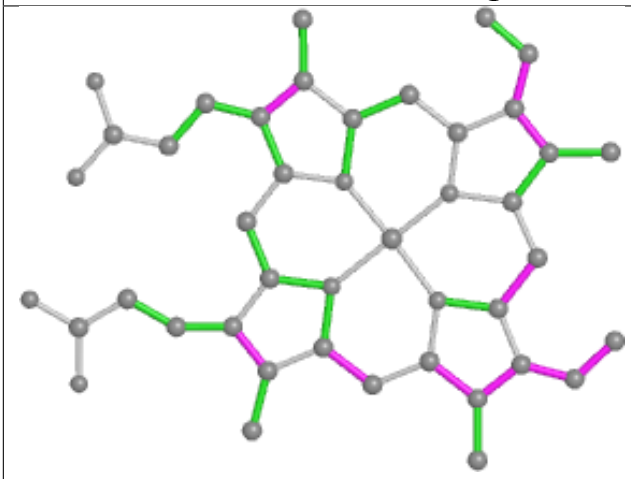


Torsions

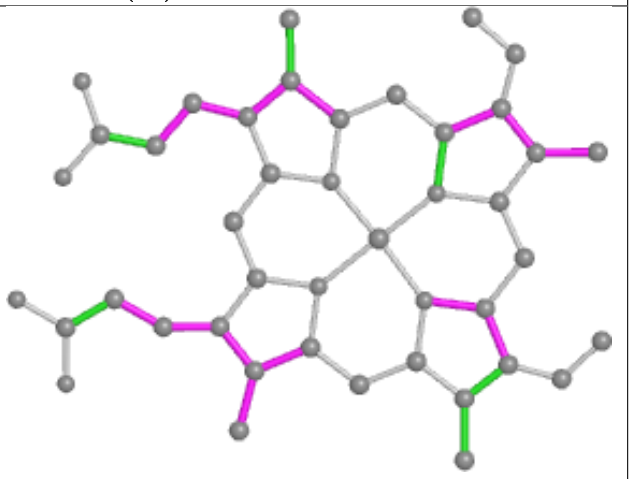


Rings

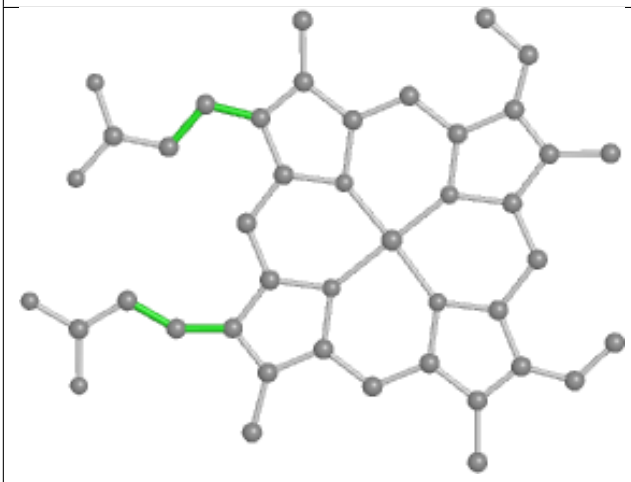
Ligand HEM C 200 (A)



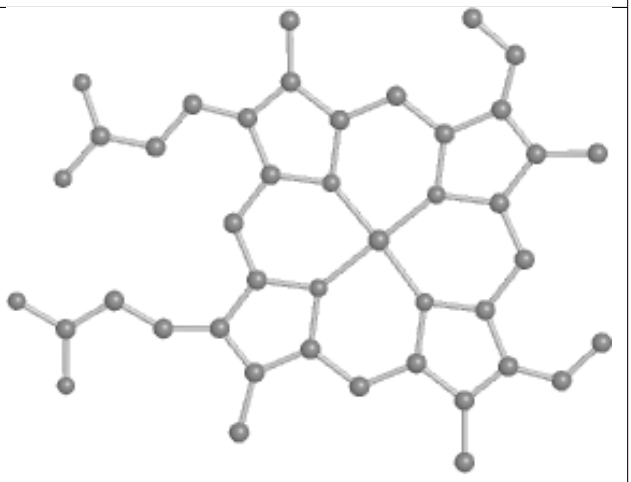
Bond lengths



Bond angles

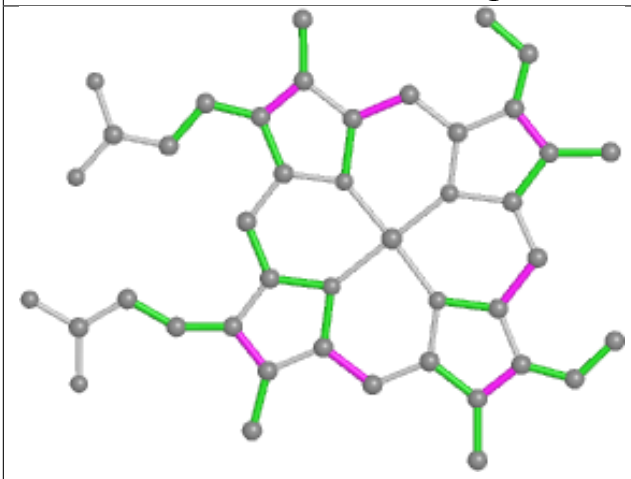


Torsions

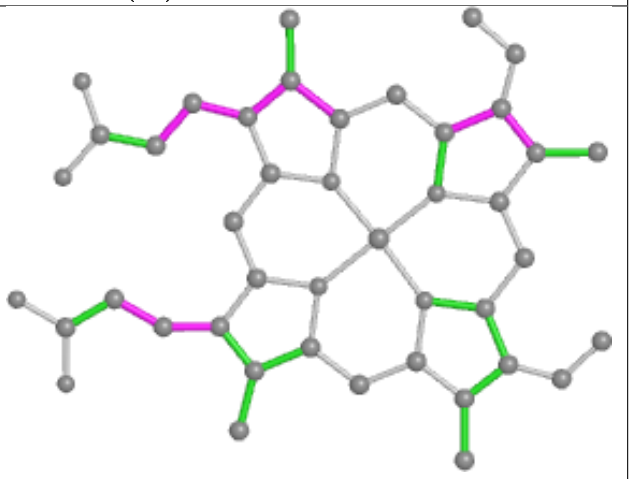


Rings

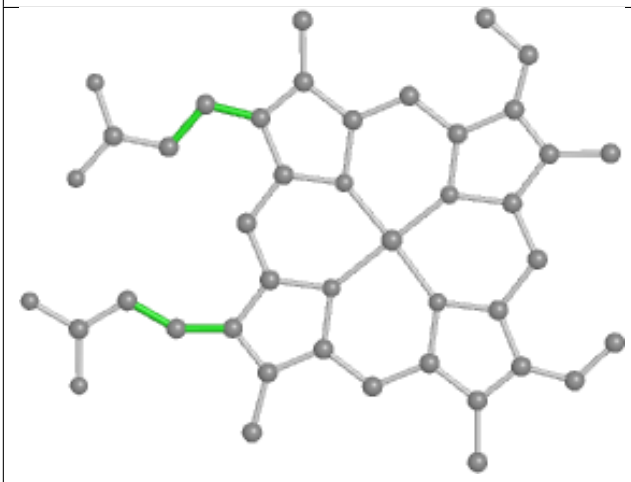
Ligand HEM A 200 (A)



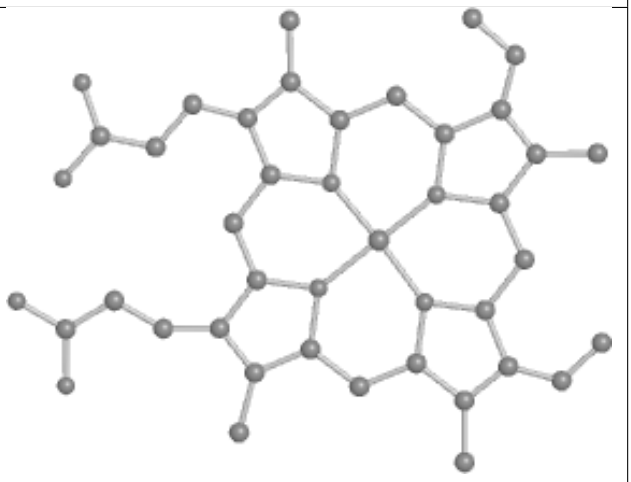
Bond lengths



Bond angles

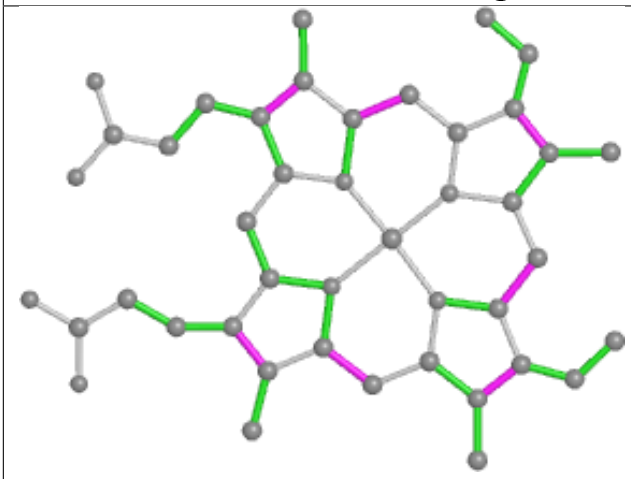


Torsions

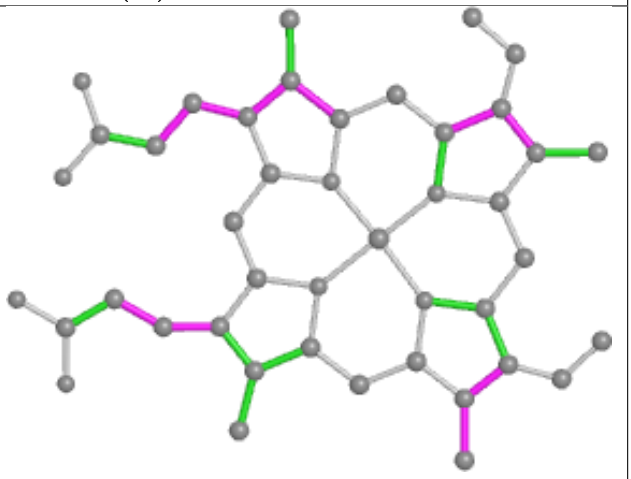


Rings

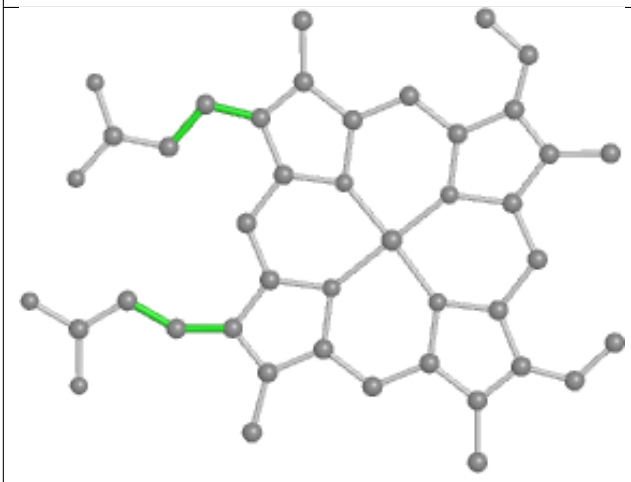
Ligand HEM I 200 (B)



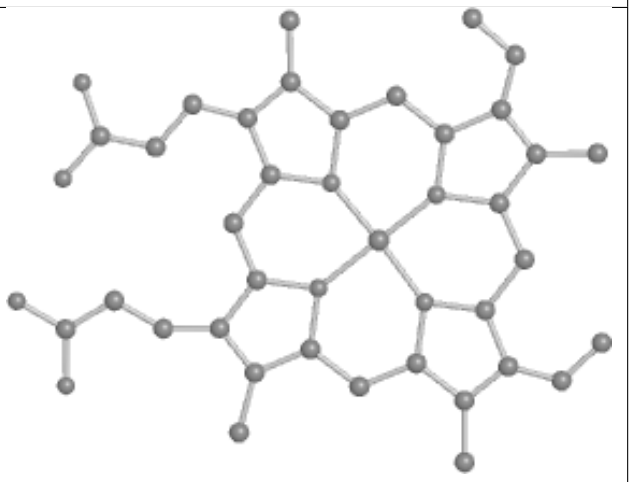
Bond lengths



Bond angles

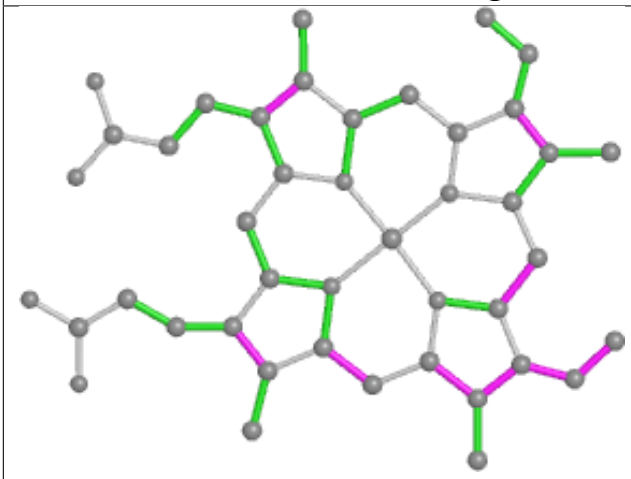


Torsions

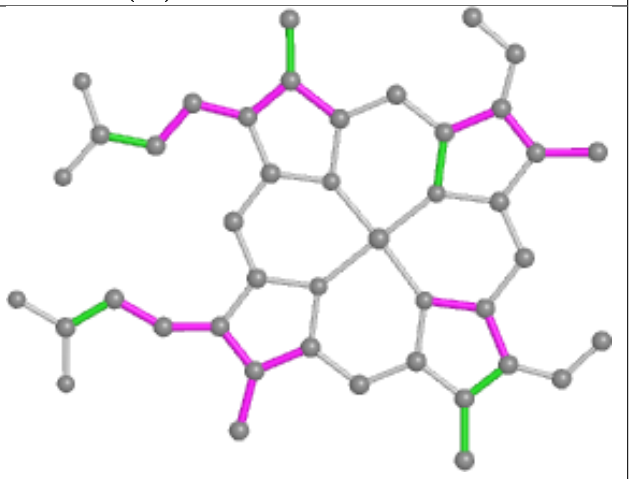


Rings

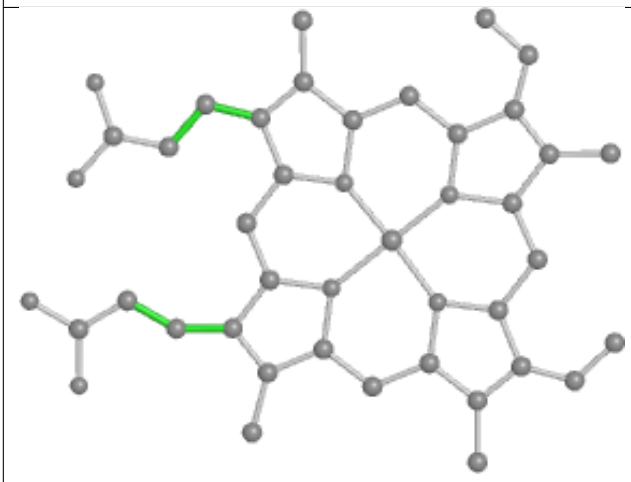
Ligand HEM C 200 (B)



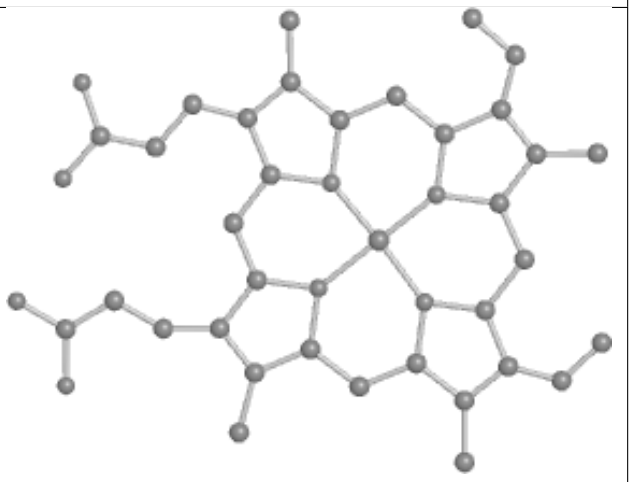
Bond lengths



Bond angles

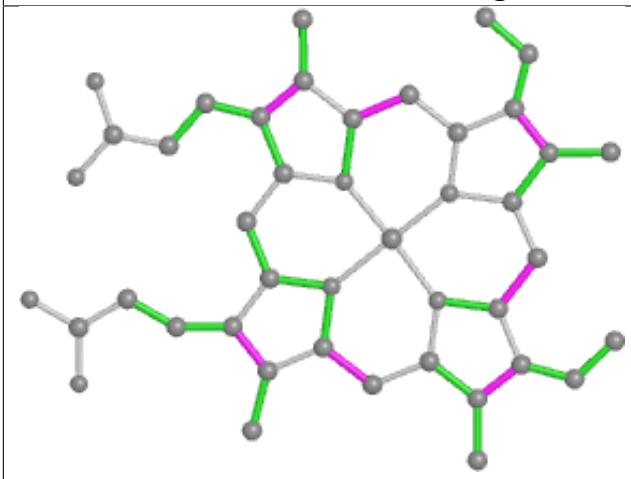


Torsions

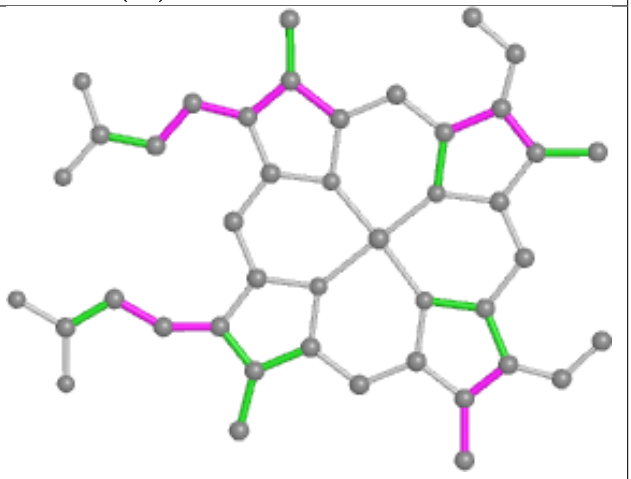


Rings

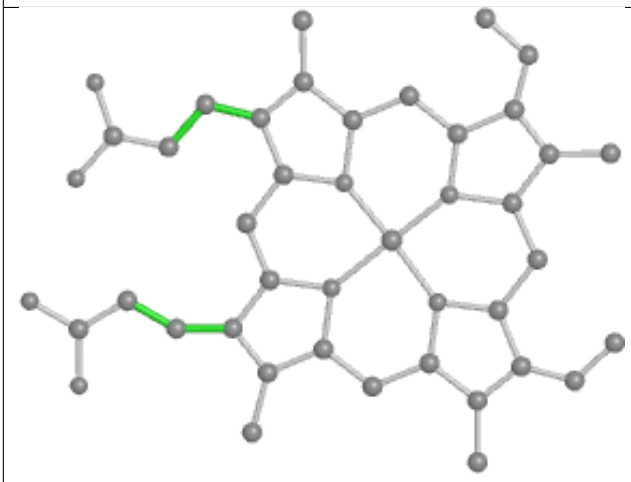
Ligand HEM I 200 (A)



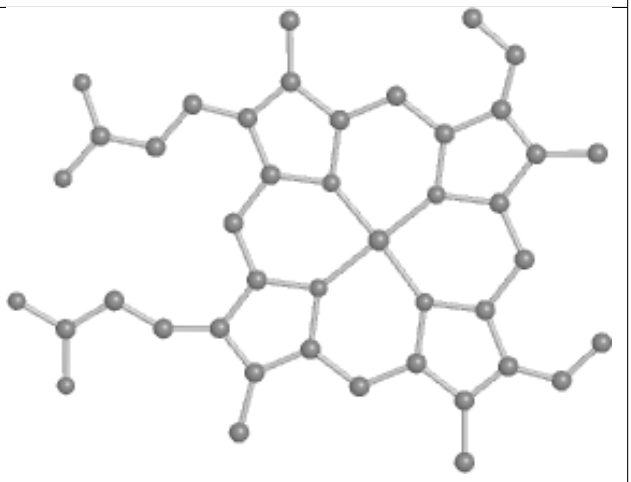
Bond lengths



Bond angles

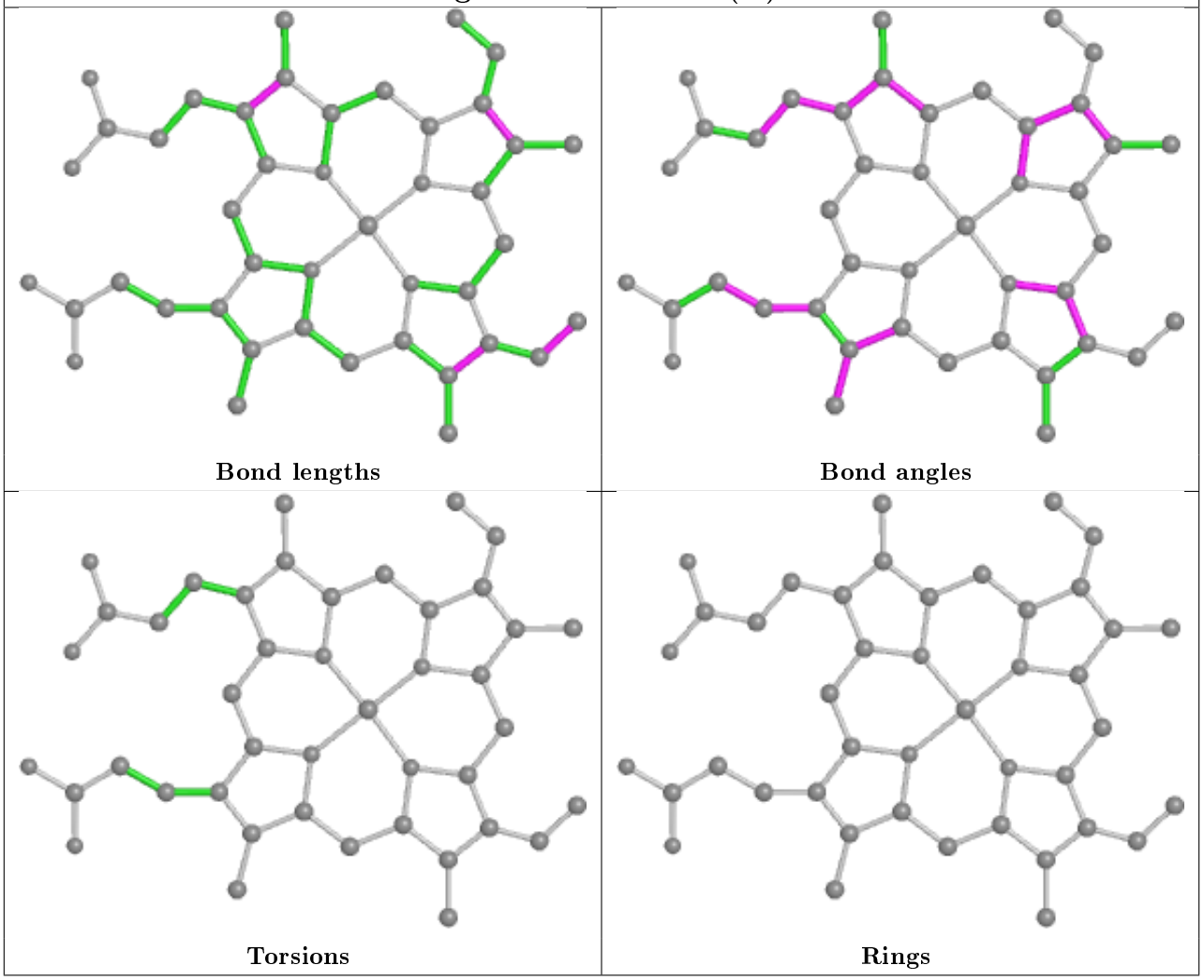


Torsions



Rings

Ligand HEM G 200 (A)



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

6.1 Protein, DNA and RNA chains

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, 95^{th} 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	157/158 (99%)	-0.21	4 (2%)	57	60	13, 20, 35, 70	1 (0%)
1	B	157/158 (99%)	-0.33	4 (2%)	57	60	13, 20, 37, 69	0
1	C	157/158 (99%)	-0.29	2 (1%)	77	79	13, 20, 36, 70	0
1	D	157/158 (99%)	-0.26	3 (1%)	66	69	13, 21, 36, 72	0
1	E	157/158 (99%)	-0.31	2 (1%)	77	79	13, 20, 35, 70	0
1	F	157/158 (99%)	-0.27	3 (1%)	66	69	13, 20, 35, 67	0
1	G	157/158 (99%)	-0.31	1 (0%)	89	90	13, 20, 36, 69	0
1	H	157/158 (99%)	-0.35	2 (1%)	77	79	12, 20, 36, 73	0
1	I	157/158 (99%)	-0.32	3 (1%)	66	69	12, 20, 36, 67	0
1	J	157/158 (99%)	-0.28	4 (2%)	57	60	13, 20, 36, 70	0
1	K	157/158 (99%)	-0.30	2 (1%)	77	79	13, 20, 36, 70	0
1	L	157/158 (99%)	-0.35	2 (1%)	77	79	12, 20, 36, 64	0
All	All	1884/1896 (99%)	-0.30	32 (1%)	70	72	12, 20, 36, 73	1 (0%)

All (32) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	D	157	GLU	3.8
1	H	157	GLU	3.6
1	F	157	GLU	3.5
1	G	157	GLU	3.4
1	I	157	GLU	3.4
1	C	157	GLU	3.4
1	I	156	GLU	3.3
1	K	157	GLU	3.2
1	E	157	GLU	3.1
1	B	157	GLU	3.0
1	K	156	GLU	2.9

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Mol	Chain	Res	Type	RSRZ
1	A	73	ASP	2.9
1	D	156	GLU	2.8
1	E	156	GLU	2.8
1	L	157	GLU	2.7
1	C	2	LYS	2.7
1	A	157	GLU	2.7
1	I	2	LYS	2.6
1	J	157	GLU	2.6
1	J	96	ASP	2.5
1	F	156	GLU	2.4
1	L	156	GLU	2.4
1	A	112	HIS	2.3
1	A	156	GLU	2.3
1	B	156	GLU	2.3
1	H	2	LYS	2.2
1	B	1	MET	2.2
1	J	1	MET	2.1
1	B	2	LYS	2.1
1	D	73	ASP	2.1
1	J	156	GLU	2.1
1	F	73	ASP	2.0

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 carbohydrates 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, 95^{th} 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
4	SO4	K	1158	5/5	0.74	0.29	84,89,92,94	0
4	SO4	G	1158	5/5	0.80	0.24	59,66,70,79	0

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	SO4	I	1158	5/5	0.85	0.27	52,65,73,75	0
4	SO4	J	1158	5/5	0.86	0.28	65,76,80,81	0
4	SO4	B	1160	5/5	0.88	0.28	63,72,78,81	0
4	SO4	B	1159	5/5	0.88	0.25	60,62,70,70	0
4	SO4	A	1158	5/5	0.89	0.27	62,73,76,78	0
4	SO4	A	1159	5/5	0.90	0.32	66,68,74,75	0
4	SO4	H	1160	5/5	0.90	0.23	47,48,54,59	0
4	SO4	J	1159	5/5	0.90	0.30	67,67,73,76	0
4	SO4	A	1160	5/5	0.90	0.32	95,96,99,105	0
4	SO4	F	1158	5/5	0.90	0.26	40,62,67,70	0
4	SO4	K	1159	5/5	0.90	0.28	58,64,69,69	0
4	SO4	D	1158	5/5	0.91	0.23	73,75,80,80	0
4	SO4	B	1161	5/5	0.91	0.28	67,73,77,81	0
4	SO4	L	1158	5/5	0.91	0.23	48,61,68,69	0
2	HEM	A	200[B]	43/43	0.92	0.14	2,15,29,37	43
2	HEM	A	200[A]	43/43	0.92	0.14	2,15,29,37	43
4	SO4	C	1160	5/5	0.92	0.20	61,63,71,72	0
2	HEM	B	200[A]	43/43	0.92	0.14	8,15,26,34	43
2	HEM	B	200[B]	43/43	0.92	0.14	8,15,26,34	43
4	SO4	L	1159	5/5	0.93	0.34	64,67,72,75	0
2	HEM	L	200[B]	43/43	0.93	0.12	6,14,29,33	43
4	SO4	E	1160	5/5	0.93	0.23	45,51,61,61	0
4	SO4	G	1159	5/5	0.93	0.31	58,61,66,73	0
2	HEM	L	200[A]	43/43	0.93	0.12	6,14,29,33	43
4	SO4	C	1159	5/5	0.93	0.20	51,57,61,62	0
2	HEM	D	200[A]	43/43	0.94	0.12	9,17,34,39	43
2	HEM	E	200[A]	43/43	0.94	0.11	6,13,30,36	43
2	HEM	K	200[A]	43/43	0.94	0.12	9,16,28,36	43
2	HEM	H	200[A]	43/43	0.94	0.12	6,12,31,36	43
2	HEM	K	200[B]	43/43	0.94	0.12	9,16,28,36	43
2	HEM	J	200[A]	43/43	0.94	0.12	4,14,31,36	43
4	SO4	E	1159	5/5	0.94	0.29	46,62,65,65	0
2	HEM	J	200[B]	43/43	0.94	0.12	4,14,31,36	43
4	SO4	I	1159	5/5	0.94	0.27	67,72,78,81	0
2	HEM	F	200[B]	43/43	0.94	0.12	5,16,26,33	43
4	SO4	H	1159	5/5	0.94	0.28	54,60,62,63	0
2	HEM	G	200[B]	43/43	0.94	0.13	10,16,30,36	43
2	HEM	C	200[A]	43/43	0.94	0.12	8,13,24,31	43
2	HEM	D	200[B]	43/43	0.94	0.12	9,17,34,39	43
2	HEM	I	200[B]	43/43	0.94	0.12	9,16,28,32	43
2	HEM	H	200[B]	43/43	0.94	0.12	6,12,31,36	43
2	HEM	C	200[B]	43/43	0.94	0.12	8,13,24,31	43

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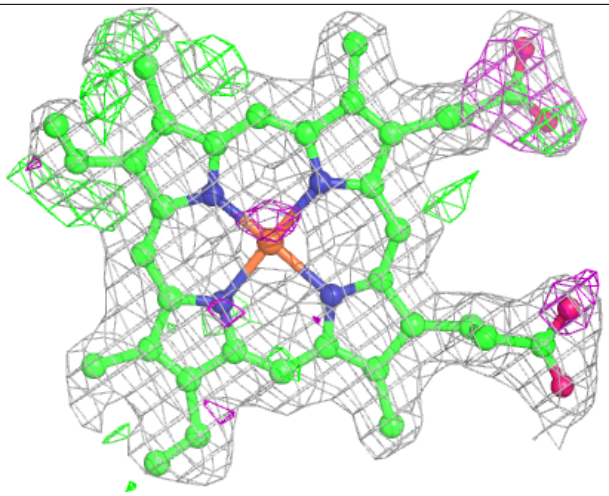
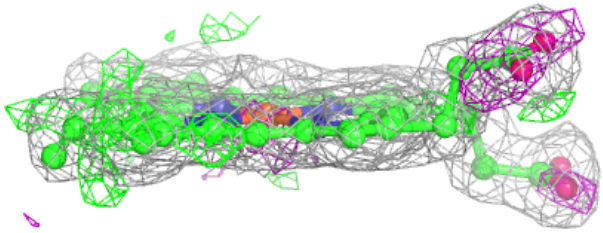
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	HEM	I	200[A]	43/43	0.94	0.12	9,16,28,32	43
2	HEM	F	200[A]	43/43	0.94	0.12	5,16,26,33	43
2	HEM	G	200[A]	43/43	0.94	0.13	10,16,30,36	43
4	SO4	D	1159	5/5	0.94	0.27	60,65,69,70	0
2	HEM	E	200[B]	43/43	0.94	0.11	6,13,30,36	43
3	ZN	L	201	1/1	0.99	0.02	24,24,24,24	0
3	ZN	H	201	1/1	0.99	0.02	24,24,24,24	0
3	ZN	B	202	1/1	0.99	0.02	21,21,21,21	0
4	SO4	E	1158	5/5	0.99	0.07	25,25,28,28	0
3	ZN	K	201	1/1	0.99	0.02	23,23,23,23	0
3	ZN	J	202	1/1	0.99	0.03	20,20,20,20	0
4	SO4	C	1158	5/5	0.99	0.06	25,25,26,29	0
3	ZN	I	202	1/1	0.99	0.03	21,21,21,21	0
3	ZN	C	201	1/1	0.99	0.03	23,23,23,23	0
3	ZN	G	202	1/1	1.00	0.02	20,20,20,20	0
3	ZN	I	201	1/1	1.00	0.02	23,23,23,23	0
3	ZN	D	202	1/1	1.00	0.02	22,22,22,22	0
3	ZN	A	202	1/1	1.00	0.03	21,21,21,21	0
3	ZN	E	202	1/1	1.00	0.03	21,21,21,21	0
3	ZN	J	201	1/1	1.00	0.02	23,23,23,23	0
3	ZN	H	202	1/1	1.00	0.03	20,20,20,20	0
3	ZN	L	202	1/1	1.00	0.02	20,20,20,20	0
3	ZN	K	202	1/1	1.00	0.03	21,21,21,21	0
4	SO4	B	1158	5/5	1.00	0.05	24,25,26,28	0
3	ZN	G	201	1/1	1.00	0.02	24,24,24,24	0
3	ZN	D	201	1/1	1.00	0.01	24,24,24,24	0
3	ZN	E	201	1/1	1.00	0.02	24,24,24,24	0
3	ZN	B	201	1/1	1.00	0.03	23,23,23,23	0
4	SO4	H	1158	5/5	1.00	0.05	24,25,26,28	0
3	ZN	F	202	1/1	1.00	0.02	21,21,21,21	0
3	ZN	F	201	1/1	1.00	0.02	24,24,24,24	0
3	ZN	C	202	1/1	1.00	0.02	21,21,21,21	0
3	ZN	A	201	1/1	1.00	0.02	24,24,24,24	0

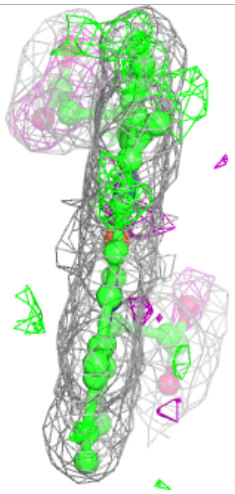
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around HEM A 200 (B):

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

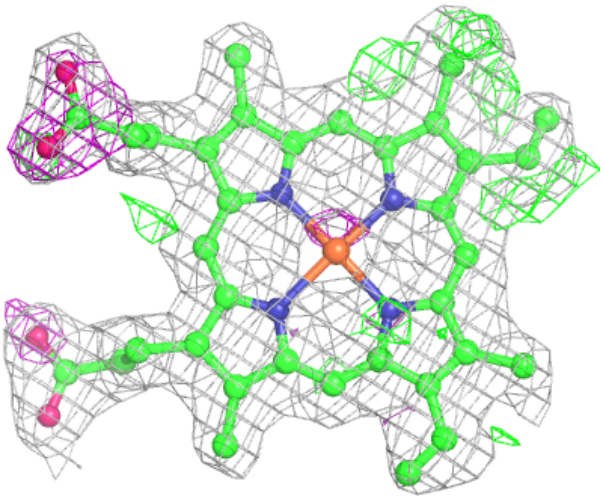


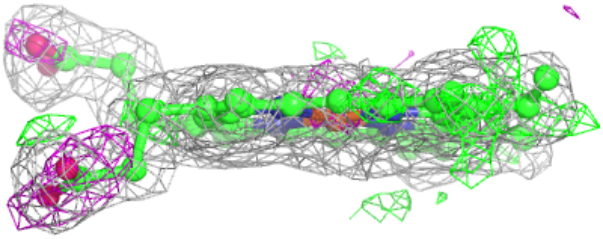


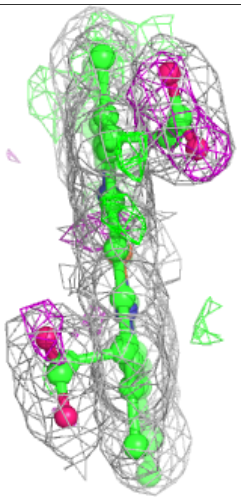


Electron density around HEM A 200 (A):

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

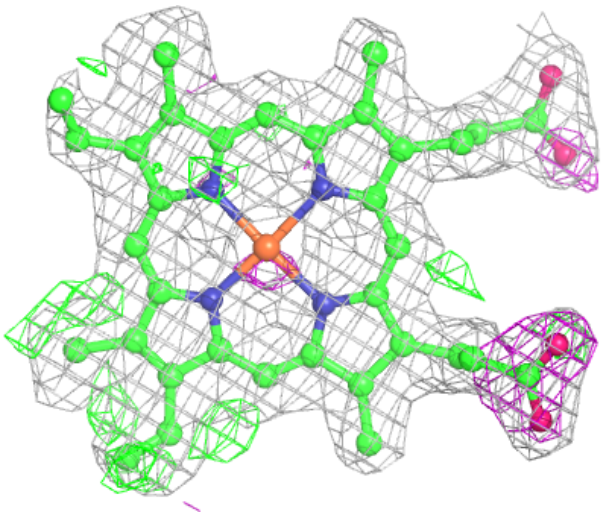


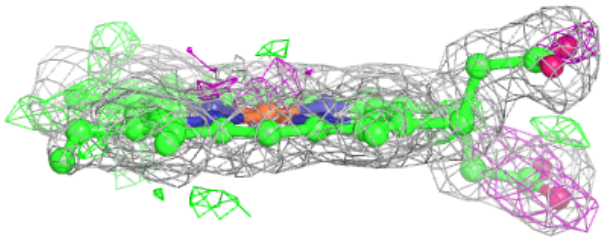


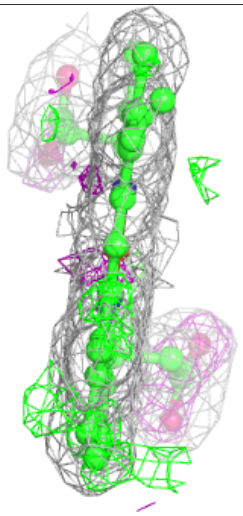


Electron density around HEM B 200 (A):

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

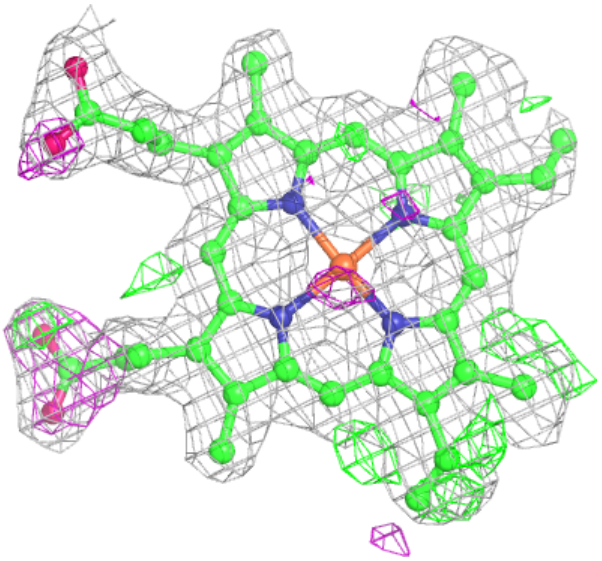


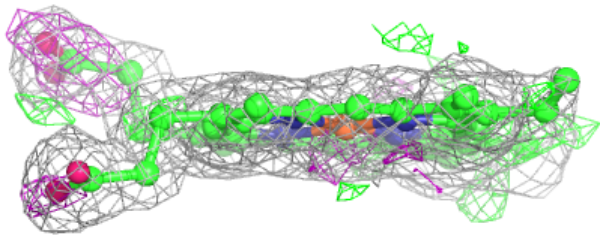


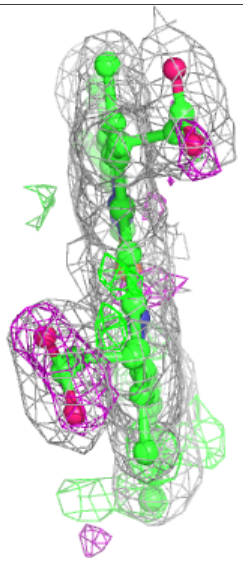


Electron density around HEM B 200 (B):

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

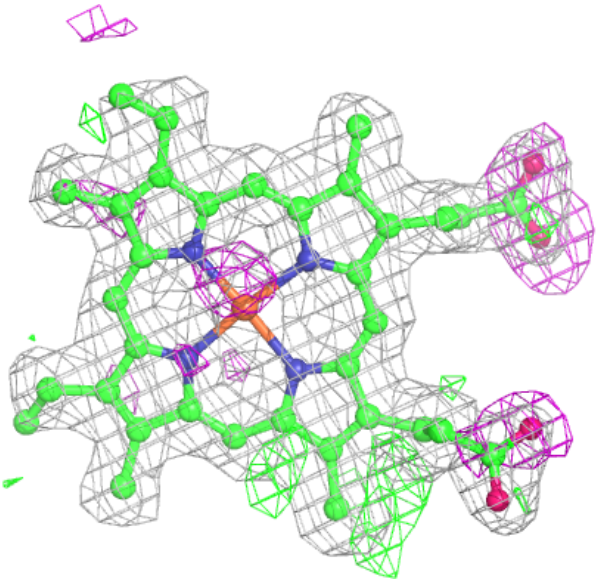


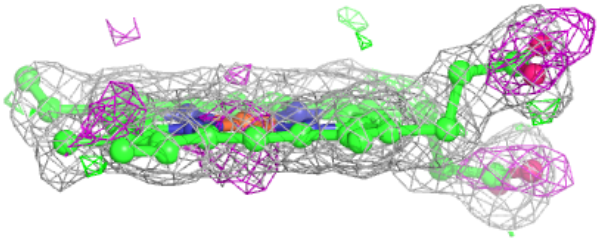


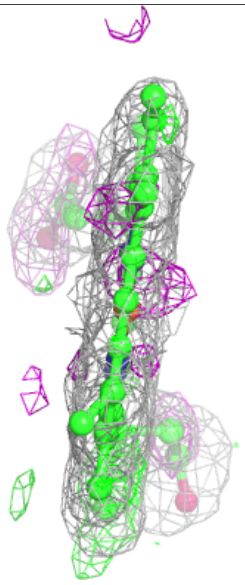


Electron density around HEM L 200 (B):

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

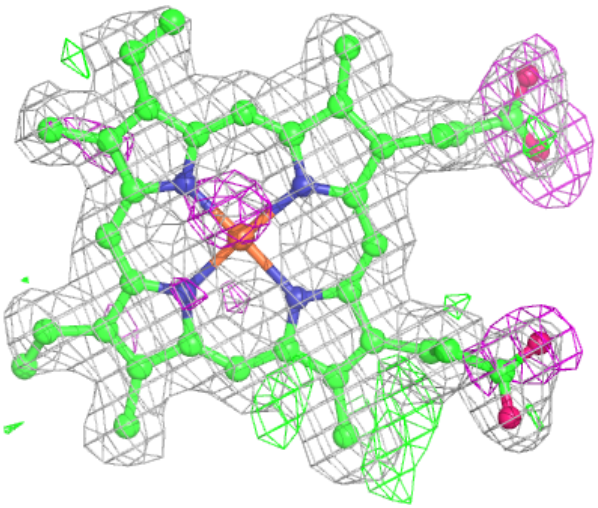


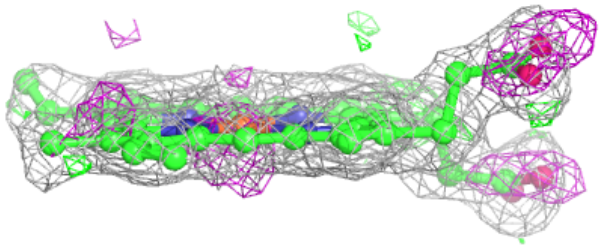


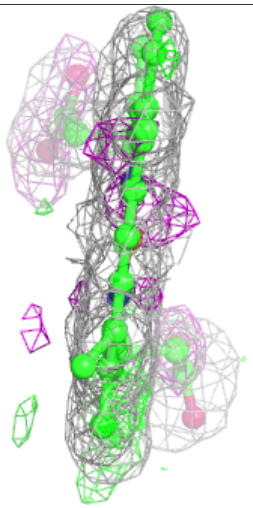


Electron density around HEM L 200 (A):

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

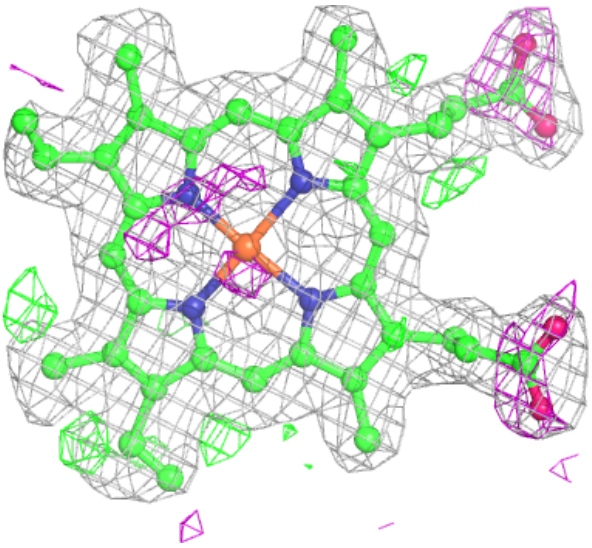


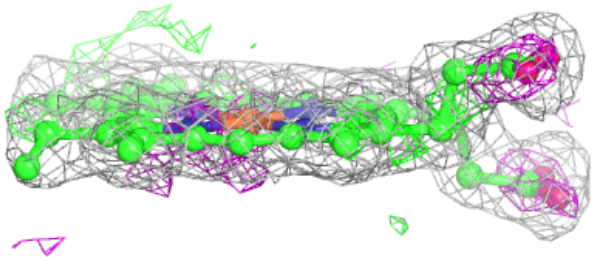


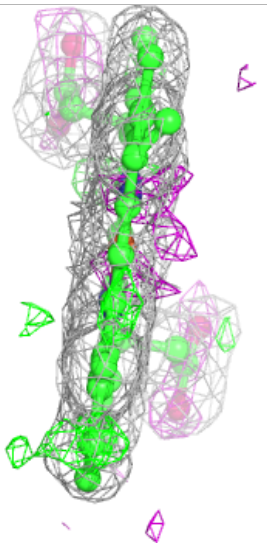


Electron density around HEM D 200 (A):

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

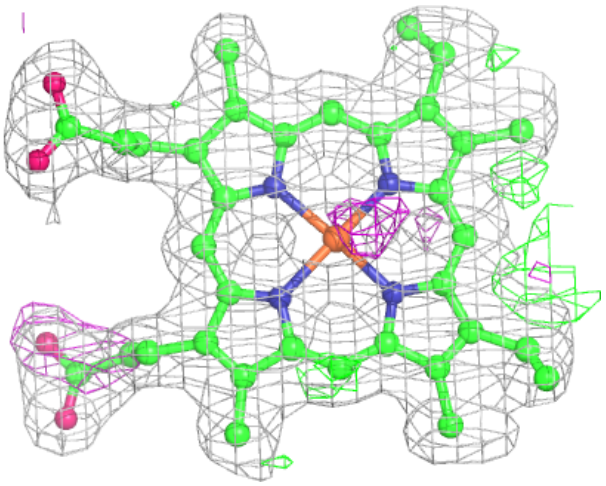


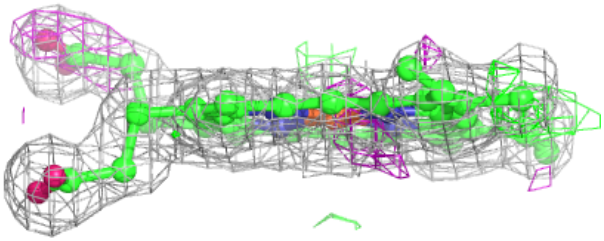


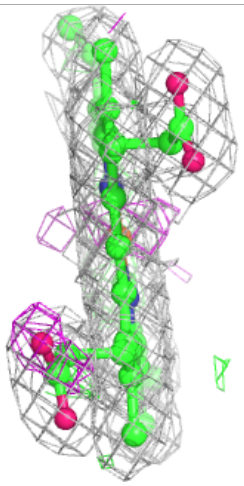


Electron density around HEM E 200 (A):

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

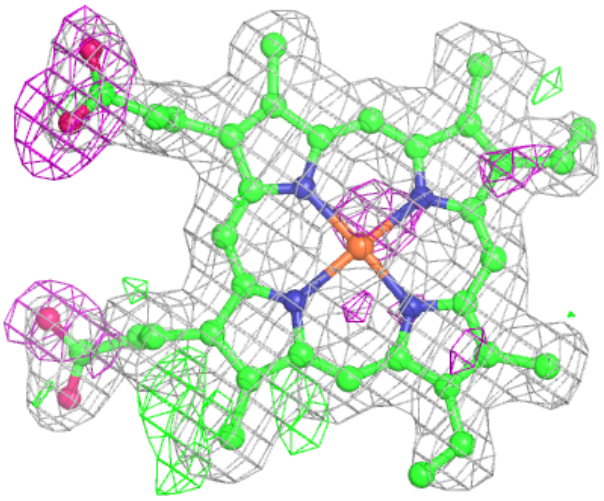


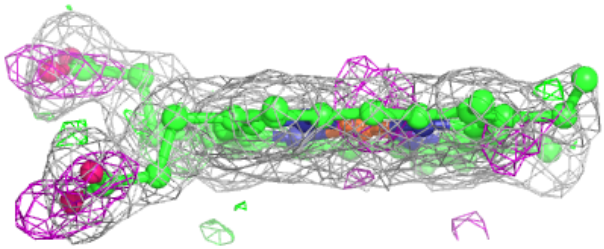


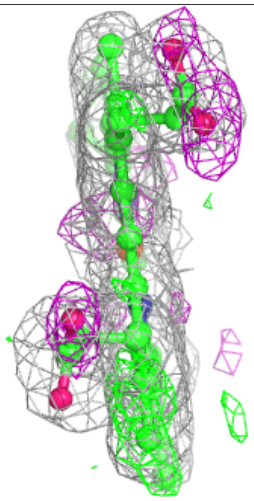


Electron density around HEM K 200 (A):

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

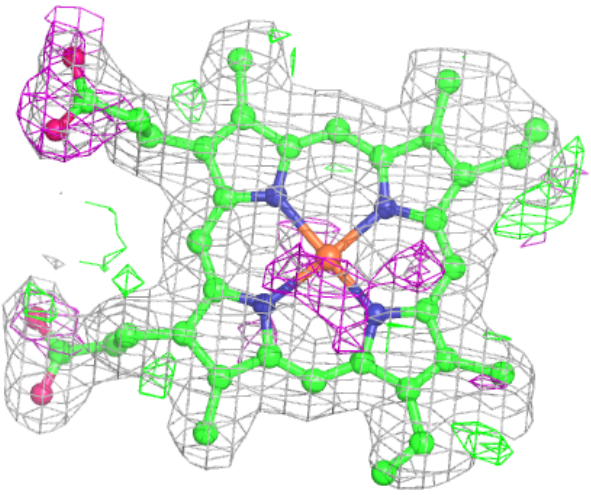


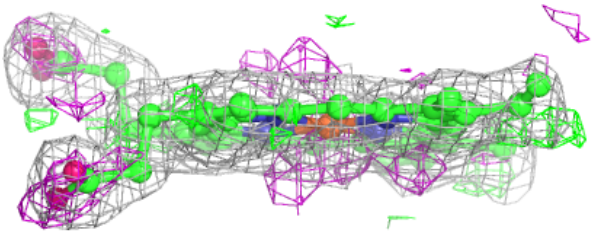


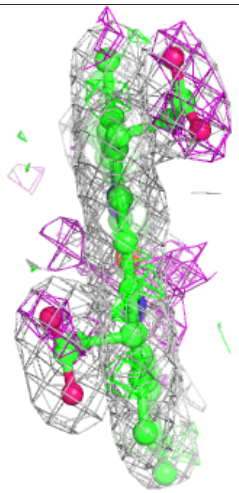


Electron density around HEM H 200 (A):

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

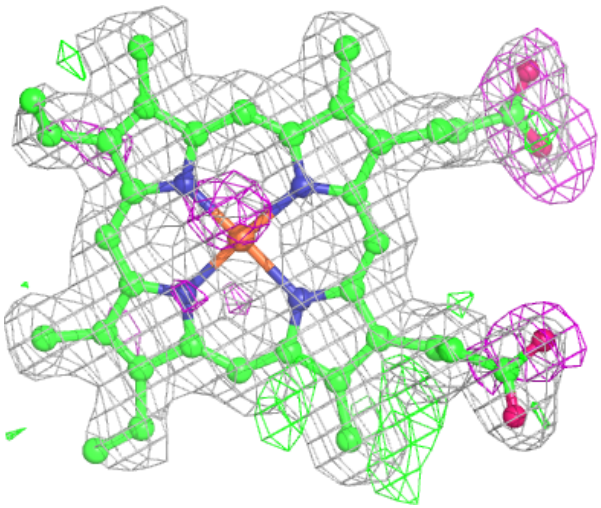


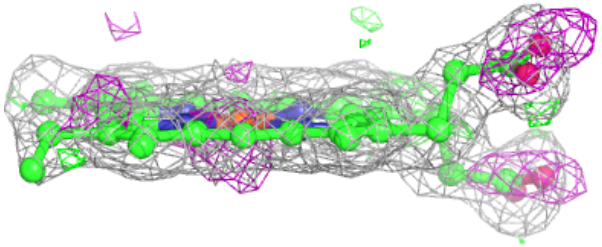


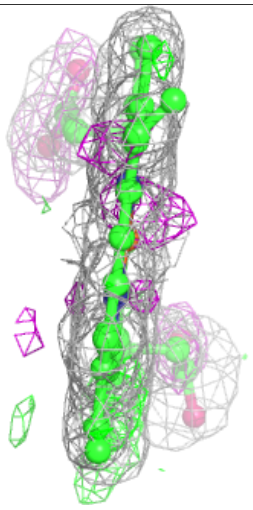


Electron density around HEM K 200 (B):

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

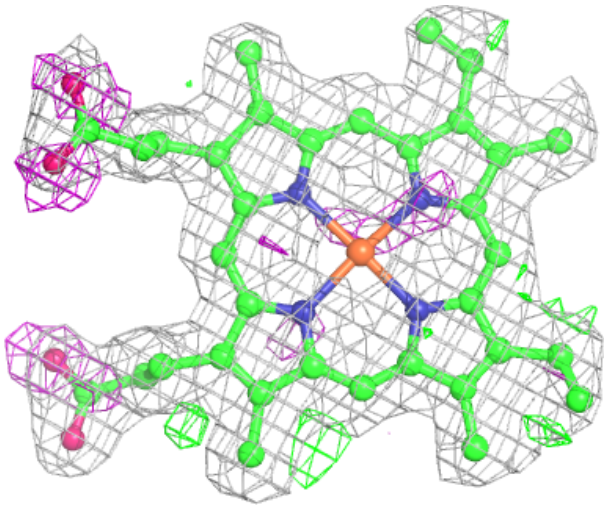


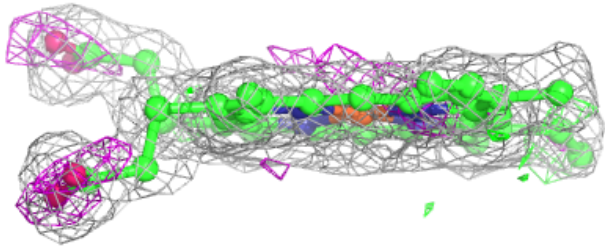


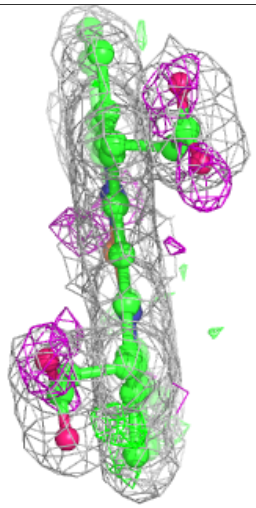


Electron density around HEM J 200 (A):

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

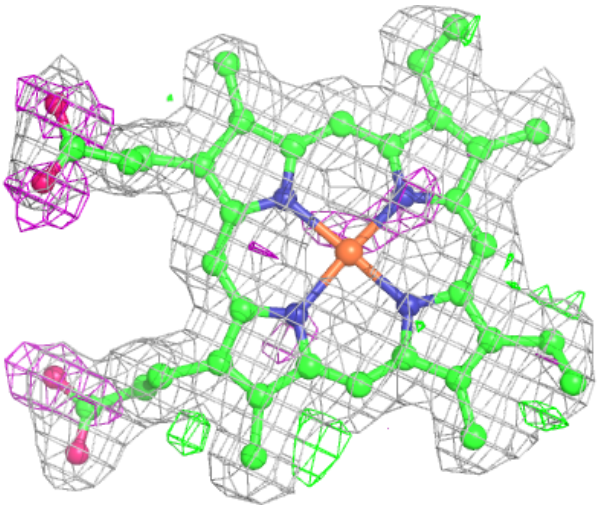


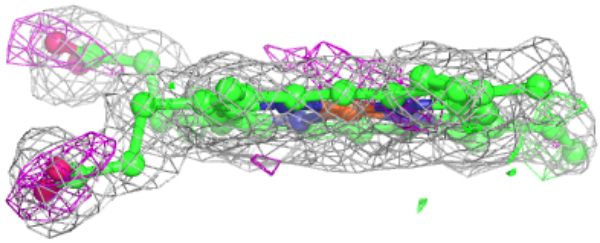


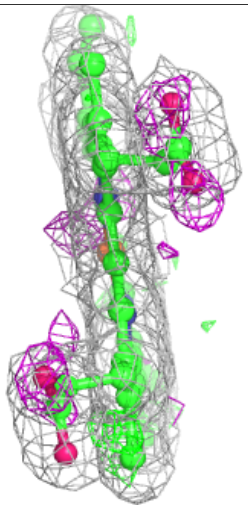


Electron density around HEM J 200 (B):

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

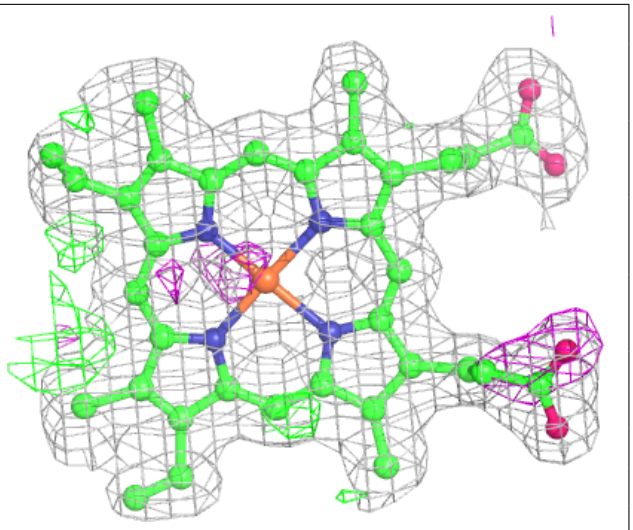


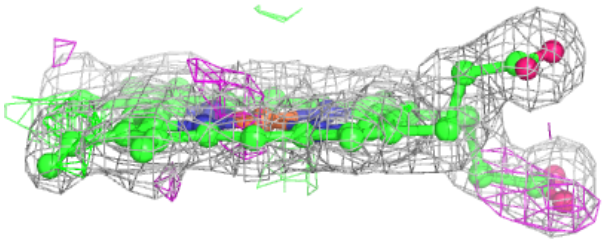


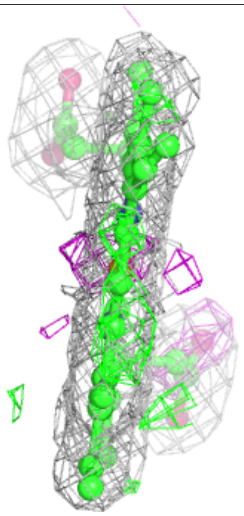


Electron density around HEM F 200 (B):

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

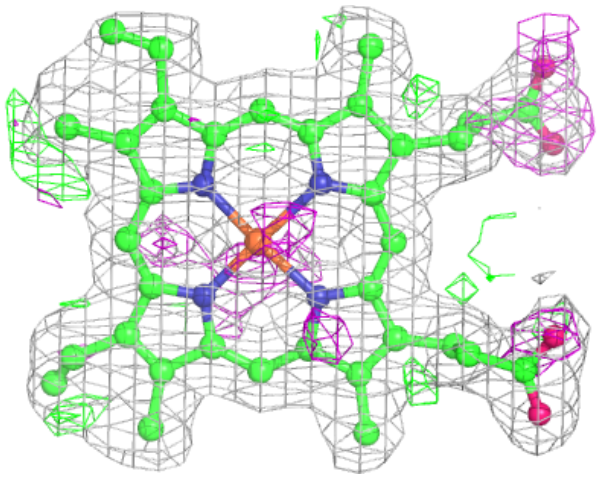


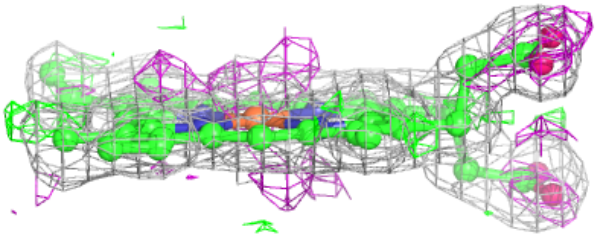


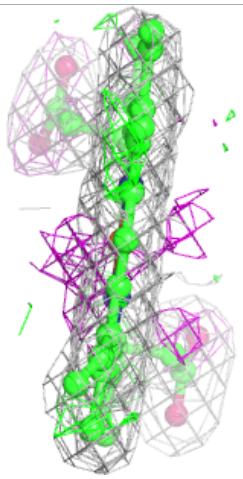


Electron density around HEM G 200 (B):

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

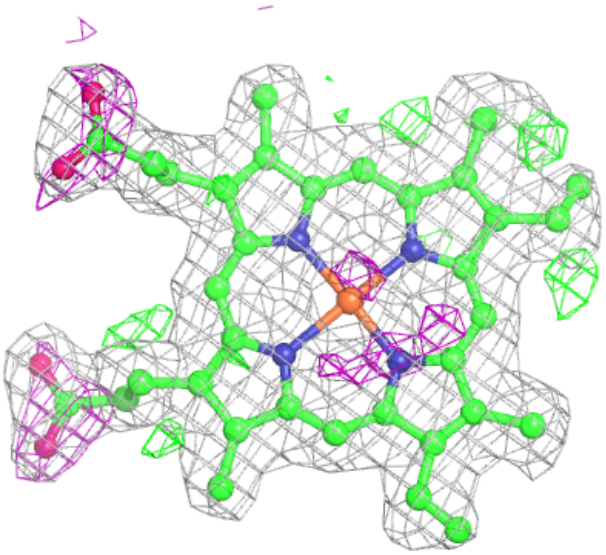


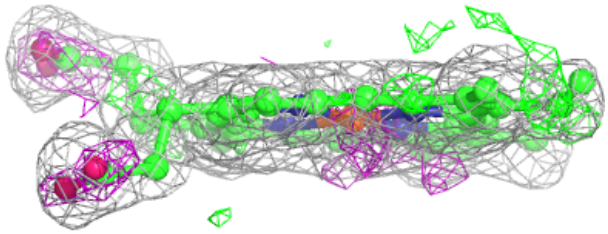


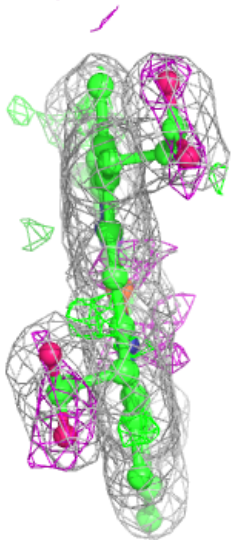


Electron density around HEM C 200 (A):

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

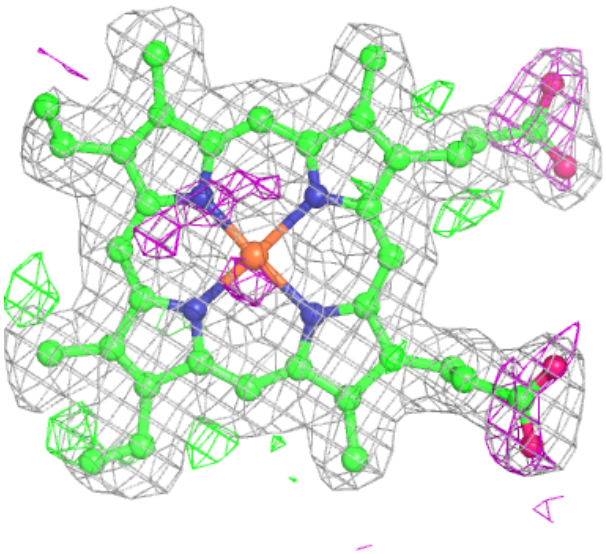


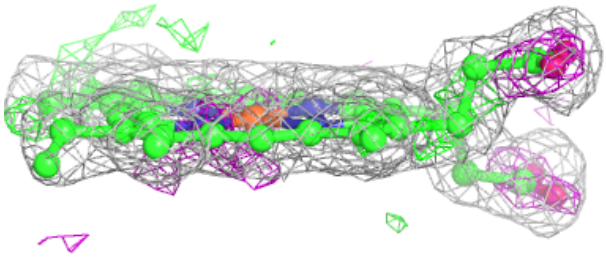


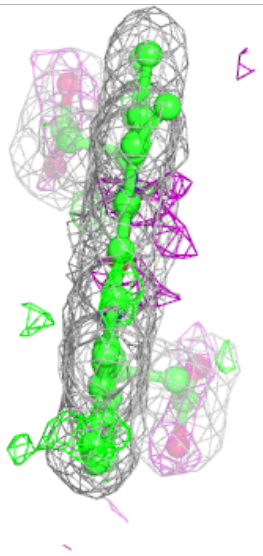


Electron density around HEM D 200 (B):

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

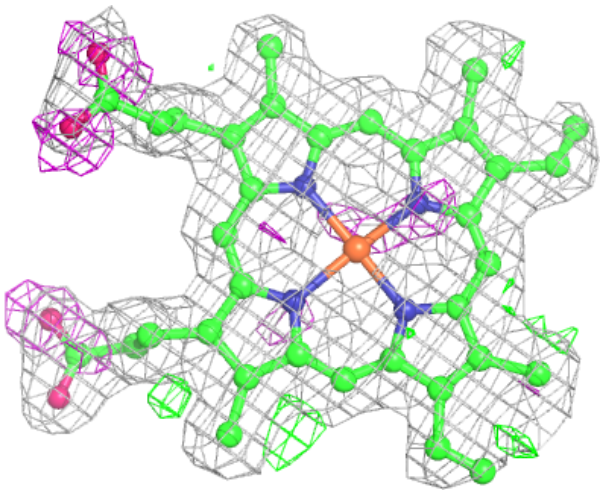


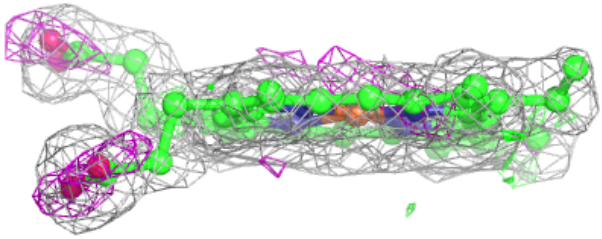


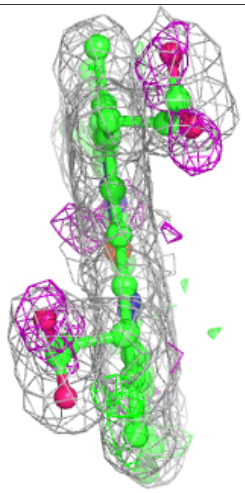


Electron density around HEM I 200 (B):

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

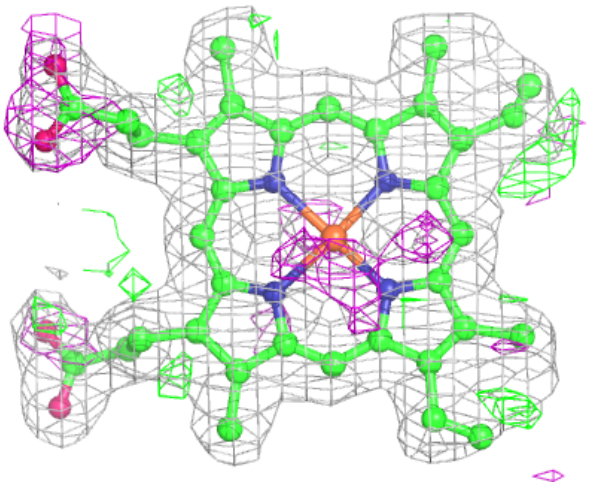


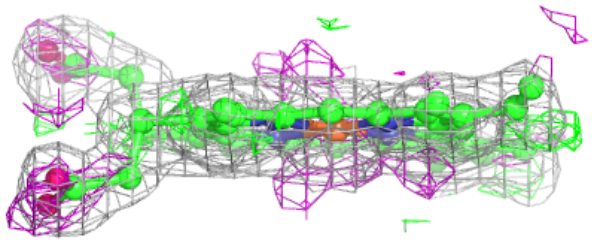


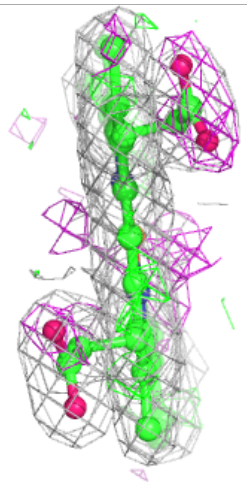


Electron density around HEM H 200 (B):

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

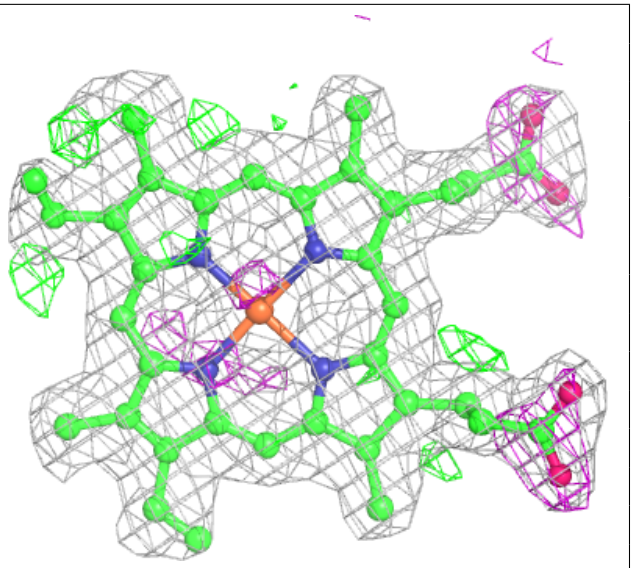


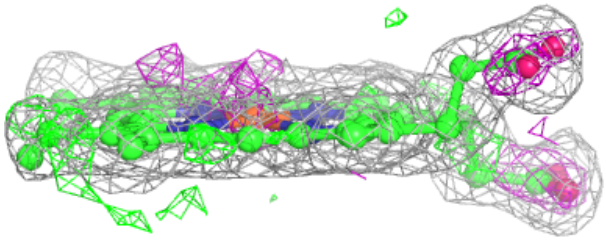


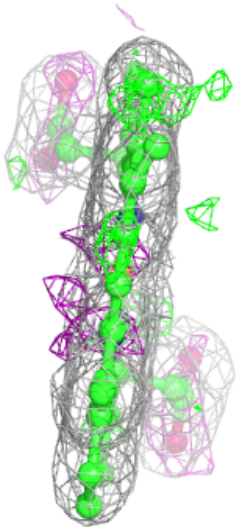


Electron density around HEM C 200 (B):

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

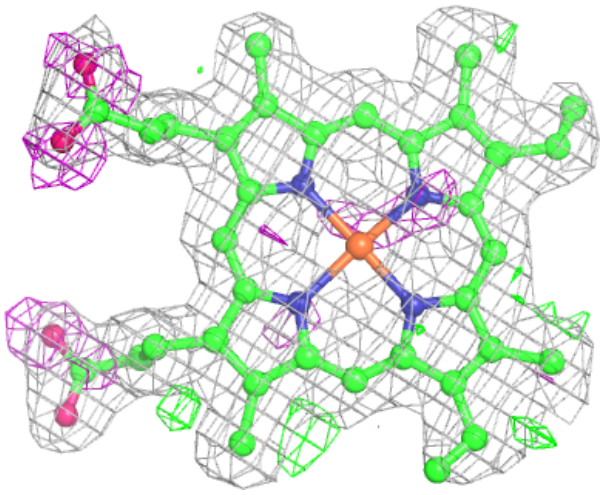


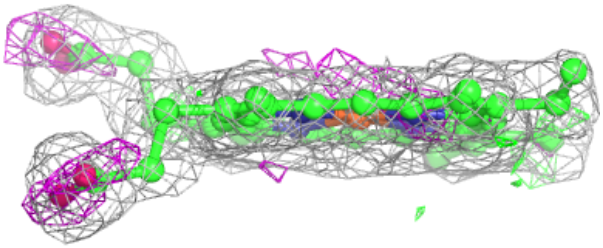


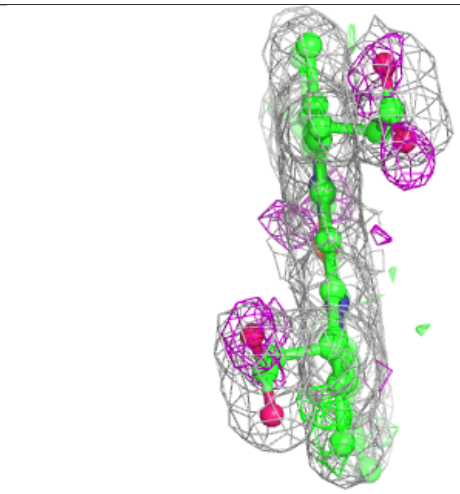


Electron density around HEM I 200 (A):

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

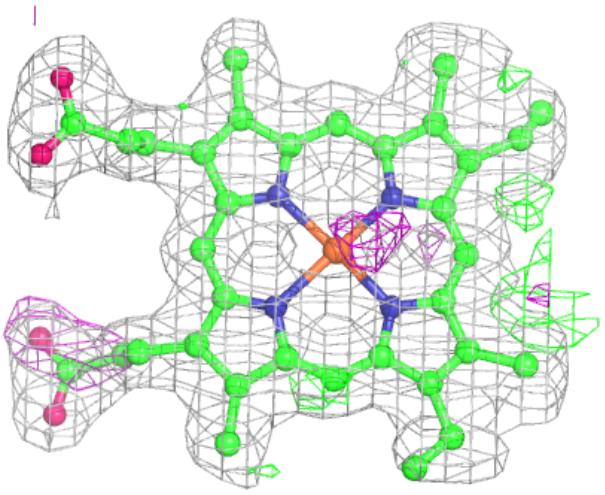


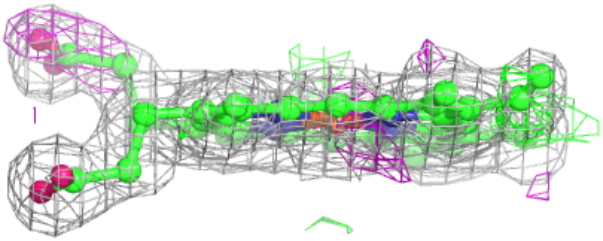


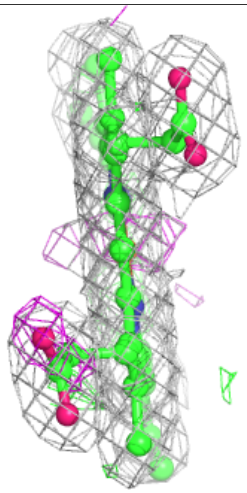


Electron density around HEM F 200 (A):

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

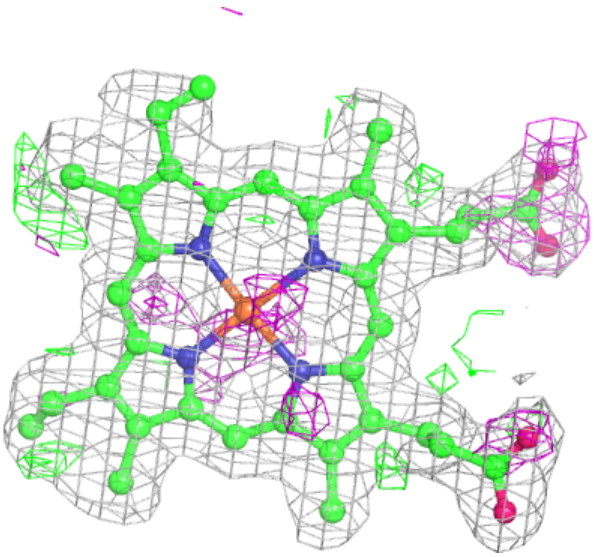


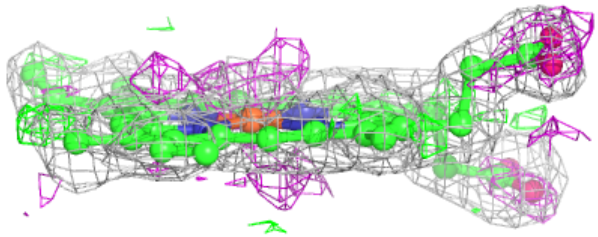


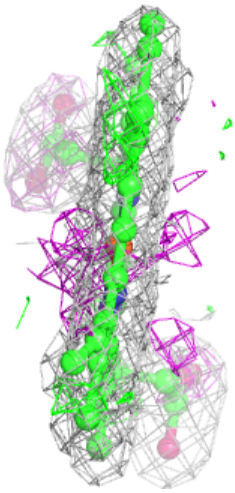


Electron density around HEM G 200 (A):

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

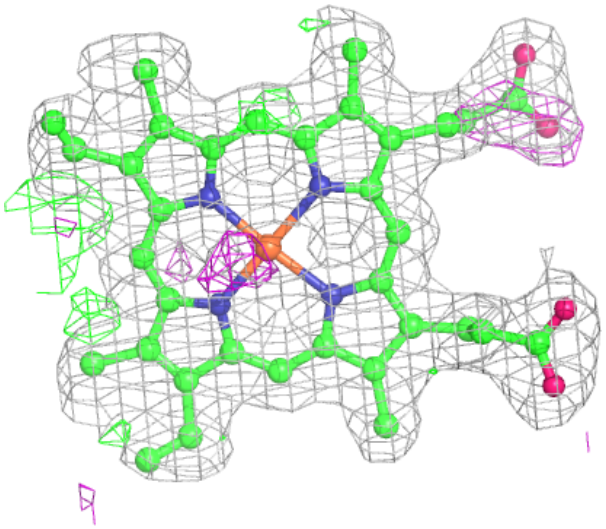


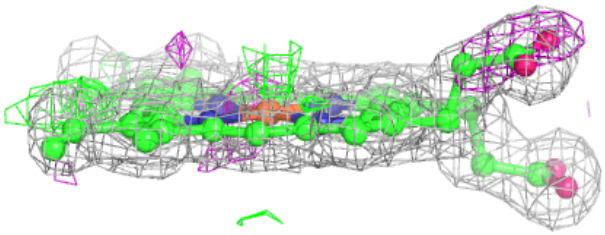


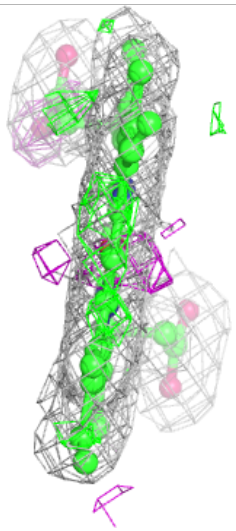


Electron density around HEM E 200 (B):

$2mF_o$-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.