Full wwPDB X-ray Structure Validation Report (i)

Aug 4, 2026 – 06:19 PM EDT

PDB ID	:	2HTN / pdb_00002htn
Title	:	E. coli bacterioferritin in its as-isolated form
Authors	:	van Eerde, A.; Wolterink-Van Loo, S.; Van Der Oost, J.; Dijkstra, B.W.
Deposited on	:	2006-07-26
Resolution	:	2.50 Å(reported)

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

We welcome your comments at *validation@mail.wwpdb.org*
A user guide is available at
https://www.wwpdb.org/validation/2017/XrayValidationReportHelp
with specific help available everywhere you see the (i) symbol.

The types of validation reports are described at
http://www.wwpdb.org/validation/2017/FAQs#types.

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

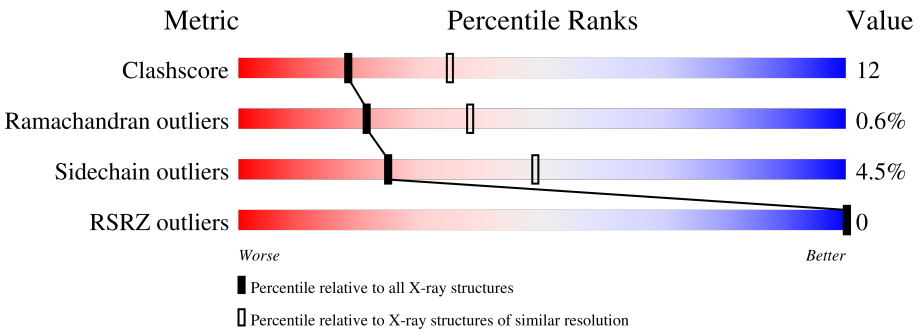
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	2022.3.0, CSD as543be (2022)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.50

1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.50 Å.

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



Metric	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
Clashscore	190562	6492 (2.50-2.50)
Ramachandran outliers	187476	6378 (2.50-2.50)
Sidechain outliers	187428	6380 (2.50-2.50)
RSRZ outliers	180081	5833 (2.50-2.50)

The table below summarises the geometric issues observed across the polymeric chains and their fit to the electron density. The red, orange, yellow and green segments of the lower bar indicate the fraction of residues that contain outliers for >=3, 2, 1 and 0 types of geometric quality criteria respectively. A grey segment represents the fraction of residues that are not modelled. The numeric value for each fraction is indicated below the corresponding segment, with a dot representing fractions <=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	67% 31% •
1	B	158	72% 26% •
1	C	158	70% 28% •
1	D	158	74% 24% •
1	E	158	72% 27% •
1	F	158	67% 31% •

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Mol	Chain	Length	Quality of chain
1	G	158	72% 26% •
1	H	158	72% 25% •

2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 10846 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
			Total	C	N	O	S			
1	A	158	1299	818	222	252	7	0	0	0
1	B	158	1299	818	222	252	7	0	0	0
1	C	158	1299	818	222	252	7	0	0	0
1	D	158	1299	818	222	252	7	0	0	0
1	E	158	1299	818	222	252	7	0	0	0
1	F	158	1299	818	222	252	7	0	0	0
1	G	158	1299	818	222	252	7	0	0	0
1	H	158	1299	818	222	252	7	0	0	0

- Molecule 2 is FE (III) ION (CCD ID: FE) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Fe		
2	A	2	2	2	0	0
2	B	2	2	2	0	0
2	C	2	2	2	0	0
2	D	2	2	2	0	0
2	E	2	2	2	0	0
2	F	2	2	2	0	0

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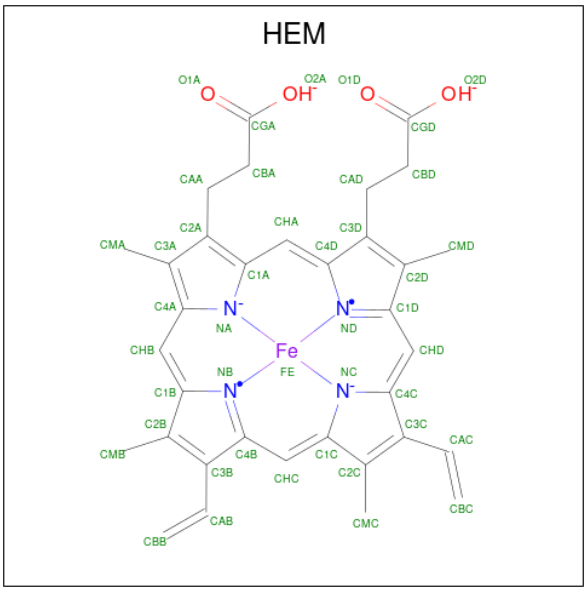
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Fe		
2	G	2	2	2	0	0
			Total	Fe		
2	H	2	2	2	0	0

- Molecule 3 is PROTOPORPHYRIN IX CONTAINING FE (CCD ID: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
3	A	1	85	68	1	8	8	0	1
			Total	C	Fe	N	O		
3	C	1	85	68	1	8	8	0	1
			Total	C	Fe	N	O		
3	E	1	85	68	1	8	8	0	1
			Total	C	Fe	N	O		
3	G	1	85	68	1	8	8	0	1

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	A	19	19	19	0	0
			Total	O		
4	B	19	19	19	0	0
			Total	O		
4	C	6	6	6	0	0

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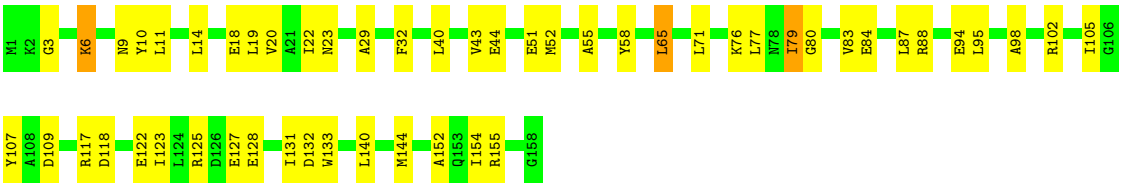
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
4	D	9	9	9	0	0
			Total	O		
4	E	12	12	12	0	0
			Total	O		
4	F	9	9	9	0	0
			Total	O		
4	G	16	16	16	0	0
			Total	O		
4	H	8	8	8	0	0

3 Residue-property plots (i)

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

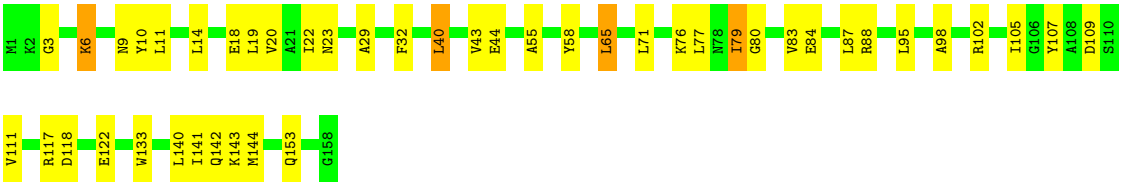
• Molecule 1: Bacterioferritin

Chain A:



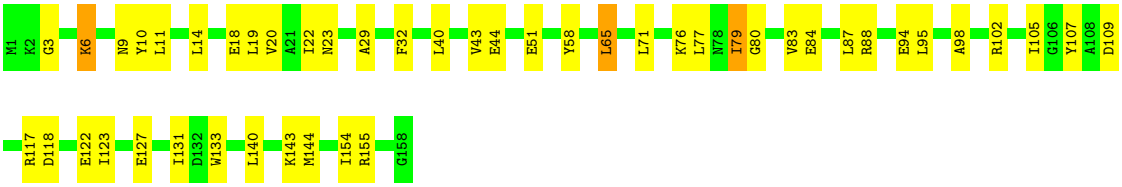
• Molecule 1: Bacterioferritin

Chain B:



• Molecule 1: Bacterioferritin

Chain C:



• Molecule 1: Bacterioferritin

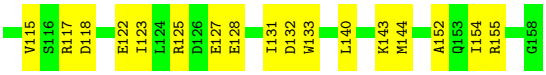
Chain D:

• Molecule 1: Bacterioferritin

Chain E:

• Molecule 1: Bacterioferritin

Chain F:



• Molecule 1: Bacterioferritin

Chain G:

• Molecule 1: Bacterioferritin

Chain H:

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants a, b, c, α, β, γ	167.89Å 167.89Å 167.89Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	59.34 – 2.50 59.34 – 2.50	Depositor EDS
% Data completeness (in resolution range)	99.8 (59.34-2.50) 99.8 (59.34-2.50)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	3.47 (at 2.51Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.171 , 0.188 0.160 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	wwPDB-VP
Wilson B-factor (Å^2)	26.0	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.36 , 19.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.38$, $< L^2 > = 0.21$	Xtriage
Estimated twinning fraction	0.477 for l,-k,h	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	10846	wwPDB-VP
Average B, all atoms (Å^2)	26.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.07% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

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

5 Model quality (i)

5.1 Standard geometry (i)

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

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	0.71	0/1318	0.94	0/1770
1	B	0.70	0/1318	0.97	2/1770 (0.1%)
1	C	0.71	0/1318	0.95	1/1770 (0.1%)
1	D	0.72	0/1318	0.98	0/1770
1	E	0.70	0/1318	0.93	1/1770 (0.1%)
1	F	0.71	0/1318	0.94	1/1770 (0.1%)
1	G	0.69	0/1318	0.97	0/1770
1	H	0.72	0/1318	0.97	1/1770 (0.1%)
All	All	0.71	0/10544	0.96	6/14160 (0.0%)

There are no bond length outliers.

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	143	LYS	N-CA-C	5.66	117.53	111.36
1	B	141	ILE	N-CA-C	-5.42	105.22	110.42
1	E	143	LYS	N-CA-C	5.36	117.20	111.36
1	H	143	LYS	N-CA-C	5.10	116.92	111.36
1	F	143	LYS	N-CA-C	5.08	116.90	111.36
1	C	143	LYS	N-CA-C	5.00	116.81	111.36

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	1299	0	1280	42	0
1	B	1299	0	1280	30	0
1	C	1299	0	1280	34	0
1	D	1299	0	1280	32	0
1	E	1299	0	1280	32	2
1	F	1299	0	1280	40	0
1	G	1299	0	1280	34	2
1	H	1299	0	1280	34	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
2	C	2	0	0	0	0
2	D	2	0	0	0	0
2	E	2	0	0	0	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
3	A	85	0	60	5	0
3	C	85	0	60	0	0
3	E	85	0	60	0	0
3	G	85	0	60	2	0
4	A	19	0	0	0	0
4	B	19	0	0	0	0
4	C	6	0	0	1	0
4	D	9	0	0	0	0
4	E	12	0	0	0	0
4	F	9	0	0	0	0
4	G	16	0	0	0	0
4	H	8	0	0	0	0
All	All	10846	0	10480	258	2

The all-atom clashscore is defined as the number of clashes found per 1000 atoms (including hydrogen atoms). The all-atom clashscore for this structure is 12.

All (258) 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:A:109:ASP:HB2	1:A:117:ARG:NH1	1.94	0.83
1:H:109:ASP:HB2	1:H:117:ARG:NH1	1.94	0.82
1:E:109:ASP:HB2	1:E:117:ARG:NH1	1.94	0.82

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:109:ASP:HB2	1:F:117:ARG:NH1	1.94	0.81
1:B:109:ASP:HB2	1:B:117:ARG:NH1	1.95	0.81
1:C:109:ASP:HB2	1:C:117:ARG:NH1	1.94	0.81
1:D:109:ASP:HB2	1:D:117:ARG:NH1	1.95	0.81
1:G:109:ASP:HB2	1:G:117:ARG:NH1	1.94	0.81
1:D:43:VAL:HG11	1:D:133:TRP:CZ2	2.24	0.72
1:H:20:VAL:HG13	1:H:77:LEU:HD23	1.76	0.68
1:B:20:VAL:HG13	1:B:77:LEU:HD23	1.76	0.68
1:B:43:VAL:HG11	1:B:133:TRP:CZ2	2.29	0.67
1:E:20:VAL:HG13	1:E:77:LEU:HD23	1.77	0.67
1:F:20:VAL:HG13	1:F:77:LEU:HD23	1.77	0.67
1:G:20:VAL:HG13	1:G:77:LEU:HD23	1.77	0.66
1:C:20:VAL:HG13	1:C:77:LEU:HD23	1.77	0.66
1:D:20:VAL:HG13	1:D:77:LEU:HD23	1.77	0.66
1:A:20:VAL:HG13	1:A:77:LEU:HD23	1.77	0.65
1:E:43:VAL:HG11	1:E:133:TRP:CZ2	2.34	0.63
1:C:43:VAL:HG11	1:C:133:TRP:CZ2	2.33	0.63
1:F:43:VAL:HG11	1:F:133:TRP:CZ2	2.33	0.63
1:A:43:VAL:HG11	1:A:133:TRP:CZ2	2.33	0.63
1:G:43:VAL:HG11	1:G:133:TRP:CZ2	2.36	0.60
1:D:18:GLU:O	1:D:22:ILE:HG13	2.02	0.59
1:D:3:GLY:HA3	1:D:65:LEU:HG	1.86	0.58
1:B:3:GLY:HA3	1:B:65:LEU:HG	1.86	0.58
1:E:3:GLY:HA3	1:E:65:LEU:HG	1.86	0.58
1:H:43:VAL:HG11	1:H:133:TRP:CZ2	2.38	0.58
1:H:3:GLY:HA3	1:H:65:LEU:HG	1.86	0.58
1:A:3:GLY:HA3	1:A:65:LEU:HG	1.85	0.58
1:G:3:GLY:HA3	1:G:65:LEU:HG	1.86	0.58
1:B:83:VAL:O	1:B:87:LEU:HD13	2.05	0.57
1:C:3:GLY:HA3	1:C:65:LEU:HG	1.85	0.57
1:C:83:VAL:O	1:C:87:LEU:HD13	2.04	0.57
1:E:83:VAL:O	1:E:87:LEU:HD13	2.05	0.57
1:F:3:GLY:HA3	1:F:65:LEU:HG	1.85	0.57
1:A:83:VAL:O	1:A:87:LEU:HD13	2.05	0.56
1:D:83:VAL:O	1:D:87:LEU:HD13	2.05	0.56
1:H:83:VAL:O	1:H:87:LEU:HD13	2.05	0.56
1:G:83:VAL:O	1:G:87:LEU:HD13	2.05	0.56
1:A:152:ALA:HB1	1:H:154:ILE:HD11	1.86	0.56
1:F:83:VAL:O	1:F:87:LEU:HD13	2.05	0.55
1:D:140:LEU:HB3	1:D:144:MET:HE3	1.89	0.55
1:C:79:ILE:HD13	4:C:304:HOH:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:140:LEU:HB3	1:H:144:MET:HE3	1.89	0.55
1:E:123:ILE:O	1:E:127:GLU:HG2	2.08	0.54
1:B:140:LEU:HB3	1:B:144:MET:HE3	1.90	0.54
1:F:123:ILE:O	1:F:127:GLU:HG2	2.08	0.53
1:A:123:ILE:O	1:A:127:GLU:HG2	2.08	0.53
1:A:118:ASP:O	1:A:122:GLU:HG3	2.09	0.53
1:C:123:ILE:O	1:C:127:GLU:HG2	2.09	0.53
1:E:118:ASP:O	1:E:122:GLU:HG3	2.09	0.52
1:B:118:ASP:O	1:B:122:GLU:HG3	2.10	0.52
1:D:128:GLU:OE1	1:F:115:VAL:HG11	2.10	0.52
1:G:118:ASP:O	1:G:122:GLU:HG3	2.09	0.52
1:H:118:ASP:O	1:H:122:GLU:HG3	2.09	0.52
1:F:118:ASP:O	1:F:122:GLU:HG3	2.09	0.52
1:F:152:ALA:HB1	1:G:154:ILE:HD11	1.91	0.52
1:C:118:ASP:O	1:C:122:GLU:HG3	2.09	0.51
1:B:40:LEU:HD13	1:B:153:GLN:OE1	2.10	0.51
1:D:118:ASP:O	1:D:122:GLU:HG3	2.10	0.51
1:B:43:VAL:HG11	1:B:133:TRP:CH2	2.45	0.51
1:D:10:TYR:O	1:D:14:LEU:HD13	2.11	0.50
1:G:10:TYR:O	1:G:14:LEU:HD13	2.12	0.50
1:H:79:ILE:HG22	1:H:80:GLY:N	2.27	0.50
1:E:10:TYR:O	1:E:14:LEU:HD13	2.12	0.50
1:E:79:ILE:HG22	1:E:80:GLY:N	2.27	0.50
1:A:152:ALA:HB1	1:H:154:ILE:CD1	2.42	0.50
1:D:79:ILE:HG22	1:D:80:GLY:N	2.26	0.50
1:A:55:ALA:HB3	3:A:200[A]:HEM:HBC2	1.94	0.49
1:G:79:ILE:HG22	1:G:80:GLY:N	2.27	0.49
1:A:79:ILE:HG22	1:A:80:GLY:N	2.27	0.49
1:F:79:ILE:HG22	1:F:80:GLY:N	2.27	0.49
1:C:10:TYR:O	1:C:14:LEU:HD13	2.11	0.49
1:E:154:ILE:HG13	1:E:155:ARG:HG3	1.95	0.49
1:F:10:TYR:O	1:F:14:LEU:HD13	2.13	0.49
1:A:154:ILE:HG13	1:A:155:ARG:HG3	1.95	0.49
1:B:10:TYR:O	1:B:14:LEU:HD13	2.12	0.49
1:C:79:ILE:HG22	1:C:80:GLY:N	2.27	0.49
1:A:55:ALA:HB3	3:A:200[A]:HEM:CBC	2.42	0.48
1:F:154:ILE:HG13	1:F:155:ARG:HG3	1.95	0.48
1:H:40:LEU:HD13	1:H:153:GLN:OE1	2.12	0.48
1:G:140:LEU:HB3	1:G:144:MET:HE3	1.96	0.48
1:C:154:ILE:HG13	1:C:155:ARG:HG3	1.95	0.48
1:A:10:TYR:O	1:A:14:LEU:HD13	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:19:LEU:HD11	1:B:71:LEU:HD23	1.95	0.48
1:C:19:LEU:HD11	1:C:71:LEU:HD23	1.96	0.48
1:H:10:TYR:O	1:H:14:LEU:HD13	2.13	0.47
1:D:128:GLU:CD	1:F:115:VAL:HG21	2.39	0.47
1:B:79:ILE:HG22	1:B:80:GLY:N	2.28	0.47
1:H:19:LEU:HD11	1:H:71:LEU:HD23	1.96	0.47
1:A:23:ASN:HD22	1:B:71:LEU:HB3	1.80	0.47
1:E:23:ASN:HD22	1:F:71:LEU:HB3	1.79	0.47
1:A:19:LEU:HD11	1:A:71:LEU:HD23	1.97	0.47
1:A:125:ARG:NH2	1:G:122:GLU:OE2	2.48	0.47
1:E:18:GLU:O	1:E:22:ILE:HG13	2.15	0.47
1:G:19:LEU:HD11	1:G:71:LEU:HD23	1.97	0.47
1:E:19:LEU:HD11	1:E:71:LEU:HD23	1.96	0.46
1:A:18:GLU:O	1:A:22:ILE:HG13	2.15	0.46
1:D:43:VAL:HG11	1:D:133:TRP:CH2	2.51	0.46
1:A:52:MET:HB3	3:A:200[B]:HEM:CHB	2.45	0.46
1:F:19:LEU:HD11	1:F:71:LEU:HD23	1.96	0.46
1:F:152:ALA:HB1	1:G:154:ILE:CD1	2.46	0.46
3:G:200[B]:HEM:HBC2	1:H:55:ALA:HB3	1.97	0.46
1:F:18:GLU:O	1:F:22:ILE:HG13	2.16	0.46
1:G:23:ASN:HD22	1:H:71:LEU:HB3	1.81	0.45
1:C:84:GLU:O	1:C:88:ARG:HG3	2.17	0.45
1:D:19:LEU:HD11	1:D:71:LEU:HD23	1.97	0.45
1:E:32:PHE:HE2	1:E:44:GLU:HG3	1.82	0.45
1:A:154:ILE:HG13	1:A:155:ARG:N	2.31	0.45
1:E:71:LEU:HB3	1:F:23:ASN:HD22	1.81	0.45
1:E:88:ARG:HH11	1:E:88:ARG:HG2	1.82	0.45
1:F:128:GLU:OE2	1:H:1:MET:HE1	2.16	0.45
1:A:6:LYS:HB3	1:A:107:TYR:CE2	2.52	0.45
3:A:200[B]:HEM:CBC	1:B:55:ALA:HB3	2.46	0.45
1:G:32:PHE:HE2	1:G:44:GLU:HG3	1.82	0.45
1:H:32:PHE:HE2	1:H:44:GLU:HG3	1.82	0.45
1:A:32:PHE:HE2	1:A:44:GLU:HG3	1.82	0.44
1:B:84:GLU:O	1:B:88:ARG:HG3	2.17	0.44
1:C:18:GLU:O	1:C:22:ILE:HG13	2.17	0.44
1:C:154:ILE:HG13	1:C:155:ARG:N	2.32	0.44
1:A:71:LEU:HB3	1:B:23:ASN:HD22	1.82	0.44
1:A:88:ARG:HG2	1:A:88:ARG:HH11	1.82	0.44
1:D:1:MET:HE1	1:H:128:GLU:OE2	2.16	0.44
1:D:32:PHE:HE2	1:D:44:GLU:HG3	1.82	0.44
1:F:88:ARG:HG2	1:F:88:ARG:HH11	1.82	0.44

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:D:6:LYS:HB3	1:D:107:TYR:CE2	2.52	0.44
1:D:84:GLU:O	1:D:88:ARG:HG3	2.17	0.44
1:F:154:ILE:HG13	1:F:155:ARG:N	2.32	0.44
1:E:6:LYS:HB3	1:E:107:TYR:CE2	2.52	0.44
1:F:6:LYS:HB3	1:F:107:TYR:CE2	2.52	0.44
1:C:6:LYS:HB3	1:C:107:TYR:CE2	2.52	0.44
1:C:43:VAL:HG11	1:C:133:TRP:CH2	2.52	0.44
1:D:29:ALA:HB2	1:D:44:GLU:HB3	2.00	0.44
1:F:132:ASP:HB2	1:H:64:PHE:CE1	2.52	0.44
1:H:84:GLU:O	1:H:88:ARG:HG3	2.17	0.44
1:E:29:ALA:HB2	1:E:44:GLU:HB3	2.00	0.44
1:F:43:VAL:HG11	1:F:133:TRP:CH2	2.53	0.44
1:F:84:GLU:O	1:F:88:ARG:HG3	2.17	0.44
1:A:51:GLU:OE1	1:A:51:GLU:HA	2.18	0.44
1:C:88:ARG:HG2	1:C:88:ARG:HH11	1.83	0.44
1:F:51:GLU:OE1	1:F:51:GLU:HA	2.17	0.44
1:H:32:PHE:CE2	1:H:44:GLU:HG3	2.53	0.44
1:A:79:ILE:HG22	1:A:80:GLY:O	2.18	0.44
1:A:84:GLU:O	1:A:88:ARG:HG3	2.17	0.44
1:B:6:LYS:HB3	1:B:107:TYR:CE2	2.52	0.44
1:B:32:PHE:HE2	1:B:44:GLU:HG3	1.83	0.44
1:E:43:VAL:HG11	1:E:133:TRP:CH2	2.53	0.44
1:F:94:GLU:HB3	1:F:131:ILE:HD11	2.00	0.44
1:H:6:LYS:HB3	1:H:107:TYR:CE2	2.53	0.44
1:C:32:PHE:HE2	1:C:44:GLU:HG3	1.82	0.44
1:C:140:LEU:HB3	1:C:144:MET:HE3	1.98	0.44
1:E:154:ILE:HG13	1:E:155:ARG:N	2.32	0.44
1:G:84:GLU:O	1:G:88:ARG:HG3	2.17	0.44
1:H:43:VAL:HG11	1:H:133:TRP:CH2	2.53	0.44
1:B:79:ILE:HG22	1:B:80:GLY:O	2.18	0.43
1:B:88:ARG:HH11	1:B:88:ARG:HG2	1.83	0.43
1:C:79:ILE:HG22	1:C:80:GLY:O	2.18	0.43
1:F:32:PHE:HE2	1:F:44:GLU:HG3	1.82	0.43
1:G:71:LEU:HB3	1:H:23:ASN:HD22	1.83	0.43
1:G:88:ARG:HH11	1:G:88:ARG:HG2	1.83	0.43
1:A:43:VAL:HG11	1:A:133:TRP:CH2	2.53	0.43
1:E:84:GLU:O	1:E:88:ARG:HG3	2.17	0.43
1:F:140:LEU:HB3	1:F:144:MET:HE3	1.99	0.43
1:G:32:PHE:CE2	1:G:44:GLU:HG3	2.53	0.43
1:H:88:ARG:HG2	1:H:88:ARG:HH11	1.83	0.43
1:H:79:ILE:HG22	1:H:80:GLY:O	2.18	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:29:ALA:HB2	1:F:44:GLU:HB3	2.00	0.43
1:G:6:LYS:HB3	1:G:107:TYR:CE2	2.53	0.43
1:B:29:ALA:HB2	1:B:44:GLU:HB3	2.00	0.43
1:C:51:GLU:HA	1:C:51:GLU:OE1	2.18	0.43
1:E:140:LEU:HB3	1:E:144:MET:HE3	2.00	0.43
1:H:6:LYS:HA	1:H:9:ASN:HD22	1.83	0.43
1:A:94:GLU:HB3	1:A:131:ILE:HD11	2.01	0.43
1:A:128:GLU:OE2	1:G:1:MET:HE1	2.18	0.43
1:B:32:PHE:CE2	1:B:44:GLU:HG3	2.54	0.43
1:D:88:ARG:HG2	1:D:88:ARG:HH11	1.83	0.43
1:E:32:PHE:CE2	1:E:44:GLU:HG3	2.54	0.43
1:E:79:ILE:HG22	1:E:80:GLY:O	2.18	0.43
1:G:6:LYS:HA	1:G:9:ASN:HD22	1.84	0.43
1:G:79:ILE:HG22	1:G:80:GLY:O	2.18	0.43
1:A:140:LEU:HB3	1:A:144:MET:HE3	1.99	0.43
1:D:79:ILE:HG22	1:D:80:GLY:O	2.19	0.43
1:D:105:ILE:HG23	1:D:117:ARG:HG3	2.01	0.43
1:C:94:GLU:HB3	1:C:131:ILE:HD11	2.00	0.43
1:D:32:PHE:CE2	1:D:44:GLU:HG3	2.53	0.43
1:F:79:ILE:HG22	1:F:80:GLY:O	2.18	0.43
1:A:6:LYS:HA	1:A:9:ASN:HD22	1.84	0.43
1:A:29:ALA:HB2	1:A:44:GLU:HB3	2.00	0.43
1:C:6:LYS:HA	1:C:9:ASN:HD22	1.84	0.43
1:C:32:PHE:CE2	1:C:44:GLU:HG3	2.53	0.43
1:F:6:LYS:HA	1:F:9:ASN:HD22	1.84	0.43
1:F:32:PHE:CE2	1:F:44:GLU:HG3	2.54	0.43
1:A:32:PHE:CE2	1:A:44:GLU:HG3	2.53	0.43
1:G:29:ALA:HB2	1:G:44:GLU:HB3	2.00	0.43
1:A:52:MET:HB3	3:A:200[B]:HEM:C4A	2.54	0.42
1:E:19:LEU:HD21	1:E:71:LEU:HD22	2.01	0.42
1:A:132:ASP:HB2	1:G:64:PHE:CE1	2.54	0.42
1:C:19:LEU:HD21	1:C:71:LEU:HD22	2.01	0.42
1:G:154:ILE:HG13	1:G:155:ARG:HG3	2.02	0.42
1:B:6:LYS:HA	1:B:9:ASN:HD22	1.84	0.42
1:H:19:LEU:HD21	1:H:71:LEU:HD22	2.01	0.42
1:E:6:LYS:HA	1:E:9:ASN:HD22	1.84	0.42
1:E:79:ILE:N	1:E:79:ILE:HD12	2.35	0.42
1:B:18:GLU:O	1:B:22:ILE:HG13	2.20	0.42
1:E:51:GLU:HA	1:E:51:GLU:OE1	2.19	0.42
1:H:29:ALA:HB2	1:H:44:GLU:HB3	2.00	0.42
1:G:19:LEU:HD21	1:G:71:LEU:HD22	2.01	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:79:ILE:N	1:H:79:ILE:HD12	2.34	0.42
1:A:19:LEU:HD21	1:A:71:LEU:HD22	2.02	0.42
1:C:105:ILE:HG23	1:C:117:ARG:HG3	2.02	0.42
1:C:14:LEU:HB3	1:C:58:TYR:CZ	2.55	0.42
1:F:19:LEU:HD21	1:F:71:LEU:HD22	2.02	0.42
1:B:79:ILE:N	1:B:79:ILE:HD12	2.35	0.41
1:E:19:LEU:HD11	1:E:71:LEU:CD2	2.51	0.41
1:F:105:ILE:HG23	1:F:117:ARG:HG3	2.02	0.41
1:A:105:ILE:HG23	1:A:117:ARG:HG3	2.02	0.41
1:E:94:GLU:HB3	1:E:131:ILE:HD11	2.02	0.41
1:F:14:LEU:HB3	1:F:58:TYR:CZ	2.55	0.41
1:A:79:ILE:HD12	1:A:79:ILE:N	2.34	0.41
1:B:98:ALA:O	1:B:102:ARG:HG3	2.21	0.41
1:C:23:ASN:HD22	1:D:71:LEU:HB3	1.85	0.41
1:C:29:ALA:HB2	1:C:44:GLU:HB3	2.00	0.41
1:D:19:LEU:HD21	1:D:71:LEU:HD22	2.02	0.41
1:G:105:ILE:HG23	1:G:117:ARG:HG3	2.02	0.41
1:A:14:LEU:HB3	1:A:58:TYR:CZ	2.55	0.41
1:D:6:LYS:HA	1:D:9:ASN:HD22	1.84	0.41
1:G:52:MET:HB3	3:G:200[B]:HEM:CHB	2.51	0.41
1:D:79:ILE:HD12	1:D:79:ILE:N	2.34	0.41
1:D:111:VAL:HG22	1:D:111:VAL:O	2.21	0.41
1:E:14:LEU:HB3	1:E:58:TYR:CZ	2.55	0.41
1:F:79:ILE:HD12	1:F:79:ILE:N	2.35	0.41
1:H:18:GLU:O	1:H:22:ILE:HG13	2.20	0.41
1:H:105:ILE:HG23	1:H:117:ARG:HG3	2.02	0.41
1:G:79:ILE:N	1:G:79:ILE:HD12	2.35	0.41
1:A:98:ALA:O	1:A:102:ARG:HG3	2.21	0.41
1:B:19:LEU:HD11	1:B:71:LEU:CD2	2.50	0.41
1:C:79:ILE:N	1:C:79:ILE:HD12	2.35	0.41
1:D:14:LEU:HB3	1:D:58:TYR:CZ	2.55	0.41
1:D:43:VAL:HG11	1:D:133:TRP:CE2	2.56	0.41
1:F:19:LEU:HD11	1:F:71:LEU:CD2	2.51	0.41
1:F:98:ALA:O	1:F:102:ARG:HG3	2.21	0.41
1:F:125:ARG:NH2	1:H:122:GLU:OE2	2.53	0.41
1:G:98:ALA:O	1:G:102:ARG:HG3	2.21	0.41
1:G:111:VAL:O	1:G:111:VAL:HG22	2.21	0.41
1:C:19:LEU:HD11	1:C:71:LEU:CD2	2.51	0.41
1:C:65:LEU:HD12	1:C:65:LEU:HA	1.92	0.40
1:A:19:LEU:HD11	1:A:71:LEU:CD2	2.51	0.40
1:D:98:ALA:O	1:D:102:ARG:HG3	2.22	0.40

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:105:ILE:HG23	1:E:117:ARG:HG3	2.02	0.40
1:G:14:LEU:HB3	1:G:58:TYR:CZ	2.56	0.40
1:G:19:LEU:HD11	1:G:71:LEU:CD2	2.52	0.40
1:H:14:LEU:HB3	1:H:58:TYR:CZ	2.56	0.40
1:B:14:LEU:HB3	1:B:58:TYR:CZ	2.56	0.40
1:B:105:ILE:HG23	1:B:117:ARG:HG3	2.03	0.40
1:B:111:VAL:O	1:B:111:VAL:HG22	2.22	0.40
1:D:105:ILE:HG22	1:D:117:ARG:NH1	2.36	0.40
1:C:98:ALA:O	1:C:102:ARG:HG3	2.21	0.40

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

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:88:ARG:NH1	1:G:73:ASP:OD1[3_545]	1.66	0.54
1:E:88:ARG:NH1	1:G:73:ASP:CG[3_545]	1.84	0.36

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	156/158 (99%)	152 (97%)	3 (2%)	1 (1%)	21	38
1	B	156/158 (99%)	152 (97%)	3 (2%)	1 (1%)	21	38
1	C	156/158 (99%)	152 (97%)	3 (2%)	1 (1%)	21	38
1	D	156/158 (99%)	152 (97%)	3 (2%)	1 (1%)	21	38
1	E	156/158 (99%)	152 (97%)	3 (2%)	1 (1%)	21	38
1	F	156/158 (99%)	152 (97%)	3 (2%)	1 (1%)	21	38
1	G	156/158 (99%)	152 (97%)	3 (2%)	1 (1%)	21	38
1	H	156/158 (99%)	152 (97%)	3 (2%)	1 (1%)	21	38

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles
All	All	1248/1264 (99%)	1216 (97%)	24 (2%)	8 (1%)	21 38

All (8) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	79	ILE
1	B	79	ILE
1	C	79	ILE
1	D	79	ILE
1	E	79	ILE
1	F	79	ILE
1	G	79	ILE
1	H	79	ILE

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	139/139 (100%)	133 (96%)	6 (4%)	26 51
1	B	139/139 (100%)	132 (95%)	7 (5%)	22 44
1	C	139/139 (100%)	133 (96%)	6 (4%)	26 51
1	D	139/139 (100%)	133 (96%)	6 (4%)	26 51
1	E	139/139 (100%)	133 (96%)	6 (4%)	26 51
1	F	139/139 (100%)	133 (96%)	6 (4%)	26 51
1	G	139/139 (100%)	133 (96%)	6 (4%)	26 51
1	H	139/139 (100%)	132 (95%)	7 (5%)	22 44
All	All	1112/1112 (100%)	1062 (96%)	50 (4%)	24 49

All (50) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	6	LYS
1	A	11	LEU

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Mol	Chain	Res	Type
1	A	40	LEU
1	A	65	LEU
1	A	76	LYS
1	A	95	LEU
1	B	6	LYS
1	B	11	LEU
1	B	40	LEU
1	B	65	LEU
1	B	76	LYS
1	B	95	LEU
1	B	142	GLN
1	C	6	LYS
1	C	11	LEU
1	C	40	LEU
1	C	65	LEU
1	C	76	LYS
1	C	95	LEU
1	D	6	LYS
1	D	11	LEU
1	D	40	LEU
1	D	65	LEU
1	D	76	LYS
1	D	95	LEU
1	E	6	LYS
1	E	11	LEU
1	E	40	LEU
1	E	65	LEU
1	E	76	LYS
1	E	95	LEU
1	F	6	LYS
1	F	11	LEU
1	F	40	LEU
1	F	65	LEU
1	F	76	LYS
1	F	95	LEU
1	G	6	LYS
1	G	11	LEU
1	G	40	LEU
1	G	65	LEU
1	G	76	LYS
1	G	95	LEU
1	H	6	LYS

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Mol	Chain	Res	Type
1	H	11	LEU
1	H	40	LEU
1	H	65	LEU
1	H	76	LYS
1	H	95	LEU
1	H	154	ILE

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (29) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	ASN
1	A	17	ASN
1	A	100	ASN
1	B	9	ASN
1	B	17	ASN
1	B	100	ASN
1	C	9	ASN
1	C	17	ASN
1	C	46	HIS
1	C	100	ASN
1	C	151	GLN
1	D	9	ASN
1	D	17	ASN
1	D	100	ASN
1	E	9	ASN
1	E	17	ASN
1	E	100	ASN
1	F	9	ASN
1	F	17	ASN
1	F	46	HIS
1	F	100	ASN
1	G	9	ASN
1	G	17	ASN
1	G	100	ASN
1	G	151	GLN
1	H	9	ASN
1	H	17	ASN
1	H	100	ASN
1	H	151	GLN

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no oligosaccharides in this entry.

5.6 Ligand geometry (i)

Of 24 ligands modelled in this entry, 16 are monoatomic - leaving 8 for Mogul analysis.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	HEM	C	200[B]	-	50,50,50	1.86	8 (16%)	67,82,82	1.59	11 (16%)
3	HEM	G	200[B]	-	50,50,50	1.91	7 (14%)	67,82,82	1.48	6 (8%)
3	HEM	A	200[B]	-	50,50,50	2.01	10 (20%)	67,82,82	1.65	10 (14%)
3	HEM	C	200[A]	-	50,50,50	1.88	10 (20%)	67,82,82	1.51	9 (13%)
3	HEM	E	200[B]	-	50,50,50	2.02	8 (16%)	67,82,82	1.51	7 (10%)
3	HEM	G	200[A]	-	50,50,50	1.93	8 (16%)	67,82,82	1.60	12 (17%)
3	HEM	A	200[A]	-	50,50,50	2.00	10 (20%)	67,82,82	1.64	13 (19%)
3	HEM	E	200[A]	-	50,50,50	2.03	7 (14%)	67,82,82	1.49	7 (10%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	HEM	C	200[B]	-	-	6/14/54/54	-
3	HEM	G	200[B]	-	-	6/14/54/54	-
3	HEM	A	200[B]	-	-	6/14/54/54	-
3	HEM	C	200[A]	-	-	8/14/54/54	-
3	HEM	E	200[B]	-	-	8/14/54/54	-
3	HEM	G	200[A]	-	-	8/14/54/54	-
3	HEM	A	200[A]	-	-	8/14/54/54	-
3	HEM	E	200[A]	-	-	8/14/54/54	-

All (68) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	E	200[B]	HEM	C3D-C2D	8.41	1.54	1.36
3	E	200[A]	HEM	C3D-C2D	8.25	1.54	1.36
3	A	200[A]	HEM	C3D-C2D	8.10	1.54	1.36
3	A	200[B]	HEM	C3D-C2D	8.01	1.54	1.36
3	G	200[B]	HEM	C3D-C2D	7.97	1.53	1.36
3	G	200[A]	HEM	C3D-C2D	7.90	1.53	1.36
3	C	200[A]	HEM	C3D-C2D	7.86	1.53	1.36
3	C	200[B]	HEM	C3D-C2D	7.65	1.53	1.36
3	A	200[B]	HEM	FE-NB	6.28	2.14	1.94
3	E	200[A]	HEM	FE-NA	6.17	2.15	1.95
3	E	200[B]	HEM	FE-NA	5.93	2.14	1.95
3	G	200[A]	HEM	FE-NA	5.47	2.13	1.95
3	G	200[B]	HEM	FE-ND	5.29	2.11	1.94
3	E	200[A]	HEM	FE-NB	5.21	2.10	1.94
3	G	200[A]	HEM	FE-ND	5.14	2.10	1.94
3	A	200[B]	HEM	FE-NC	5.06	2.11	1.95
3	C	200[A]	HEM	FE-NB	4.90	2.10	1.94
3	C	200[B]	HEM	FE-NB	4.88	2.09	1.94
3	E	200[B]	HEM	FE-ND	4.84	2.09	1.94
3	A	200[A]	HEM	FE-NA	4.79	2.11	1.95
3	A	200[A]	HEM	FE-NB	4.56	2.08	1.94
3	G	200[B]	HEM	FE-NA	4.34	2.09	1.95
3	A	200[A]	HEM	FE-ND	4.15	2.07	1.94
3	E	200[A]	HEM	FE-ND	4.10	2.07	1.94
3	E	200[B]	HEM	FE-NB	4.01	2.07	1.94
3	A	200[A]	HEM	FE-NC	3.84	2.07	1.95
3	C	200[A]	HEM	FE-NC	3.76	2.07	1.95
3	C	200[B]	HEM	FE-NC	3.61	2.07	1.95
3	G	200[A]	HEM	FE-NB	3.54	2.05	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	C	200[A]	HEM	FE-NA	3.53	2.06	1.95
3	G	200[B]	HEM	FE-NB	3.52	2.05	1.94
3	C	200[B]	HEM	FE-NA	3.37	2.06	1.95
3	C	200[B]	HEM	FE-ND	3.06	2.04	1.94
3	C	200[A]	HEM	FE-ND	3.06	2.04	1.94
3	G	200[B]	HEM	CAC-C3C	2.87	1.55	1.47
3	G	200[A]	HEM	CAB-C3B	2.85	1.55	1.47
3	C	200[B]	HEM	CAC-C3C	2.82	1.54	1.47
3	E	200[A]	HEM	CAB-C3B	2.79	1.54	1.47
3	E	200[A]	HEM	CAC-C3C	2.74	1.54	1.47
3	A	200[B]	HEM	CAC-C3C	2.71	1.54	1.47
3	A	200[B]	HEM	CAB-C3B	2.70	1.54	1.47
3	A	200[A]	HEM	CAB-C3B	2.67	1.54	1.47
3	C	200[B]	HEM	CAB-C3B	2.65	1.54	1.47
3	G	200[A]	HEM	CAC-C3C	2.65	1.54	1.47
3	E	200[B]	HEM	CAC-C3C	2.62	1.54	1.47
3	G	200[B]	HEM	CAB-C3B	2.61	1.54	1.47
3	C	200[A]	HEM	CAB-C3B	2.56	1.54	1.47
3	C	200[B]	HEM	CMA-C3A	2.48	1.55	1.50
3	A	200[B]	HEM	FE-ND	2.48	2.02	1.94
3	A	200[A]	HEM	CAC-C3C	2.42	1.53	1.47
3	C	200[A]	HEM	CMD-C2D	2.36	1.55	1.50
3	E	200[B]	HEM	CAB-C3B	2.34	1.53	1.47
3	C	200[A]	HEM	CAC-C3C	2.31	1.53	1.47
3	A	200[A]	HEM	CMA-C3A	2.29	1.55	1.50
3	E	200[B]	HEM	CMD-C2D	2.28	1.55	1.50
3	A	200[B]	HEM	CMA-C3A	2.27	1.55	1.50
3	A	200[B]	HEM	FE-NA	2.23	2.02	1.95
3	A	200[B]	HEM	C3B-C2B	-2.19	1.32	1.37
3	E	200[A]	HEM	CMA-C3A	2.14	1.55	1.50
3	A	200[B]	HEM	CMD-C2D	2.13	1.55	1.50
3	A	200[A]	HEM	CMD-C2D	2.13	1.55	1.50
3	E	200[B]	HEM	CMB-C2B	2.10	1.55	1.50
3	G	200[B]	HEM	CMC-C2C	2.07	1.55	1.50
3	G	200[A]	HEM	CMB-C2B	2.06	1.55	1.50
3	G	200[A]	HEM	C2A-C3A	-2.03	1.33	1.38
3	A	200[A]	HEM	CMC-C2C	2.03	1.54	1.50
3	C	200[A]	HEM	C2A-C3A	-2.03	1.33	1.38
3	C	200[A]	HEM	CMB-C2B	2.01	1.54	1.50

All (75) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	A	200[B]	HEM	C4D-ND-C1D	7.57	114.18	105.21
3	E	200[A]	HEM	C4D-ND-C1D	6.81	113.27	105.21
3	C	200[B]	HEM	C4D-ND-C1D	6.79	113.25	105.21
3	G	200[B]	HEM	C4D-ND-C1D	6.66	113.10	105.21
3	A	200[A]	HEM	C4D-ND-C1D	6.56	112.97	105.21
3	C	200[A]	HEM	C4D-ND-C1D	6.21	112.56	105.21
3	E	200[B]	HEM	C4D-ND-C1D	6.06	112.38	105.21
3	G	200[A]	HEM	C4D-ND-C1D	5.73	111.99	105.21
3	A	200[B]	HEM	C3B-C2B-C1B	4.04	109.45	106.41
3	C	200[B]	HEM	C3B-C2B-C1B	3.78	109.25	106.41
3	E	200[B]	HEM	C3B-C4B-NB	-3.66	106.84	109.47
3	C	200[B]	HEM	CAD-CBD-CGD	-3.15	105.30	113.67
3	E	200[B]	HEM	C1B-NB-C4B	3.15	108.94	105.21
3	C	200[B]	HEM	C3D-C4D-ND	-3.15	106.72	110.17
3	A	200[A]	HEM	C3B-C4B-NB	-3.12	107.23	109.47
3	A	200[B]	HEM	C3D-C4D-ND	-3.04	106.83	110.17
3	A	200[A]	HEM	C4C-C3C-C2C	2.99	109.40	106.81
3	C	200[B]	HEM	CHA-C4D-ND	2.96	128.03	124.37
3	G	200[A]	HEM	C2A-C1A-NA	-2.87	106.96	110.15
3	C	200[A]	HEM	C2A-C1A-NA	-2.80	107.04	110.15
3	A	200[A]	HEM	C1B-NB-C4B	2.74	108.46	105.21
3	A	200[A]	HEM	C2A-C1A-NA	-2.73	107.12	110.15
3	A	200[B]	HEM	CHA-C4D-ND	2.70	127.71	124.37
3	E	200[A]	HEM	C3B-C2B-C1B	2.69	108.43	106.41
3	G	200[A]	HEM	C1B-NB-C4B	2.68	108.38	105.21
3	G	200[A]	HEM	C3B-C4B-NB	-2.67	107.55	109.47
3	E	200[A]	HEM	CHD-C4C-NC	2.67	127.36	124.45
3	G	200[A]	HEM	CAA-CBA-CGA	-2.57	106.84	113.67
3	C	200[A]	HEM	C4C-C3C-C2C	2.53	109.01	106.81
3	G	200[A]	HEM	CAD-C3D-C4D	2.51	129.08	124.70
3	E	200[A]	HEM	C4C-C3C-C2C	2.48	108.96	106.81
3	C	200[B]	HEM	C1D-C2D-C3D	-2.48	104.37	106.98
3	C	200[A]	HEM	C1D-C2D-C3D	-2.45	104.40	106.98
3	A	200[B]	HEM	CHD-C1D-ND	2.44	127.05	124.42
3	E	200[B]	HEM	C2B-C1B-NB	-2.44	107.04	109.84
3	C	200[B]	HEM	CHC-C4B-NB	2.42	127.03	124.42
3	C	200[A]	HEM	CHC-C4B-NB	2.41	127.02	124.42
3	A	200[B]	HEM	C2B-C1B-NB	-2.41	107.08	109.84
3	C	200[B]	HEM	C2B-C1B-NB	-2.41	107.08	109.84
3	G	200[B]	HEM	CHD-C1D-ND	2.40	127.00	124.42
3	E	200[B]	HEM	CHC-C4B-NB	2.40	127.00	124.42
3	A	200[B]	HEM	CAD-CBD-CGD	-2.35	107.42	113.67
3	G	200[A]	HEM	C3B-C2B-C1B	2.32	108.16	106.41

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	G	200[B]	HEM	CAD-CBD-CGD	-2.29	107.60	113.67
3	C	200[A]	HEM	C3B-C4B-NB	-2.28	107.83	109.47
3	G	200[A]	HEM	CHC-C4B-NB	2.27	126.87	124.42
3	A	200[B]	HEM	C2A-C1A-NA	-2.26	107.65	110.15
3	G	200[A]	HEM	O1A-CGA-CBA	-2.24	115.97	123.09
3	G	200[B]	HEM	C3D-C4D-ND	-2.24	107.71	110.17
3	E	200[B]	HEM	C3B-C2B-C1B	2.24	108.09	106.41
3	C	200[A]	HEM	C3D-C4D-ND	-2.23	107.72	110.17
3	G	200[A]	HEM	CMD-C2D-C1D	2.23	128.52	125.03
3	E	200[A]	HEM	C2D-C1D-ND	-2.22	107.34	109.90
3	A	200[A]	HEM	CHA-C4D-ND	2.21	127.11	124.37
3	E	200[B]	HEM	CHB-C1B-NB	2.19	127.07	124.37
3	G	200[A]	HEM	CBC-CAC-C3C	-2.19	116.60	127.53
3	A	200[A]	HEM	CHC-C4B-NB	2.18	126.77	124.42
3	A	200[B]	HEM	C2D-C1D-ND	-2.17	107.39	109.90
3	G	200[A]	HEM	C4C-C3C-C2C	2.16	108.69	106.81
3	C	200[A]	HEM	CBD-CAD-C3D	-2.15	106.58	112.53
3	E	200[A]	HEM	C4C-NC-C1C	2.15	109.32	105.82
3	C	200[B]	HEM	C1B-NB-C4B	2.15	107.75	105.21
3	E	200[A]	HEM	CHD-C1D-ND	2.14	126.72	124.42
3	G	200[B]	HEM	CHA-C4D-ND	2.12	126.99	124.37
3	A	200[A]	HEM	C3B-C2B-C1B	2.10	107.99	106.41
3	C	200[B]	HEM	CHD-C1D-ND	2.10	126.68	124.42
3	C	200[A]	HEM	CMD-C2D-C1D	2.08	128.29	125.03
3	A	200[A]	HEM	CHD-C1D-ND	2.08	126.66	124.42
3	A	200[A]	HEM	CAD-CBD-CGD	-2.07	108.16	113.67
3	G	200[B]	HEM	C4C-C3C-C2C	2.06	108.60	106.81
3	A	200[A]	HEM	C3D-C4D-ND	-2.06	107.92	110.17
3	A	200[B]	HEM	C1B-NB-C4B	2.05	107.63	105.21
3	C	200[B]	HEM	C2A-C1A-NA	-2.04	107.89	110.15
3	A	200[A]	HEM	CBC-CAC-C3C	-2.02	117.44	127.53
3	A	200[A]	HEM	CMD-C2D-C1D	2.02	128.18	125.03

There are no chirality outliers.

All (58) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	200[A]	HEM	C2B-C3B-CAB-CBB
3	A	200[A]	HEM	C2C-C3C-CAC-CBC
3	A	200[A]	HEM	C4C-C3C-CAC-CBC
3	A	200[B]	HEM	C2B-C3B-CAB-CBB
3	A	200[B]	HEM	C4B-C3B-CAB-CBB

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Mol	Chain	Res	Type	Atoms
3	A	200[B]	HEM	C2C-C3C-CAC-CBC
3	C	200[A]	HEM	C2B-C3B-CAB-CBB
3	C	200[A]	HEM	C4B-C3B-CAB-CBB
3	C	200[A]	HEM	C2C-C3C-CAC-CBC
3	C	200[A]	HEM	C4C-C3C-CAC-CBC
3	C	200[B]	HEM	C2B-C3B-CAB-CBB
3	C	200[B]	HEM	C4B-C3B-CAB-CBB
3	C	200[B]	HEM	C2C-C3C-CAC-CBC
3	E	200[A]	HEM	C2B-C3B-CAB-CBB
3	E	200[A]	HEM	C4B-C3B-CAB-CBB
3	E	200[A]	HEM	C2C-C3C-CAC-CBC
3	E	200[B]	HEM	C2B-C3B-CAB-CBB
3	E	200[B]	HEM	C2C-C3C-CAC-CBC
3	E	200[B]	HEM	C4C-C3C-CAC-CBC
3	G	200[A]	HEM	C2B-C3B-CAB-CBB
3	G	200[A]	HEM	C4B-C3B-CAB-CBB
3	G	200[A]	HEM	C2C-C3C-CAC-CBC
3	G	200[A]	HEM	C4C-C3C-CAC-CBC
3	G	200[B]	HEM	C2B-C3B-CAB-CBB
3	G	200[B]	HEM	C4B-C3B-CAB-CBB
3	G	200[B]	HEM	C2C-C3C-CAC-CBC
3	G	200[B]	HEM	C4C-C3C-CAC-CBC
3	A	200[A]	HEM	C4B-C3B-CAB-CBB
3	C	200[B]	HEM	C4C-C3C-CAC-CBC
3	E	200[A]	HEM	C4C-C3C-CAC-CBC
3	E	200[B]	HEM	C4B-C3B-CAB-CBB
3	A	200[B]	HEM	C4C-C3C-CAC-CBC
3	E	200[A]	HEM	CAD-CBD-CGD-O2D
3	C	200[B]	HEM	CAD-CBD-CGD-O1D
3	C	200[A]	HEM	CAD-CBD-CGD-O1D
3	E	200[A]	HEM	CAD-CBD-CGD-O1D
3	E	200[B]	HEM	CAD-CBD-CGD-O1D
3	G	200[B]	HEM	CAD-CBD-CGD-O2D
3	G	200[B]	HEM	CAD-CBD-CGD-O1D
3	A	200[B]	HEM	CAD-CBD-CGD-O1D
3	C	200[A]	HEM	CAD-CBD-CGD-O2D
3	A	200[A]	HEM	CAD-CBD-CGD-O2D
3	A	200[B]	HEM	CAD-CBD-CGD-O2D
3	C	200[B]	HEM	CAD-CBD-CGD-O2D
3	E	200[A]	HEM	CAA-CBA-CGA-O1A
3	E	200[A]	HEM	CAA-CBA-CGA-O2A
3	E	200[B]	HEM	CAD-CBD-CGD-O2D

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Mol	Chain	Res	Type	Atoms
3	E	200[B]	HEM	CAA-CBA-CGA-O2A
3	A	200[A]	HEM	CAD-CBD-CGD-O1D
3	G	200[A]	HEM	CAA-CBA-CGA-O2A
3	C	200[A]	HEM	CAA-CBA-CGA-O2A
3	E	200[B]	HEM	CAA-CBA-CGA-O1A
3	C	200[A]	HEM	CAA-CBA-CGA-O1A
3	G	200[A]	HEM	CAA-CBA-CGA-O1A
3	G	200[A]	HEM	CAD-CBD-CGD-O2D
3	G	200[A]	HEM	CAD-CBD-CGD-O1D
3	A	200[A]	HEM	CAA-CBA-CGA-O2A
3	A	200[A]	HEM	CAA-CBA-CGA-O1A

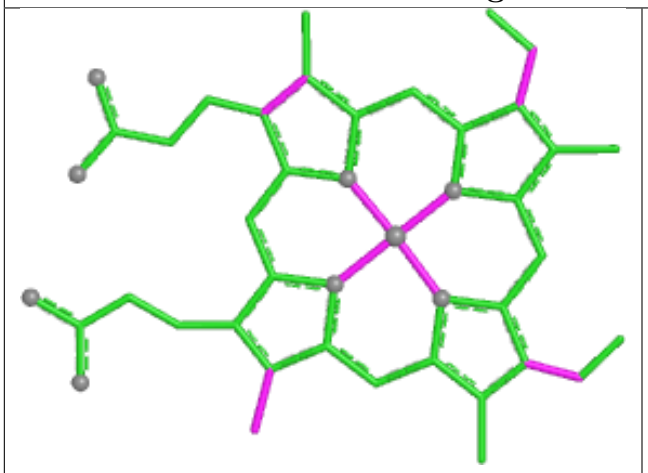
There are no ring outliers.

3 monomers are involved in 7 short contacts:

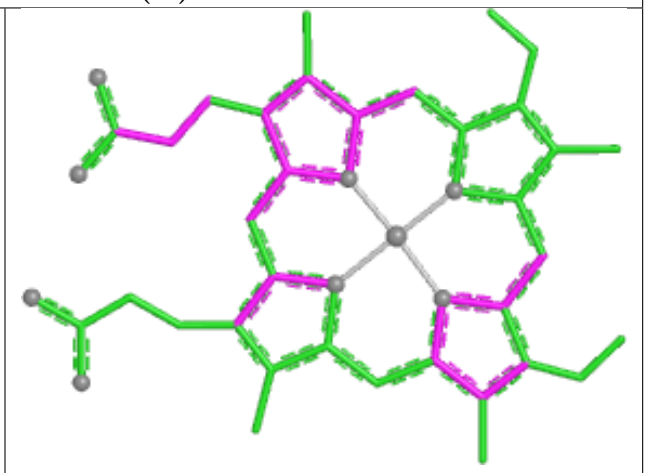
Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	G	200[B]	HEM	2	0
3	A	200[B]	HEM	3	0
3	A	200[A]	HEM	2	0

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.

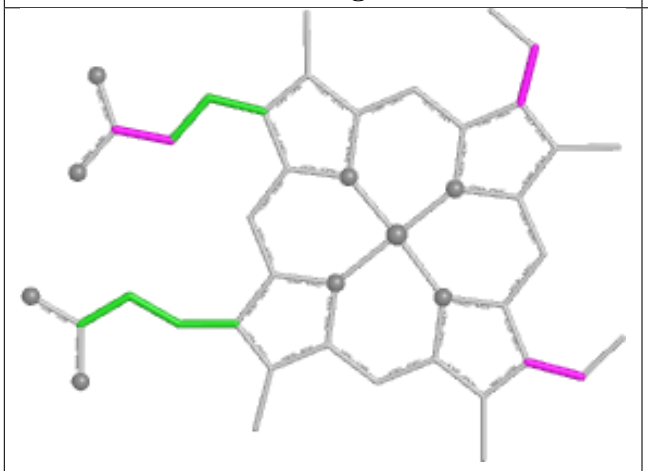
Ligand HEM C 200 (B)



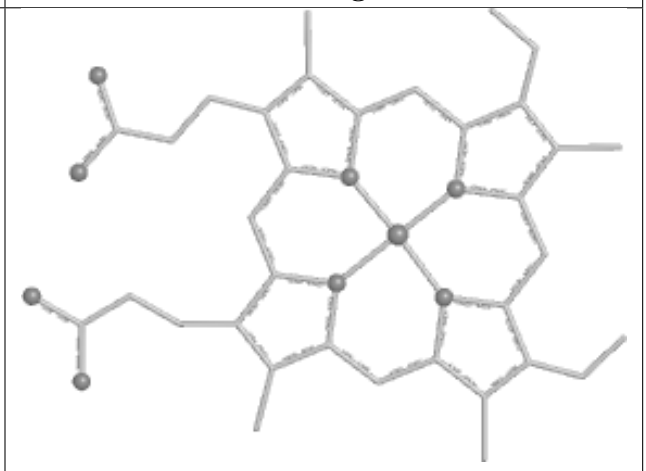
Bond lengths



Bond angles

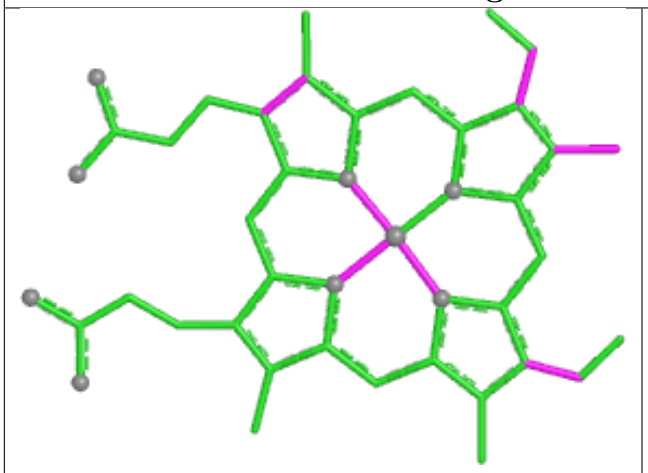


Torsions

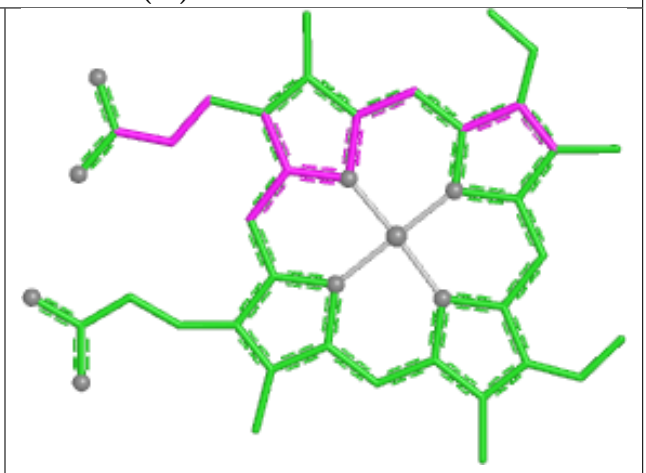


Rings

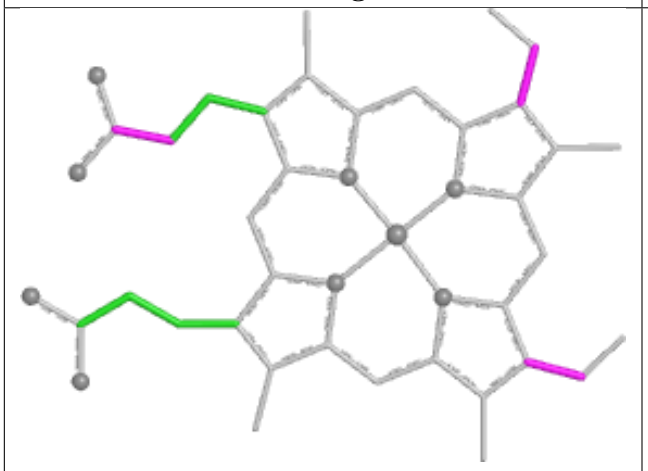
Ligand HEM G 200 (B)



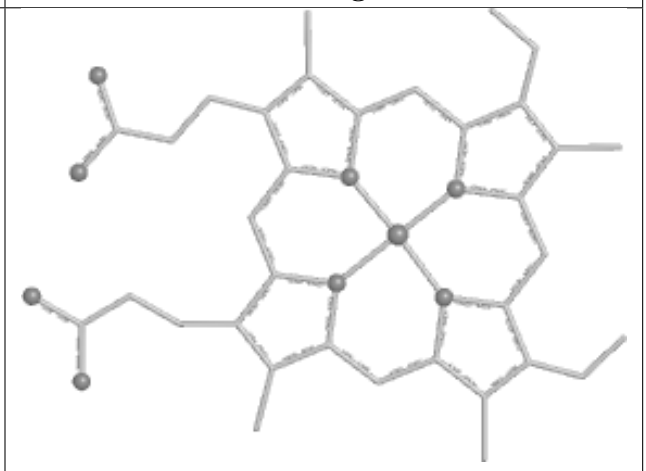
Bond lengths



Bond angles

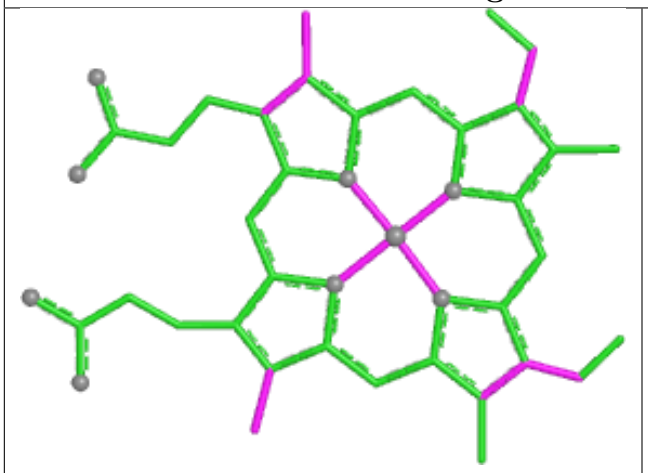


Torsions

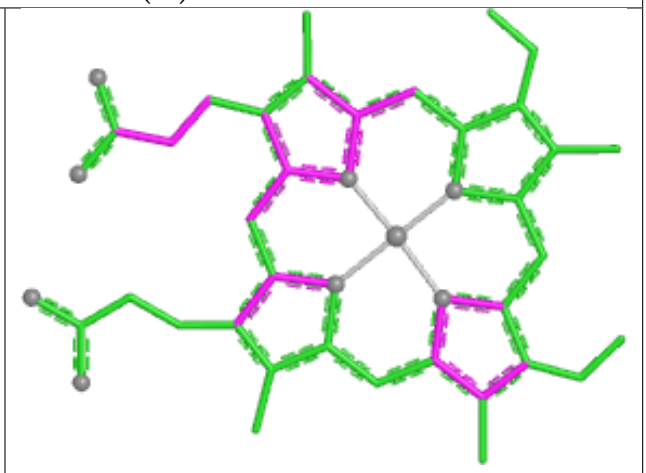


Rings

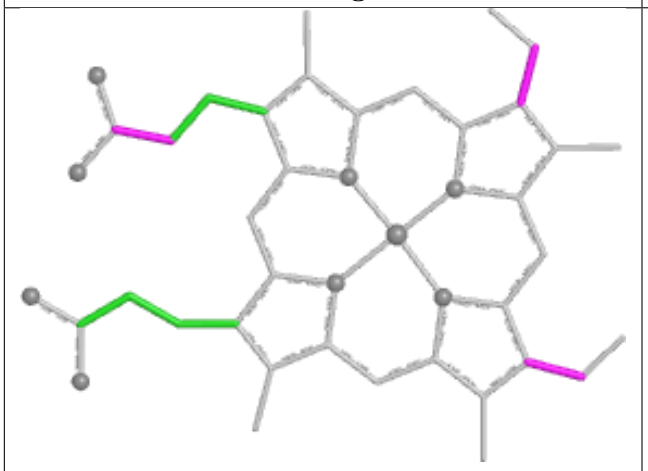
Ligand HEM A 200 (B)



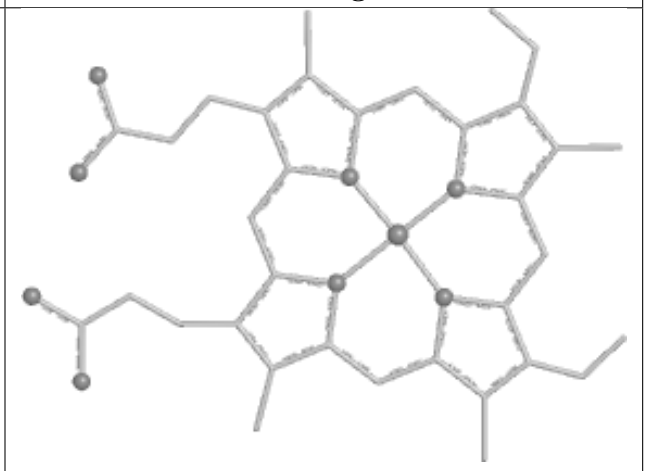
Bond lengths



Bond angles

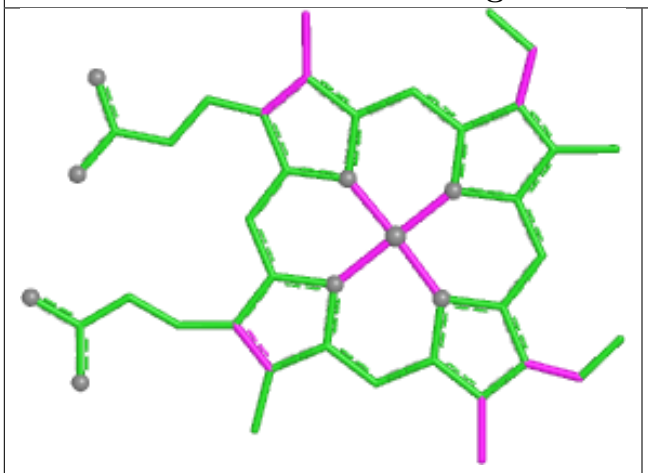


Torsions

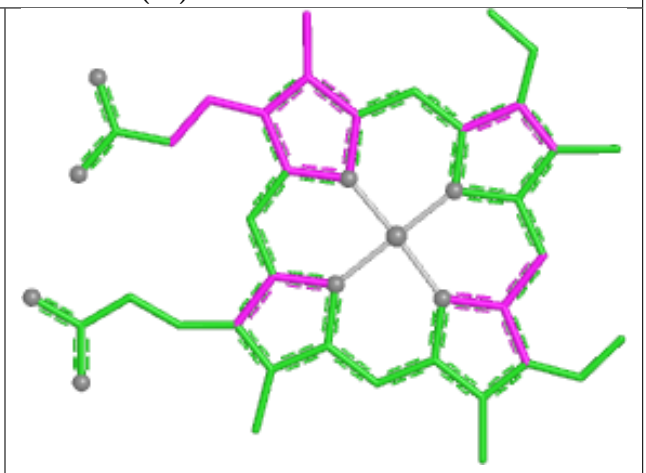


Rings

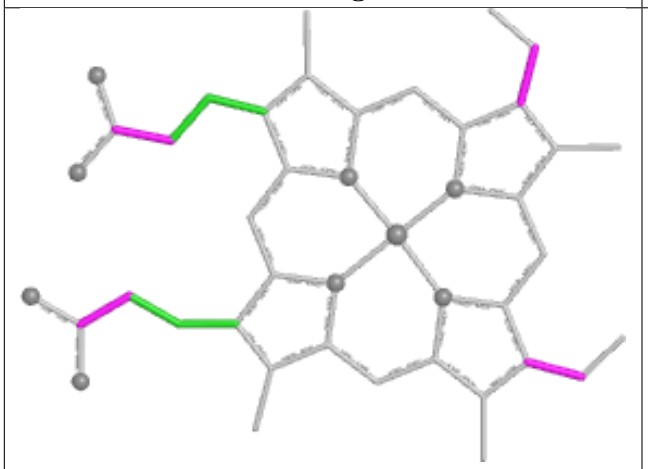
Ligand HEM C 200 (A)



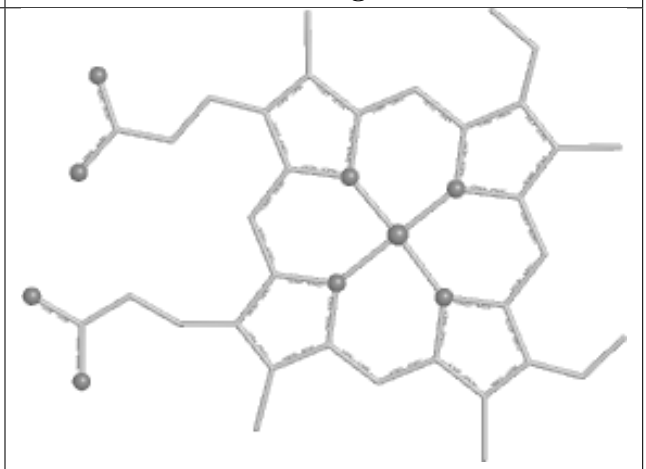
Bond lengths



Bond angles

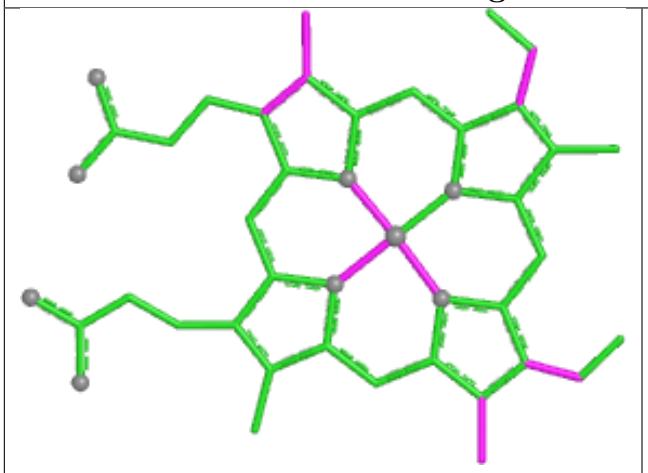


Torsions

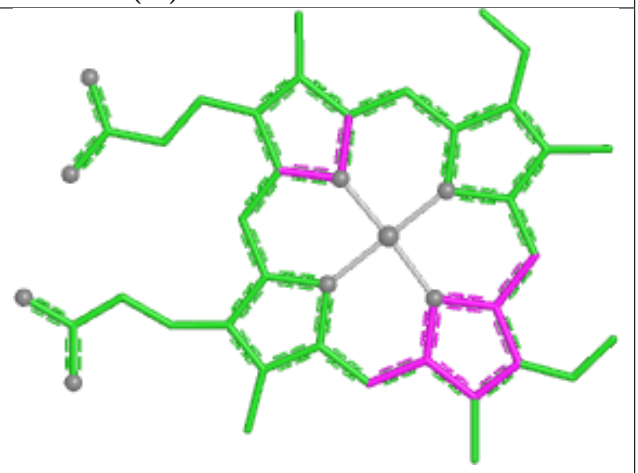


Rings

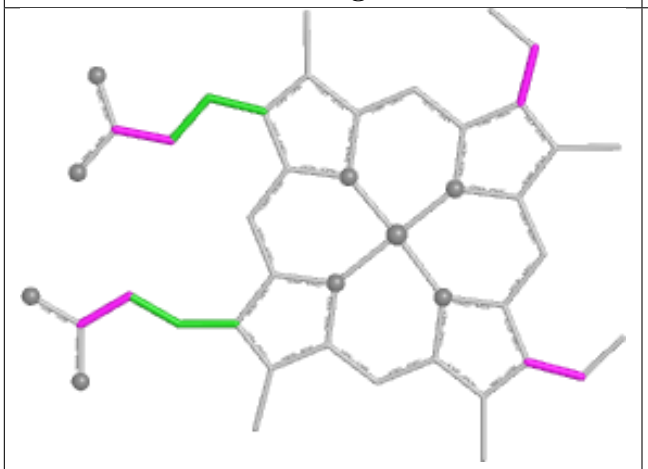
Ligand HEM E 200 (B)



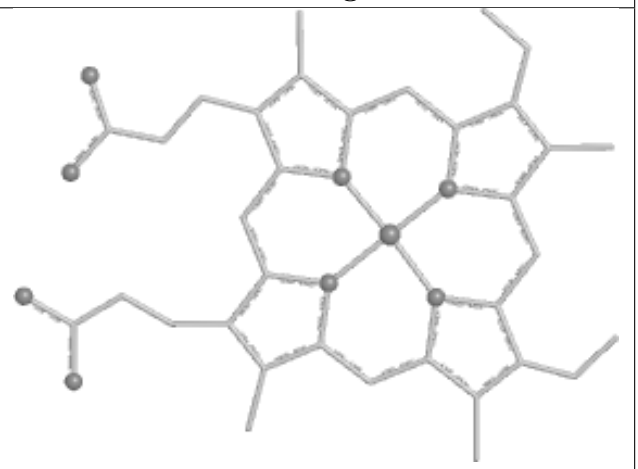
Bond lengths



Bond angles

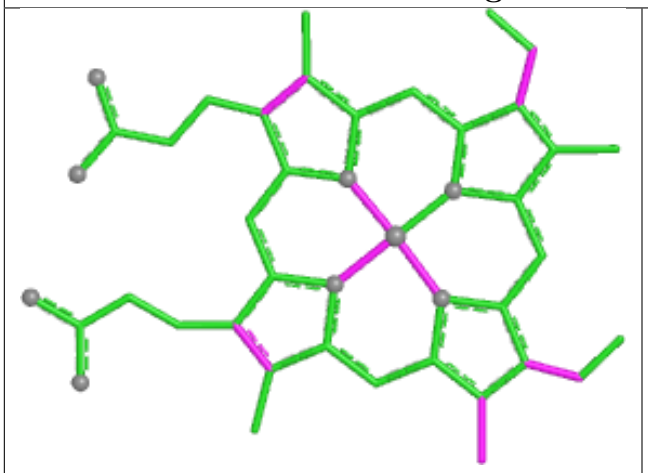


Torsions

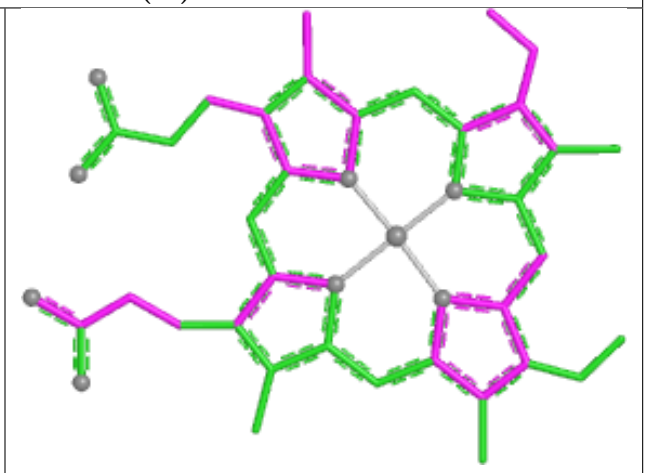


Rings

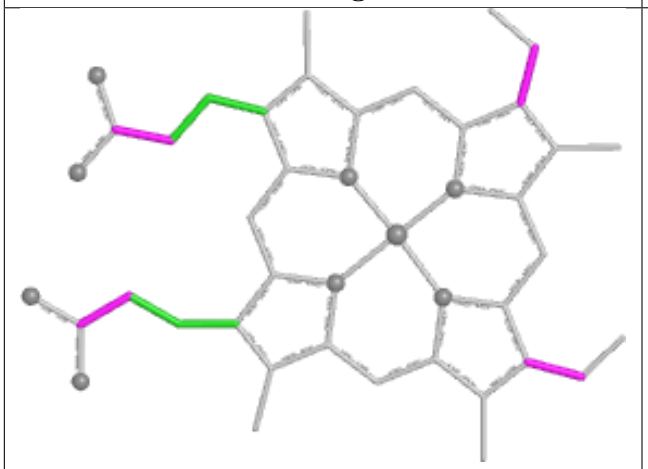
Ligand HEM G 200 (A)



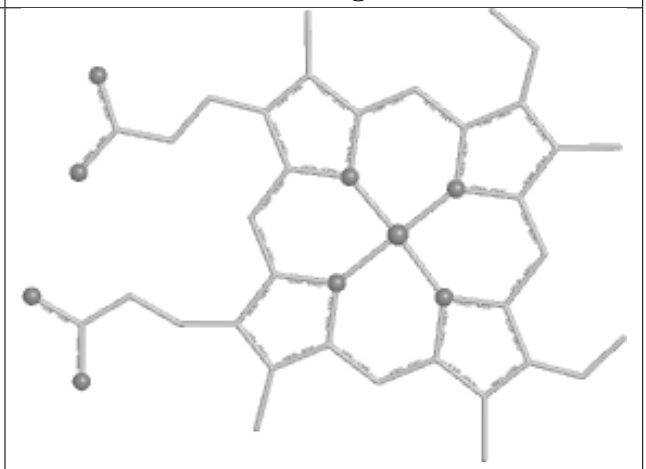
Bond lengths



Bond angles

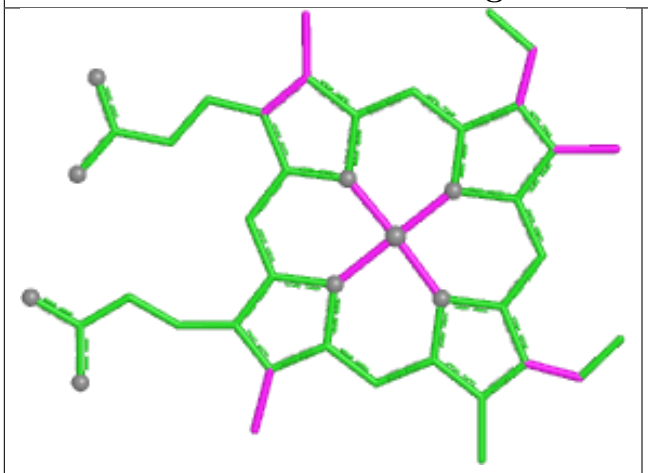


Torsions

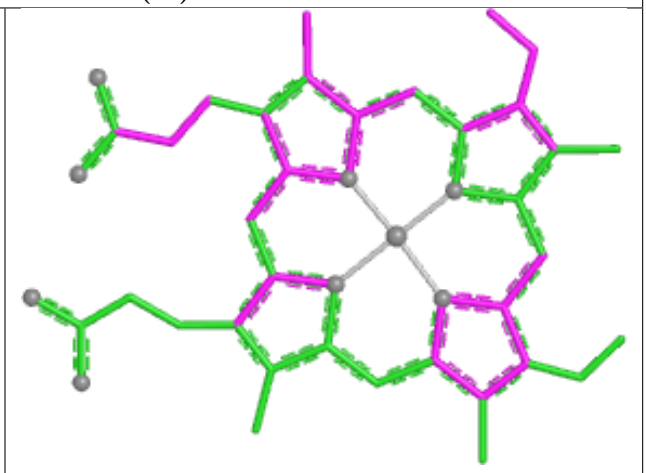


Rings

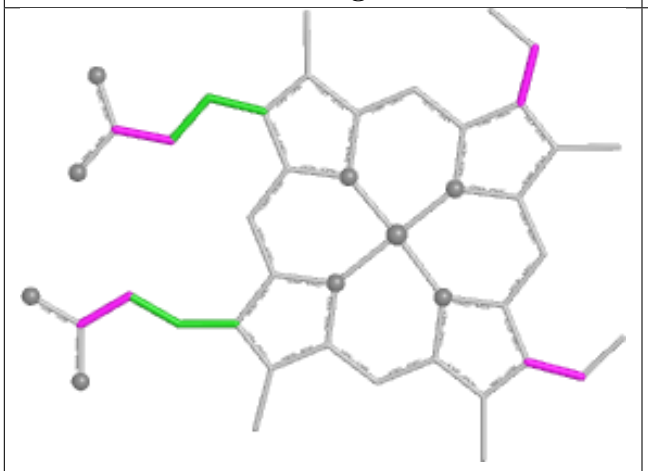
Ligand HEM A 200 (A)



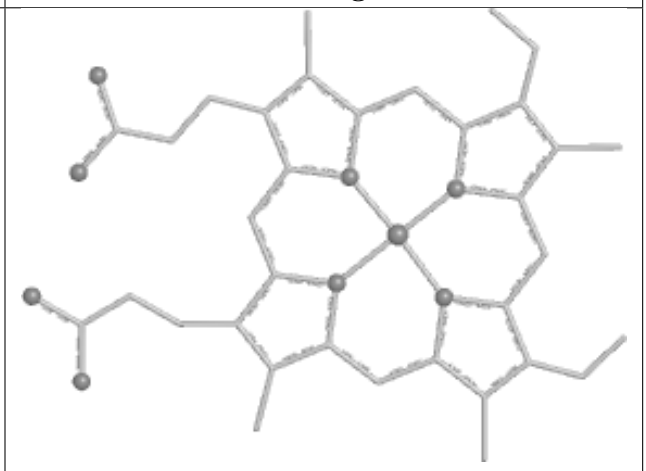
Bond lengths



Bond angles

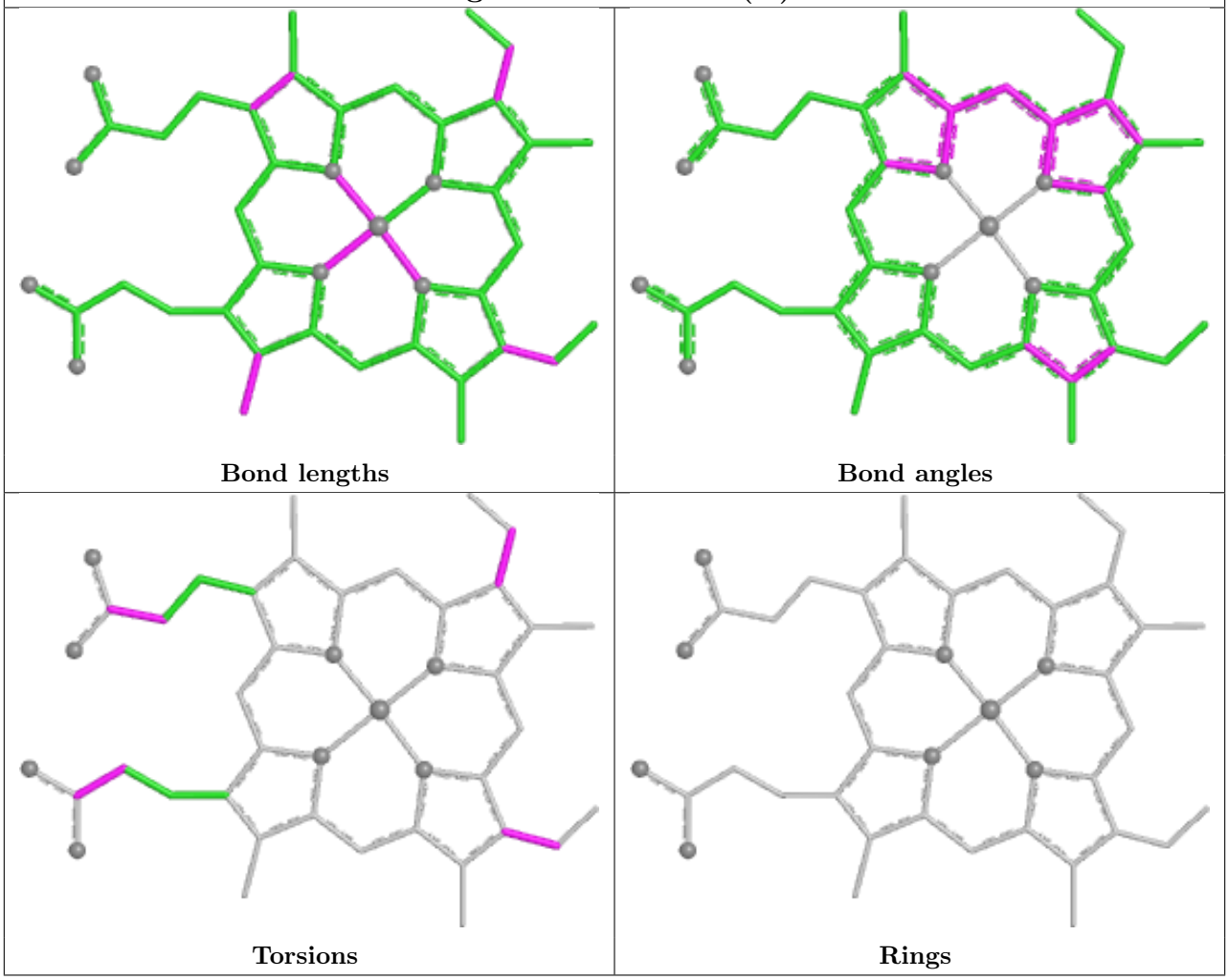


Torsions



Rings

Ligand HEM E 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 (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	158/158 (100%)	-1.62	0	100	100	14, 24, 41, 65	0
1	B	158/158 (100%)	-1.64	0	100	100	8, 23, 40, 63	0
1	C	158/158 (100%)	-1.68	0	100	100	14, 24, 40, 66	0
1	D	158/158 (100%)	-1.63	0	100	100	13, 23, 41, 63	0
1	E	158/158 (100%)	-1.64	0	100	100	14, 24, 41, 65	0
1	F	158/158 (100%)	-1.63	0	100	100	13, 24, 40, 65	0
1	G	158/158 (100%)	-1.66	0	100	100	13, 23, 40, 79	0
1	H	158/158 (100%)	-1.64	0	100	100	13, 23, 43, 75	0
All	All	1264/1264 (100%)	-1.64	0	100	100	8, 23, 42, 79	0

There are no RSRZ outliers to report.

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 oligosaccharides in this entry.

6.4 Ligands (i)

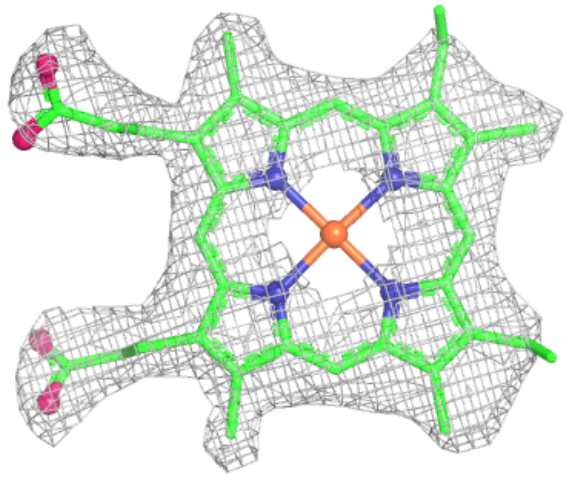
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

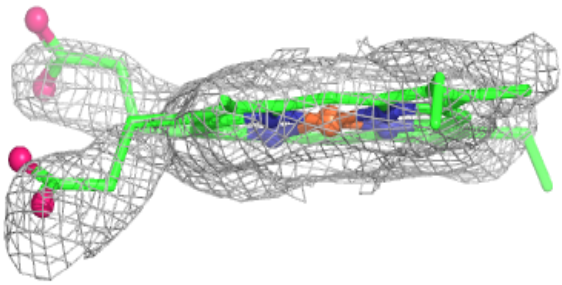
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
2	FE	A	301	1/1	1.00	0.00	23,23,23,23	0
2	FE	A	302	1/1	1.00	0.01	44,44,44,44	1
2	FE	B	301	1/1	1.00	0.01	30,30,30,30	0
2	FE	B	302	1/1	1.00	0.03	47,47,47,47	1
2	FE	C	301	1/1	1.00	0.01	27,27,27,27	0
2	FE	C	302	1/1	1.00	0.02	37,37,37,37	1
2	FE	D	301	1/1	1.00	0.01	26,26,26,26	0
2	FE	D	302	1/1	1.00	0.02	36,36,36,36	1
2	FE	E	301	1/1	1.00	0.01	27,27,27,27	0
2	FE	E	302	1/1	1.00	0.01	37,37,37,37	1
2	FE	F	301	1/1	1.00	0.00	29,29,29,29	0
2	FE	F	302	1/1	1.00	0.01	48,48,48,48	1
2	FE	G	301	1/1	1.00	0.01	28,28,28,28	0
2	FE	G	302	1/1	1.00	0.01	40,40,40,40	1
2	FE	H	301	1/1	1.00	0.01	28,28,28,28	0
2	FE	H	302	1/1	1.00	0.02	34,34,34,34	1
3	HEM	A	200[A]	43/43	1.00	0.04	16,48,76,100	42
3	HEM	A	200[B]	43/43	1.00	0.04	2,12,30,65	42
3	HEM	C	200[A]	43/43	1.00	0.04	2,24,62,81	42
3	HEM	C	200[B]	43/43	1.00	0.04	2,19,33,49	42
3	HEM	E	200[A]	43/43	1.00	0.02	7,25,46,55	42
3	HEM	E	200[B]	43/43	1.00	0.02	2,23,42,79	42
3	HEM	G	200[A]	43/43	1.00	0.03	5,31,57,64	42
3	HEM	G	200[B]	43/43	1.00	0.03	2,20,39,60	42

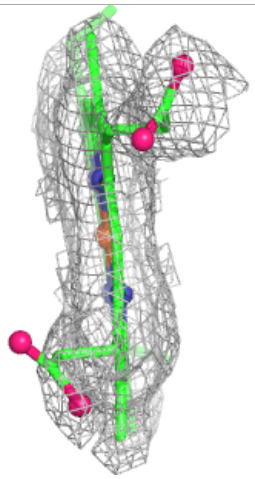
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around HEM A 200 (A):

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

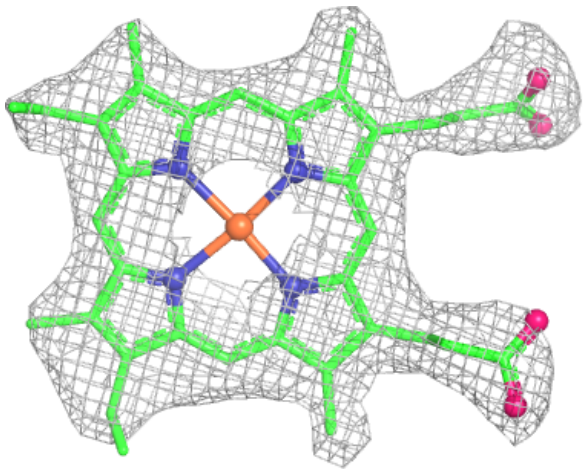


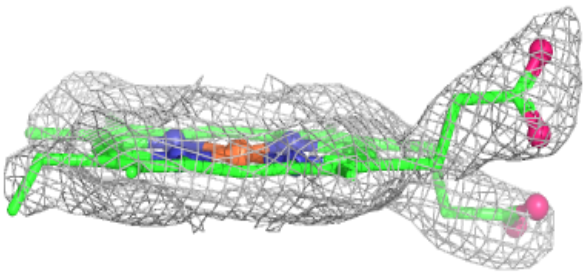


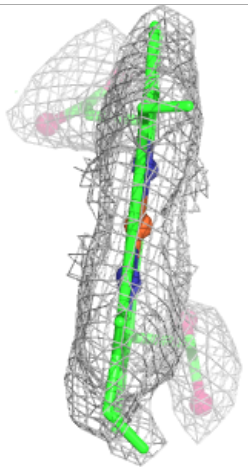


Electron density around HEM A 200 (B):

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

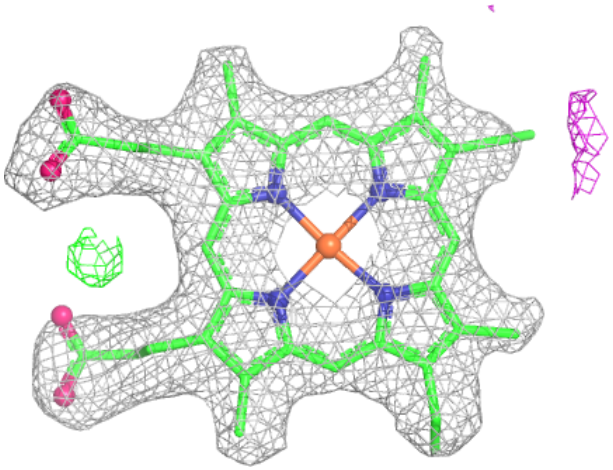


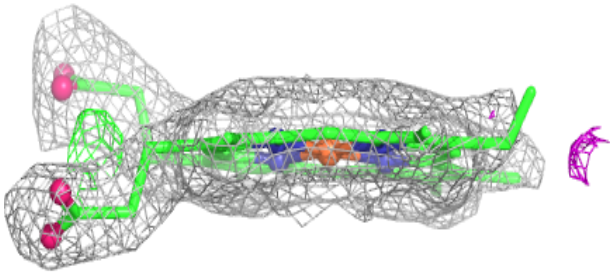


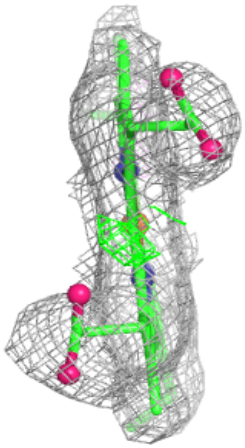


Electron density around HEM C 200 (A):

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

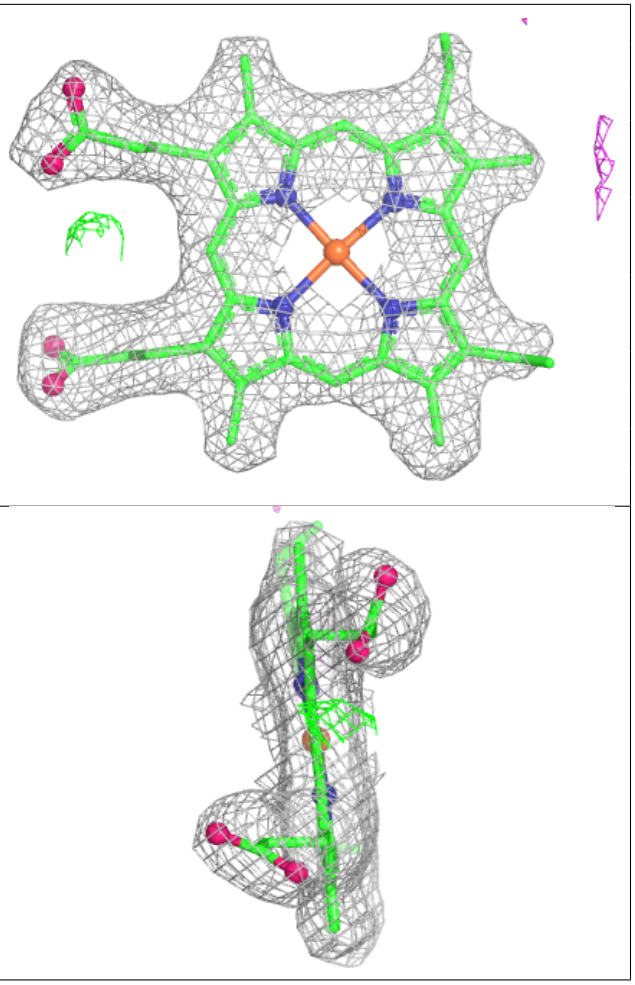






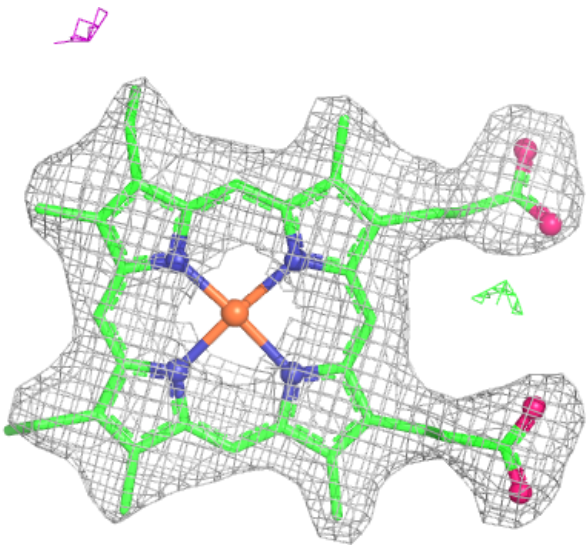
Electron density around HEM C 200 (B):

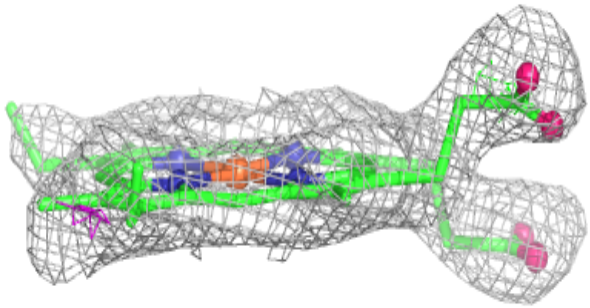
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

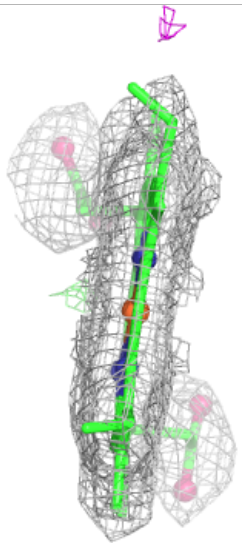


Electron density around HEM E 200 (A):

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

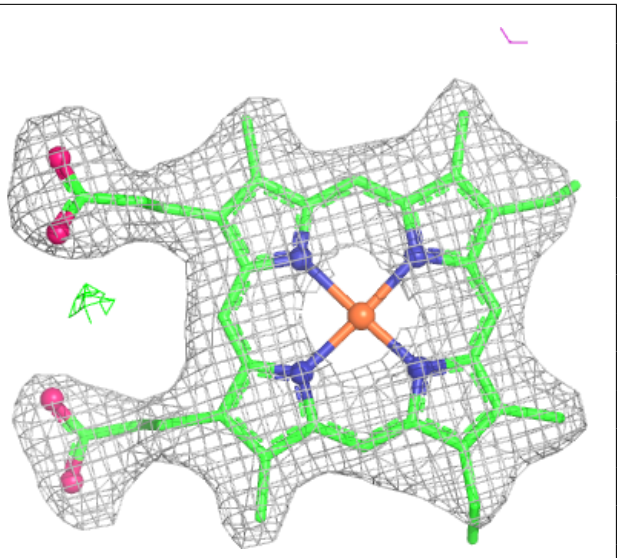


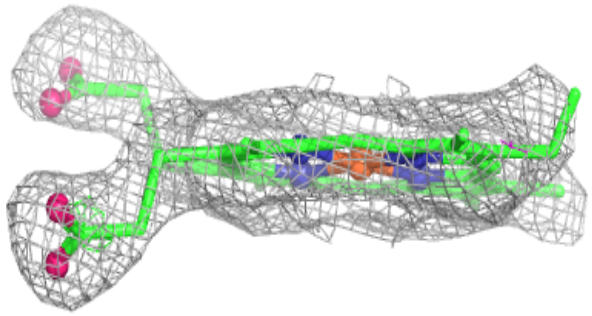


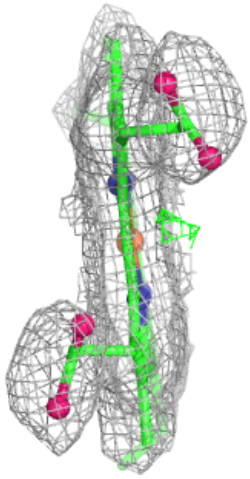


Electron density around HEM E 200 (B):

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

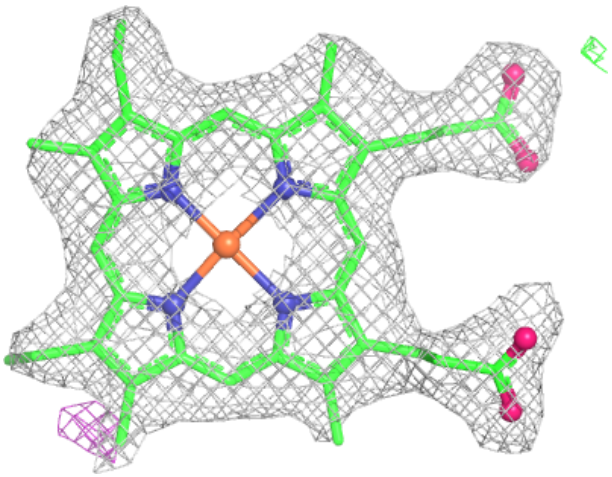


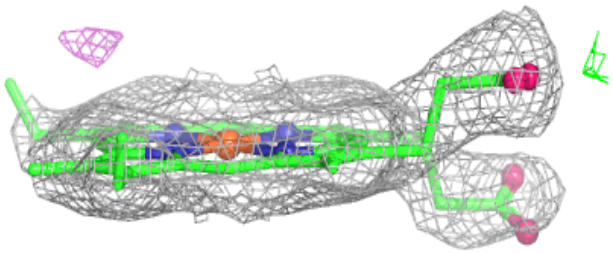


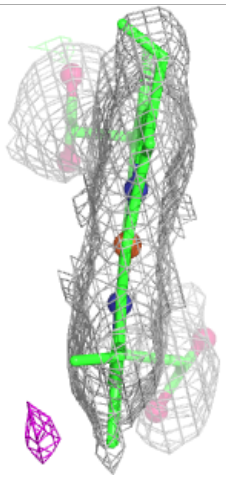


Electron density around HEM G 200 (A):

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

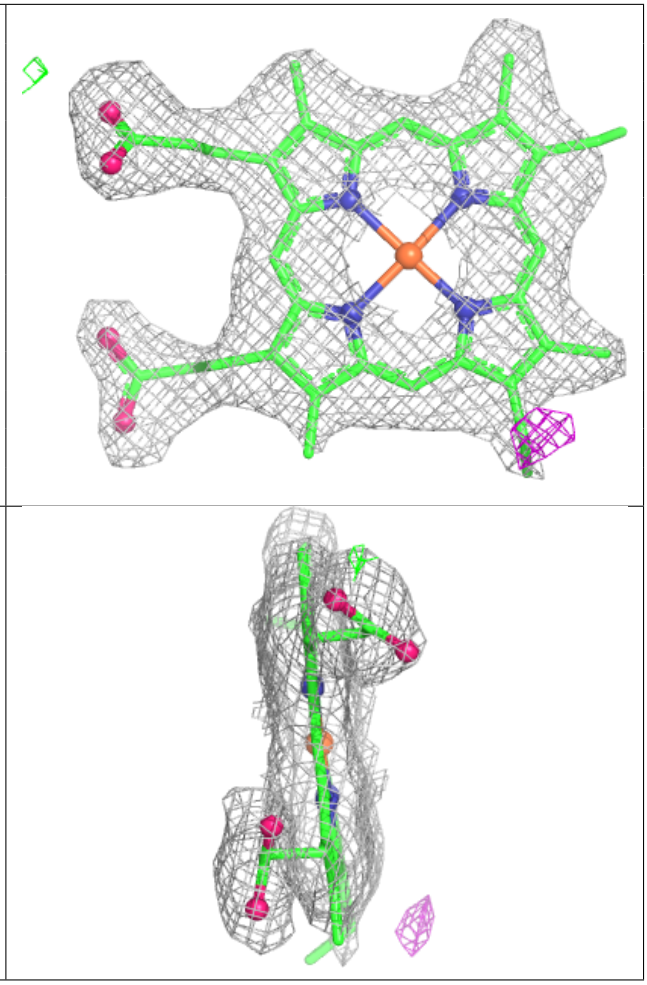






Electron density around HEM G 200 (B):

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



6.5 Other polymers (i)

There are no such residues in this entry.