Full wwPDB X-ray Structure Validation Report (i)

Aug 22, 2023 – 09:47 AM EDT

PDB ID	:	2QRW
Title	:	Crystal structure of Mycobacterium tuberculosis trHbO WG8F mutant
Authors	:	Milani, M.; Bolognesi, M.
Deposited on	:	2007-07-30
Resolution	:	1.93 Å(reported)

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

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

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

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

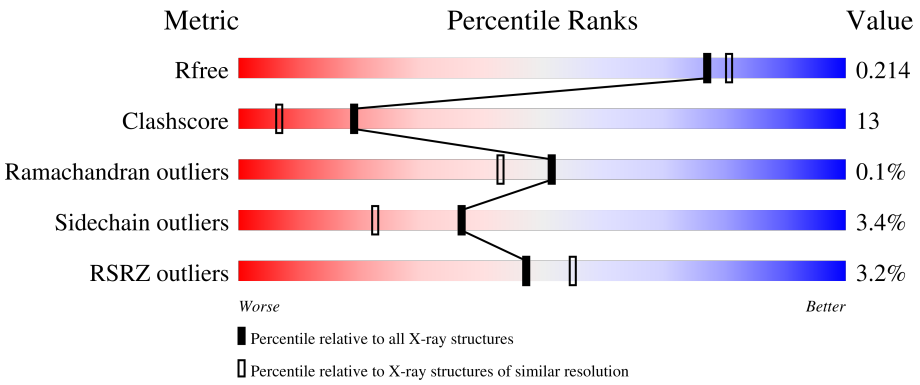
MolProbity	:	4.02b-467
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.35
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.35

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 1.93 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	4310 (1.96-1.92)
Clashscore	141614	1023 (1.94-1.94)
Ramachandran outliers	138981	1007 (1.94-1.94)
Sidechain outliers	138945	1007 (1.94-1.94)
RSRZ outliers	127900	4250 (1.96-1.92)

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	128	3% 74% 23% • •
1	B	128	2% 82% 15% • •
1	C	128	% 90% 9% • •
1	D	128	3% 78% 17% • •
1	E	128	5% 80% 16% • •

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Mol	Chain	Length	Quality of chain
1	F	128	5% 74% 22%
1	G	128	% 82% 15%
1	H	128	2% 80% 17%
1	I	128	83% 16%
1	J	128	4% 78% 20%
1	K	128	6% 80% 18%
1	L	128	8% 72% 25%

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

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
3	SO4	E	802	-	-	-	X
3	SO4	H	801	-	-	X	-

2 Entry composition (i)

There are 5 unique types of molecules in this entry. The entry contains 15001 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 Hemoglobin-like protein HbO.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	126	1049	657	192	195	5	0	2	0
1	B	127	1042	654	189	194	5	8	0	0
1	C	127	1048	657	190	196	5	7	1	0
1	D	126	1044	654	190	195	5	0	1	0
1	E	126	1044	654	190	195	5	0	1	0
1	F	126	1052	659	192	196	5	0	2	0
1	G	127	1062	665	192	199	6	0	3	0
1	H	127	1050	659	190	195	6	0	1	0
1	I	127	1064	666	194	199	5	0	3	0
1	J	126	1057	662	195	195	5	0	3	0
1	K	127	1075	674	198	198	5	0	4	0
1	L	126	1063	664	197	197	5	0	3	0

There are 12 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	88	PHE	TRP	engineered mutation	UNP P0A595
B	88	PHE	TRP	engineered mutation	UNP P0A595
C	88	PHE	TRP	engineered mutation	UNP P0A595
D	88	PHE	TRP	engineered mutation	UNP P0A595
E	88	PHE	TRP	engineered mutation	UNP P0A595

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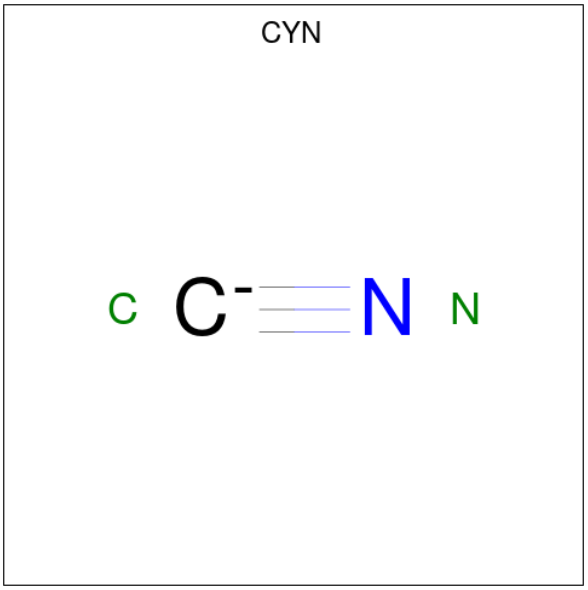
Chain	Residue	Modelled	Actual	Comment	Reference
F	88	PHE	TRP	engineered mutation	UNP P0A595
G	88	PHE	TRP	engineered mutation	UNP P0A595
H	88	PHE	TRP	engineered mutation	UNP P0A595
I	88	PHE	TRP	engineered mutation	UNP P0A595
J	88	PHE	TRP	engineered mutation	UNP P0A595
K	88	PHE	TRP	engineered mutation	UNP P0A595
L	88	PHE	TRP	engineered mutation	UNP P0A595

- Molecule 2 is CYANIDE ION (three-letter code: CYN) (formula: CN).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N			
2	A	1	2	1	1		0	0
2	B	1	2	1	1		0	0
2	C	1	2	1	1		0	0
2	D	1	2	1	1		0	0
2	E	1	2	1	1		0	0
2	F	1	2	1	1		0	0
2	G	1	2	1	1		0	0
2	H	1	2	1	1		0	0

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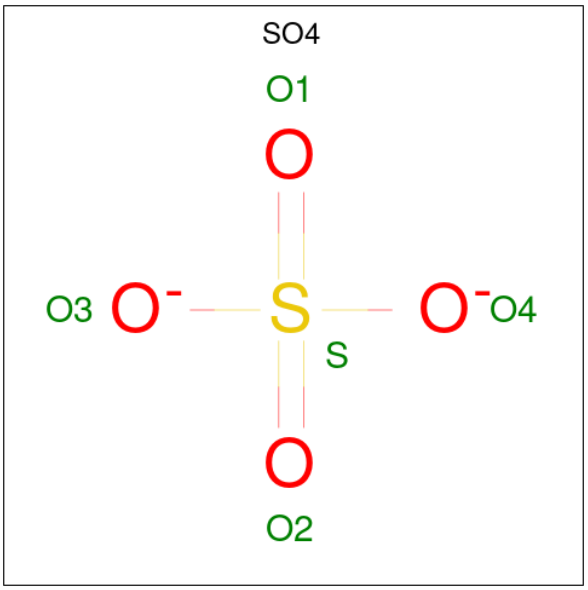
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	N		
2	I	1	2	1	1	0	0
			Total	C	N		
2	J	1	2	1	1	0	0
			Total	C	N		
2	K	1	2	1	1	0	0
			Total	C	N		
2	L	1	2	1	1	0	0

- Molecule 3 is SULFATE ION (three-letter code: SO4) (formula: O_4S).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
3	A	1	5	4	1	0	0
			Total	O	S		
3	A	1	5	4	1	0	0
			Total	O	S		
3	A	1	5	4	1	0	0
			Total	O	S		
3	A	1	5	4	1	0	0
			Total	O	S		
3	B	1	5	4	1	0	0
			Total	O	S		
3	B	1	5	4	1	0	0
			Total	O	S		
3	B	1	5	4	1	0	0

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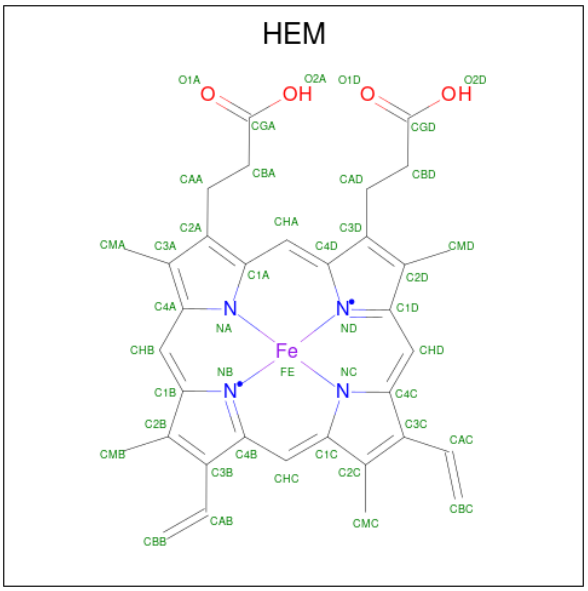
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	O	S		
3	D	1	5	4	1	0	0
			Total	O	S		
3	D	1	5	4	1	0	0
			Total	O	S		
3	E	1	5	4	1	0	0
			Total	O	S		
3	E	1	5	4	1	0	0
			Total	O	S		
3	F	1	5	4	1	0	0
			Total	O	S		
3	F	1	5	4	1	0	0
			Total	O	S		
3	F	1	5	4	1	0	0
			Total	O	S		
3	G	1	5	4	1	0	0
			Total	O	S		
3	G	1	5	4	1	0	0
			Total	O	S		
3	H	1	5	4	1	0	0
			Total	O	S		
3	I	1	5	4	1	0	0
			Total	O	S		
3	J	1	5	4	1	0	0
			Total	O	S		
3	K	1	5	4	1	0	0
			Total	O	S		
3	L	1	5	4	1	0	0
			Total	O	S		
3	L	1	5	4	1	0	0
			Total	O	S		
3	L	1	5	4	1	0	0

- Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
4	A	1	43	34	1	4	4	0	0
4	B	1	43	34	1	4	4	0	0
4	C	1	43	34	1	4	4	0	0
4	D	1	43	34	1	4	4	0	0
4	E	1	43	34	1	4	4	0	0
4	F	1	43	34	1	4	4	0	0
4	G	1	43	34	1	4	4	0	0
4	H	1	43	34	1	4	4	0	0
4	I	1	43	34	1	4	4	0	0
4	J	1	43	34	1	4	4	0	0
4	K	1	43	34	1	4	4	0	0
4	L	1	43	34	1	4	4	0	0

- Molecule 5 is water.

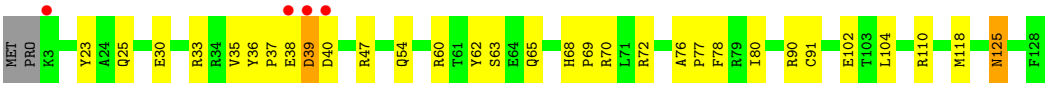
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
5	A	145	145	145	0	0
			Total	O		
5	B	129	129	129	0	0
			Total	O		
5	C	155	155	155	0	0
			Total	O		
5	D	135	135	135	0	0
			Total	O		
5	E	117	117	117	0	0
			Total	O		
5	F	137	137	137	0	0
			Total	O		
5	G	172	172	172	0	0
			Total	O		
5	H	131	131	131	0	0
			Total	O		
5	I	176	176	176	0	0
			Total	O		
5	J	123	123	123	0	0
			Total	O		
5	K	143	143	143	0	0
			Total	O		
5	L	133	133	133	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.

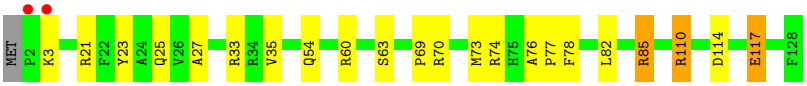
• Molecule 1: Hemoglobin-like protein HbO

Chain A:



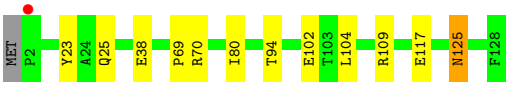
• Molecule 1: Hemoglobin-like protein HbO

Chain B:



• Molecule 1: Hemoglobin-like protein HbO

Chain C:

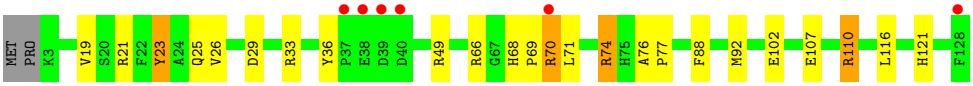


• Molecule 1: Hemoglobin-like protein HbO

Chain D:

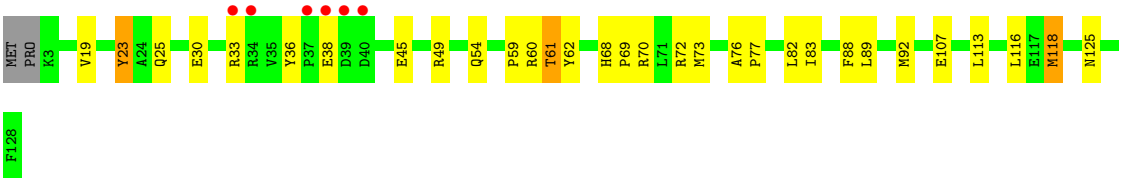
• Molecule 1: Hemoglobin-like protein HbO

Chain E:



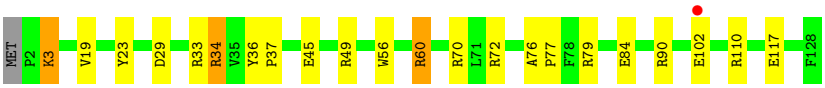
• Molecule 1: Hemoglobin-like protein HbO

Chain F: 5% 74% 22%



• Molecule 1: Hemoglobin-like protein HbO

Chain G: % 82% 15%

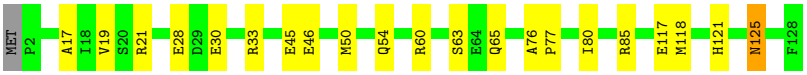


• Molecule 1: Hemoglobin-like protein HbO

Chain H: 2% 80% 17%

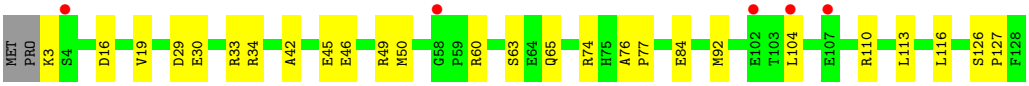
• Molecule 1: Hemoglobin-like protein HbO

Chain I: 83% 16%



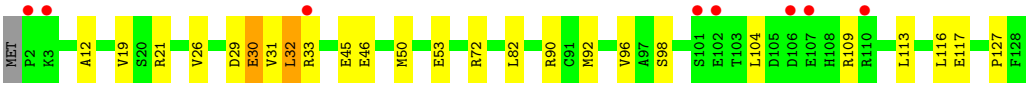
• Molecule 1: Hemoglobin-like protein HbO

Chain J: 4% 78% 20%



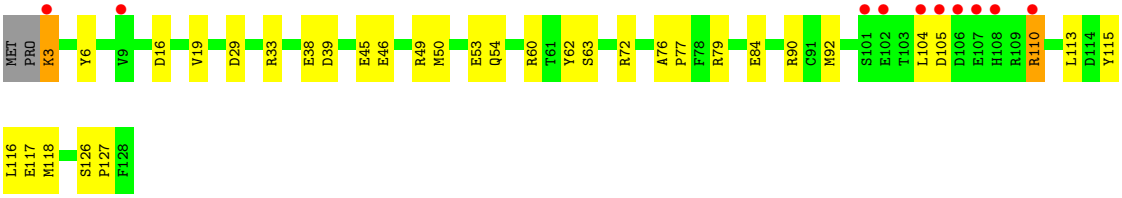
• Molecule 1: Hemoglobin-like protein HbO

Chain K: 6% 80% 18%



• Molecule 1: Hemoglobin-like protein HbO

Chain L: 8% 72% 25%


MET
PRO
K3
Y6
V9
D16
V19
D29
R33
E38
D39
E45
E46
R49
M50
E53
Q54
R60
T61
Y62
S63
R72
A76
P77
F78
R79
E84
R90
C91
M92
S101
E102
T103
L104
D105
D106
E107
H108
R109
R110
L113
D114
Y115

L116
E117
M118
S126
P127
F128

4 Data and refinement statistics (i)

Property	Value	Source
Space group	I 41 2 2	Depositor
Cell constants a, b, c, α, β, γ	187.31Å 187.31Å 274.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	40.00 – 1.93 47.65 – 1.93	Depositor EDS
% Data completeness (in resolution range)	98.7 (40.00-1.93) 98.7 (47.65-1.93)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	2.23 (at 1.92Å)	Xtriage
Refinement program	REFMAC 5.2.0019	Depositor
R, R_{free}	0.176 , 0.215 0.176 , 0.214	Depositor DCC
R_{free} test set	8925 reflections (5.00%)	wwPDB-VP
Wilson B-factor (Å^2)	26.2	Xtriage
Anisotropy	0.035	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 58.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.49$, $< L^2 > = 0.33$	Xtriage
Estimated twinning fraction	0.011 for -1/2*h-1/2*k-1/2*l,-1/2*h-1/2*k+1/2*l,-h+k 0.008 for -1/2*h+1/2*k-1/2*l,1/2*h-1/2*k-1/2*l,-h-k	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	15001	wwPDB-VP
Average B, all atoms (Å^2)	28.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 2.32% 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: SO4, CYN, HEM

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	0.53	0/1076	0.60	0/1449
1	B	0.47	0/1067	0.58	1/1438 (0.1%)
1	C	0.54	0/1073	0.60	0/1446
1	D	0.48	0/1068	0.57	0/1439
1	E	0.44	0/1068	0.55	0/1439
1	F	0.51	0/1079	0.63	0/1453
1	G	0.52	0/1087	0.56	0/1465
1	H	0.47	0/1075	0.58	0/1448
1	I	0.55	0/1090	0.58	0/1470
1	J	0.44	0/1087	0.55	0/1464
1	K	0.45	0/1103	0.53	0/1486
1	L	0.40	0/1087	0.57	0/1464
All	All	0.49	0/12960	0.57	1/17461 (0.0%)

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

Mol	Chain	#Chirality outliers	#Planarity outliers
1	E	0	1

There are no bond length outliers.

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	B	85	ARG	NE-CZ-NH1	5.21	122.90	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	E	70	ARG	Peptide

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	1049	0	1006	50	0
1	B	1042	0	997	17	0
1	C	1048	0	1001	13	0
1	D	1044	0	996	39	0
1	E	1044	0	997	34	0
1	F	1052	0	1008	37	0
1	G	1062	0	1015	31	0
1	H	1050	0	1006	32	0
1	I	1064	0	1013	22	0
1	J	1057	0	1021	21	0
1	K	1075	0	1038	24	0
1	L	1063	0	1019	38	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	1	0
2	F	2	0	0	0	0
2	G	2	0	0	0	0
2	H	2	0	0	0	0
2	I	2	0	0	0	0
2	J	2	0	0	0	0
2	K	2	0	0	0	0
2	L	2	0	0	0	0
3	A	20	0	0	1	0
3	B	15	0	0	1	0
3	D	10	0	0	0	0
3	E	10	0	0	0	0
3	F	15	0	0	0	0
3	G	10	0	0	2	0
3	H	5	0	0	6	0
3	I	5	0	0	0	0
3	J	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	K	5	0	0	0	0
3	L	15	0	0	0	0
4	A	43	0	30	2	0
4	B	43	0	30	0	0
4	C	43	0	30	2	0
4	D	43	0	30	0	0
4	E	43	0	30	2	0
4	F	43	0	30	3	0
4	G	43	0	30	2	0
4	H	43	0	30	1	0
4	I	43	0	30	2	0
4	J	43	0	30	3	0
4	K	43	0	30	3	0
4	L	43	0	30	3	0
5	A	145	0	0	15	0
5	B	129	0	0	2	0
5	C	155	0	0	2	0
5	D	135	0	0	7	0
5	E	117	0	0	11	0
5	F	137	0	0	16	0
5	G	172	0	0	16	0
5	H	131	0	0	7	0
5	I	176	0	0	6	0
5	J	123	0	0	5	0
5	K	143	0	0	3	0
5	L	133	0	0	11	0
All	All	15001	0	12477	335	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 13.

All (335) 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:D:23:TYR:OH	1:D:36:TYR:CE2	1.65	1.48
1:D:60:ARG:HD2	5:D:925:HOH:O	1.30	1.27
1:A:72:ARG:HH22	1:L:79:ARG:CG	1.47	1.26
1:A:72:ARG:NH2	1:L:79:ARG:HG2	1.51	1.23
1:F:33:ARG:HD3	5:F:919:HOH:O	1.36	1.23
1:C:117:GLU:HB2	5:C:933:HOH:O	1.41	1.21
1:A:69:PRO:HG2	5:G:862:HOH:O	1.36	1.20

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:F:23:TYR:OH	1:F:36:TYR:HE2	0.87	1.19
1:L:49:ARG:HD2	5:L:837:HOH:O	1.43	1.16
1:F:30:GLU:HG2	5:F:936:HOH:O	1.42	1.16
1:E:23:TYR:OH	1:E:36:TYR:HE2	0.81	1.14
1:E:110:ARG:HH11	1:E:110:ARG:HG2	1.10	1.14
1:K:96:VAL:HG11	1:K:113[B]:LEU:HD11	1.28	1.13
1:D:23:TYR:OH	1:D:36:TYR:HE2	0.79	1.12
1:H:72:ARG:HD3	5:H:851:HOH:O	1.45	1.12
1:B:117:GLU:HG2	5:E:860:HOH:O	1.49	1.11
1:J:34:ARG:HB2	5:J:896:HOH:O	1.54	1.08
1:F:23:TYR:OH	1:F:36:TYR:CE2	1.70	1.07
1:D:117:GLU:HG2	5:D:908:HOH:O	1.56	1.06
1:L:72:ARG:HD3	5:L:812:HOH:O	1.55	1.04
1:A:72:ARG:HH22	1:L:79:ARG:HG2	1.09	1.02
1:B:25:GLN:NE2	5:B:919:HOH:O	1.92	1.01
1:L:117:GLU:HG3	5:L:887:HOH:O	1.62	0.97
1:H:16:ASP:HB2	5:H:880:HOH:O	1.69	0.93
1:D:60:ARG:HD3	1:L:90:ARG:CZ	1.99	0.92
1:E:23:TYR:CZ	1:E:36:TYR:CE2	2.57	0.92
1:F:107:GLU:HG3	5:F:831:HOH:O	1.68	0.92
1:A:118:MET:CE	1:D:82:LEU:HD21	2.00	0.91
1:F:61:THR:HG21	5:F:841:HOH:O	1.69	0.90
1:G:110:ARG:HD2	5:G:962:HOH:O	1.72	0.90
1:A:118:MET:CE	1:D:82:LEU:CD2	2.50	0.90
1:E:69:PRO:HA	5:E:892:HOH:O	1.73	0.89
4:I:700:HEM:HMC1	4:I:700:HEM:HBC2	1.55	0.88
1:A:72:ARG:NH2	1:L:79:ARG:CG	2.21	0.88
1:C:80:ILE:H	1:C:125:ASN:HD21	1.22	0.88
1:A:25:GLN:HG2	5:A:913:HOH:O	1.73	0.87
1:D:46:GLU:HG2	5:D:931:HOH:O	1.74	0.87
1:E:110:ARG:HG2	1:E:110:ARG:NH1	1.81	0.86
1:D:23:TYR:CZ	1:D:36:TYR:CE2	2.62	0.86
1:I:117:GLU:CG	5:I:936:HOH:O	2.23	0.85
1:J:29:ASP:O	1:J:33:ARG:HG3	1.77	0.84
1:F:60:ARG:HD3	1:K:90:ARG:CZ	2.08	0.84
1:A:80:ILE:H	1:A:125:ASN:HD21	1.25	0.83
1:F:23:TYR:CZ	1:F:36:TYR:CE2	2.67	0.83
4:K:700:HEM:HMC1	4:K:700:HEM:HBC2	1.60	0.82
1:L:29:ASP:O	1:L:33[A]:ARG:HG2	1.80	0.81
1:H:92:MET:HG3	1:H:116:LEU:HD13	1.62	0.81
1:L:110:ARG:HE	1:L:110:ARG:HA	1.44	0.81

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:E:23:TYR:CZ	1:E:36:TYR:HE2	1.92	0.80
1:A:72:ARG:HH22	1:L:79:ARG:HG3	1.47	0.80
1:A:118:MET:HE1	1:D:82:LEU:CD2	2.12	0.80
1:J:92:MET:HG3	1:J:116:LEU:HD13	1.63	0.80
1:A:47[A]:ARG:NH2	3:A:802:SO4:O2	2.15	0.79
5:A:937:HOH:O	1:G:90:ARG:HD2	1.82	0.79
1:I:117:GLU:HG3	5:I:936:HOH:O	1.84	0.78
1:G:102:GLU:HG3	5:G:906:HOH:O	1.83	0.78
1:I:54:GLN:HE22	1:I:63:SER:H	1.28	0.78
1:I:80:ILE:H	1:I:125:ASN:HD21	1.32	0.78
1:E:110:ARG:HH11	1:E:110:ARG:CG	1.95	0.77
1:L:92:MET:HG3	1:L:116:LEU:HD13	1.67	0.76
1:C:38:GLU:CD	1:G:60:ARG:NH1	2.39	0.76
1:H:54:GLN:HE22	1:H:63:SER:H	1.34	0.75
4:J:700:HEM:HMC1	4:J:700:HEM:HBC2	1.66	0.75
1:A:70:ARG:HD3	5:A:934:HOH:O	1.85	0.75
1:A:54:GLN:HE22	1:A:63:SER:H	1.34	0.75
1:D:69:PRO:HB2	5:L:906:HOH:O	1.86	0.75
1:K:92:MET:HG3	1:K:116:LEU:HD13	1.70	0.74
4:E:700:HEM:HMC1	4:E:700:HEM:HBC2	1.69	0.74
1:K:29:ASP:O	1:K:33[A]:ARG:HG2	1.87	0.74
1:E:92:MET:CG	1:E:116:LEU:HD13	2.17	0.74
1:H:49:ARG:CD	3:H:801:SO4:O2	2.36	0.74
1:A:118:MET:HE1	1:D:82:LEU:HD21	1.67	0.73
1:I:118:MET:SD	5:I:960:HOH:O	2.47	0.72
1:E:102:GLU:HG3	5:E:897:HOH:O	1.88	0.72
1:H:80:ILE:H	1:H:125:ASN:HD21	1.36	0.72
1:D:46:GLU:HG2	5:D:893:HOH:O	1.90	0.71
1:B:54:GLN:HE22	1:B:63:SER:H	1.38	0.71
1:A:118:MET:HE3	1:D:82:LEU:HD21	1.74	0.70
1:K:109:ARG:O	1:K:113[B]:LEU:HD13	1.90	0.70
1:J:46:GLU:O	1:J:50:MET:HG3	1.92	0.70
1:L:54:GLN:HE22	1:L:63:SER:H	1.37	0.70
1:E:23:TYR:OH	1:E:36:TYR:CE2	1.70	0.69
4:L:700:HEM:HMC1	4:L:700:HEM:HBC2	1.74	0.69
1:F:118:MET:SD	1:K:127:PRO:HB3	2.33	0.69
1:D:92:MET:CG	1:D:116:LEU:HD13	2.23	0.69
1:G:34:ARG:NH1	5:G:954:HOH:O	2.23	0.68
1:D:92:MET:HG2	1:D:116:LEU:HD13	1.76	0.68
1:J:49:ARG:HD3	5:J:900:HOH:O	1.93	0.68
1:I:117:GLU:O	1:I:121[B]:HIS:CD2	2.46	0.68

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:60:ARG:NH1	5:A:937:HOH:O	2.25	0.68
1:I:85:ARG:HD2	1:I:121[B]:HIS:CD2	2.29	0.68
4:K:700:HEM:HBC2	4:K:700:HEM:CMC	2.24	0.67
1:K:92:MET:CG	1:K:116:LEU:HD13	2.24	0.67
1:F:83:ILE:HD12	5:F:934:HOH:O	1.94	0.67
1:L:92:MET:CG	1:L:116:LEU:HD13	2.25	0.67
1:B:69:PRO:HB2	5:H:910:HOH:O	1.95	0.67
1:G:49:ARG:NE	5:G:941:HOH:O	2.28	0.67
1:H:49:ARG:HG2	3:H:801:SO4:O2	1.95	0.67
1:H:85:ARG:HH11	1:H:121:HIS:HD2	1.42	0.66
1:H:85:ARG:HH11	1:H:121:HIS:CD2	2.14	0.66
4:G:700:HEM:HBC2	4:G:700:HEM:HMC2	1.76	0.66
1:C:38:GLU:HG3	1:G:60:ARG:HH12	1.62	0.65
1:E:68:HIS:HE1	1:J:84:GLU:OE1	1.78	0.65
1:A:72:ARG:HH21	1:L:79:ARG:HG2	1.53	0.65
1:A:125:ASN:C	1:A:125:ASN:HD22	1.99	0.64
1:F:59:PRO:HB2	1:F:61:THR:CG2	2.27	0.64
1:E:92:MET:HG2	1:E:116:LEU:HD13	1.78	0.64
1:K:46:GLU:O	1:K:50:MET:HG3	1.98	0.63
1:K:113[B]:LEU:O	1:K:117:GLU:HG3	1.98	0.63
1:F:92:MET:HG3	1:F:116:LEU:HD13	1.80	0.63
1:G:49:ARG:HD3	5:G:945:HOH:O	1.98	0.63
1:D:54:GLN:HE22	1:D:63:SER:H	1.44	0.63
4:I:700:HEM:HBC2	4:I:700:HEM:CMC	2.27	0.62
1:J:16:ASP:HB2	5:J:882:HOH:O	1.98	0.62
4:J:700:HEM:HBC2	4:J:700:HEM:CMC	2.29	0.62
1:C:125:ASN:C	1:C:125:ASN:HD22	2.03	0.62
1:L:63:SER:HB3	5:L:859:HOH:O	1.99	0.62
1:B:60:ARG:HD3	1:H:90:ARG:CZ	2.30	0.61
1:B:70:ARG:NH2	1:I:46:GLU:OE1	2.22	0.61
1:F:92:MET:CG	1:F:116:LEU:HD13	2.29	0.61
1:J:92:MET:CG	1:J:116:LEU:HD13	2.30	0.61
1:D:46:GLU:CG	5:D:931:HOH:O	2.40	0.61
1:C:80:ILE:H	1:C:125:ASN:ND2	1.94	0.61
1:E:92:MET:HG3	1:E:116:LEU:HD13	1.81	0.61
1:H:92:MET:CG	1:H:116:LEU:HD13	2.30	0.61
1:I:80:ILE:H	1:I:125:ASN:ND2	1.97	0.61
1:G:29:ASP:O	1:G:33:ARG:HG2	2.00	0.61
1:F:38:GLU:HB3	5:F:916:HOH:O	1.99	0.61
1:D:3:LYS:HD3	1:D:3:LYS:C	2.21	0.60
1:G:60:ARG:NH1	3:G:802:SO4:O4	2.34	0.60

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:H:80:ILE:H	1:H:125:ASN:ND2	1.99	0.60
1:G:19:VAL:HG12	1:G:45:GLU:HG3	1.84	0.60
1:H:49:ARG:CG	3:H:801:SO4:O2	2.50	0.60
1:L:19:VAL:HG12	1:L:45:GLU:HG3	1.84	0.59
1:I:17:ALA:O	1:I:21:ARG:HD3	2.02	0.59
1:A:118:MET:HE1	1:D:82:LEU:HD23	1.82	0.59
1:H:70:ARG:HD2	5:H:817:HOH:O	2.03	0.59
1:A:90:ARG:NH2	5:A:848:HOH:O	2.30	0.58
4:E:700:HEM:HBC2	4:E:700:HEM:CMC	2.33	0.58
1:J:34:ARG:HH22	1:L:46:GLU:CD	2.06	0.58
1:D:3:LYS:HD3	1:D:3:LYS:O	2.04	0.58
1:G:3:LYS:HD3	5:G:881:HOH:O	2.04	0.57
1:K:113[A]:LEU:O	1:K:117:GLU:HG3	2.04	0.57
1:F:30:GLU:CG	5:F:936:HOH:O	2.21	0.57
4:G:700:HEM:HBC2	4:G:700:HEM:CMC	2.34	0.57
1:G:49:ARG:NH2	5:G:901:HOH:O	2.25	0.57
1:F:60:ARG:HD3	1:K:90:ARG:NE	2.19	0.57
1:F:19:VAL:HG12	1:F:45:GLU:HG3	1.87	0.57
1:H:93:HIS:CE1	1:H:117:GLU:OE1	2.57	0.56
1:C:70:ARG:NE	5:C:935:HOH:O	2.39	0.56
1:D:68:HIS:CD2	1:D:70:ARG:NH1	2.74	0.56
1:E:23:TYR:CE1	1:E:36:TYR:CE2	2.93	0.56
1:J:126[A]:SER:OG	1:J:127:PRO:HD2	2.05	0.56
3:B:801:SO4:O4	1:I:50:MET:CG	2.53	0.56
1:I:19:VAL:HG12	1:I:45:GLU:HG3	1.87	0.56
4:L:700:HEM:HBC2	4:L:700:HEM:CMC	2.36	0.56
1:A:47[A]:ARG:HH21	1:H:59:PRO:HG3	1.71	0.56
1:G:34:ARG:NH2	1:H:46:GLU:OE1	2.39	0.56
1:D:69:PRO:HG2	5:L:906:HOH:O	2.06	0.56
1:A:118:MET:CE	1:D:82:LEU:HD23	2.31	0.55
1:L:54:GLN:HE21	1:L:60:ARG:HA	1.71	0.55
1:B:21:ARG:NH1	1:B:25:GLN:HG2	2.22	0.55
1:J:63:SER:HB3	5:J:816:HOH:O	2.07	0.55
1:F:70:ARG:HD2	5:F:885:HOH:O	2.07	0.55
1:L:3:LYS:N	5:L:917:HOH:O	2.40	0.55
1:H:49:ARG:NE	3:H:801:SO4:O2	2.39	0.54
1:A:72:ARG:CZ	5:A:888:HOH:O	2.56	0.54
1:G:70:ARG:HD2	5:G:931:HOH:O	2.08	0.54
1:I:54:GLN:NE2	1:I:63:SER:H	2.02	0.54
1:A:68:HIS:CE1	1:G:84:GLU:OE1	2.61	0.54
1:E:68:HIS:ND1	1:E:69:PRO:HD2	2.22	0.54

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:19:VAL:HG12	1:K:45:GLU:HG3	1.90	0.54
1:D:69:PRO:CB	5:L:906:HOH:O	2.48	0.54
1:G:33:ARG:NE	5:G:935:HOH:O	2.40	0.53
1:I:125:ASN:C	1:I:125:ASN:HD22	2.11	0.53
1:K:30:GLU:CD	1:K:30:GLU:H	2.10	0.53
1:E:25[B]:GLN:O	1:E:26:VAL:C	2.46	0.53
1:A:90:ARG:NE	5:A:848:HOH:O	2.24	0.53
1:D:68:HIS:NE2	1:D:70:ARG:NH1	2.57	0.52
1:J:74[B]:ARG:HD3	4:J:700:HEM:CGD	2.39	0.52
1:K:72:ARG:NE	5:K:907:HOH:O	2.22	0.52
1:F:70:ARG:HH11	1:F:70:ARG:HG2	1.75	0.52
1:A:80:ILE:H	1:A:125:ASN:ND2	2.00	0.52
1:H:125:ASN:C	1:H:125:ASN:HD22	2.13	0.52
1:B:70:ARG:NH1	1:I:46:GLU:OE2	2.33	0.52
1:D:38:GLU:HB3	5:D:920:HOH:O	2.09	0.52
1:K:113[B]:LEU:N	1:K:113[B]:LEU:CD1	2.73	0.51
1:L:50:MET:HE2	1:L:62:TYR:HA	1.92	0.51
1:G:49:ARG:CD	5:G:941:HOH:O	2.58	0.51
1:E:76:ALA:N	1:E:77:PRO:CD	2.74	0.51
1:E:121:HIS:HE1	5:E:916:HOH:O	1.94	0.51
1:E:68:HIS:CE1	1:J:84:GLU:OE1	2.61	0.51
1:H:49:ARG:HD2	5:H:880:HOH:O	2.11	0.51
1:A:25:GLN:CG	5:A:913:HOH:O	2.45	0.51
1:E:121:HIS:CD2	5:E:880:HOH:O	2.64	0.51
1:G:49:ARG:CZ	5:G:941:HOH:O	2.57	0.51
1:L:46:GLU:O	1:L:50:MET:HG3	2.10	0.50
1:A:125:ASN:C	1:A:125:ASN:ND2	2.65	0.50
1:K:21:ARG:HD3	1:K:98:SER:OG	2.11	0.50
1:K:113[B]:LEU:N	1:K:113[B]:LEU:HD12	2.25	0.50
4:F:700:HEM:HBC2	4:F:700:HEM:HMC1	1.94	0.50
1:E:21:ARG:HH12	1:E:25[A]:GLN:NE2	2.09	0.49
1:L:110:ARG:HE	1:L:110:ARG:CA	2.13	0.49
1:C:25:GLN:HE22	1:C:94:THR:HB	1.77	0.49
1:H:54:GLN:NE2	1:H:63:SER:H	2.08	0.49
4:A:700:HEM:HBC2	4:A:700:HEM:HMC1	1.95	0.49
1:F:33:ARG:HG3	5:F:910:HOH:O	2.11	0.49
5:F:939:HOH:O	1:J:60:ARG:HG3	2.12	0.49
1:J:30:GLU:CD	1:J:30:GLU:H	2.16	0.49
1:D:42:ALA:O	1:D:46:GLU:HG3	2.12	0.49
1:F:68:HIS:HB3	5:F:862:HOH:O	2.13	0.49
1:J:3:LYS:O	1:J:3:LYS:HG3	2.12	0.49

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:39:ASP:HA	5:A:924:HOH:O	2.13	0.48
1:D:38:GLU:HA	1:D:38:GLU:OE1	2.12	0.48
1:E:107:GLU:HG2	5:E:861:HOH:O	2.11	0.48
1:F:72:ARG:NH1	5:F:896:HOH:O	2.46	0.48
1:H:63:SER:HB3	5:H:901:HOH:O	2.13	0.48
1:E:66:ARG:NH2	1:E:74:ARG:HH21	2.11	0.48
1:G:56:TRP:CZ3	3:G:801:SO4:O1	2.66	0.48
1:E:74:ARG:HD3	5:E:843:HOH:O	2.12	0.48
1:C:69:PRO:HG2	5:I:947:HOH:O	2.13	0.48
1:A:102:GLU:HG3	5:A:917:HOH:O	2.13	0.48
1:L:72:ARG:CD	5:L:812:HOH:O	2.34	0.48
1:A:40:ASP:HB3	5:A:928:HOH:O	2.13	0.48
1:H:49:ARG:HD2	3:H:801:SO4:O2	2.13	0.47
4:C:700:HEM:HBC2	4:C:700:HEM:CMC	2.44	0.47
1:B:74:ARG:HG3	5:B:869:HOH:O	2.13	0.47
1:D:23:TYR:CE1	1:D:36:TYR:CE2	3.03	0.47
1:F:23:TYR:OH	1:F:36:TYR:CD2	2.54	0.47
4:A:700:HEM:HBC2	4:A:700:HEM:CMC	2.44	0.47
5:D:933:HOH:O	1:L:90:ARG:HD2	2.13	0.47
1:A:72:ARG:HH22	1:L:79:ARG:CD	2.21	0.47
1:B:60:ARG:NH1	1:H:90:ARG:HD3	2.30	0.47
1:F:69:PRO:HA	5:F:887:HOH:O	2.15	0.47
1:J:19:VAL:HG12	1:J:45:GLU:HG3	1.96	0.47
1:E:29:ASP:O	1:E:33:ARG:HG2	2.15	0.47
1:D:25[B]:GLN:O	1:D:26:VAL:C	2.52	0.47
1:H:76:ALA:N	1:H:77:PRO:CD	2.78	0.47
1:D:69:PRO:CG	5:L:906:HOH:O	2.63	0.46
1:G:49:ARG:HD2	5:G:956:HOH:O	2.16	0.46
1:D:30:GLU:HG3	1:D:33:ARG:NH2	2.31	0.46
1:D:60:ARG:HD3	1:L:90:ARG:NE	2.30	0.46
1:J:110:ARG:HD3	5:J:923:HOH:O	2.15	0.46
1:A:110:ARG:HD3	5:A:923:HOH:O	2.14	0.46
1:E:21:ARG:NH1	1:E:25[A]:GLN:CG	2.79	0.46
1:F:25:GLN:OE1	5:F:907:HOH:O	2.21	0.46
1:H:54:GLN:HE22	1:H:63:SER:N	2.09	0.46
1:G:37:PRO:HG2	1:G:37:PRO:O	2.16	0.46
1:A:39:ASP:HB3	5:A:924:HOH:O	2.15	0.46
1:A:39:ASP:CB	5:A:924:HOH:O	2.63	0.46
1:H:93:HIS:HD2	5:H:847:HOH:O	1.99	0.46
1:A:68:HIS:CG	1:A:69:PRO:HD2	2.51	0.45
4:C:700:HEM:HBC2	4:C:700:HEM:HMC1	1.98	0.45

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:K:109:ARG:HD3	5:K:847:HOH:O	2.15	0.45
1:B:73:MET:HG3	1:I:65:GLN:HE22	1.81	0.45
1:L:6:TYR:HE1	1:L:53:GLU:HG2	1.82	0.45
1:B:82:LEU:HD22	1:B:85:ARG:NH2	2.32	0.45
1:E:70:ARG:N	5:E:892:HOH:O	2.28	0.45
1:H:54:GLN:HE21	1:H:60:ARG:HA	1.82	0.45
1:A:38:GLU:HA	1:A:38:GLU:OE1	2.17	0.45
1:F:88:PHE:CB	4:F:700:HEM:HAB	2.46	0.45
1:L:3:LYS:HD3	1:L:3:LYS:HA	1.39	0.45
1:F:59:PRO:HB2	1:F:61:THR:HG22	1.99	0.45
1:G:79:ARG:NH2	5:G:909:HOH:O	2.39	0.45
1:K:72:ARG:NH2	5:K:907:HOH:O	2.47	0.45
1:A:76:ALA:N	1:A:77:PRO:CD	2.81	0.45
1:A:37:PRO:HG3	1:A:78:PHE:CE2	2.52	0.44
1:F:49[B]:ARG:HD3	5:F:809:HOH:O	2.16	0.44
1:K:92:MET:HG2	1:K:116:LEU:HD13	1.99	0.44
4:F:700:HEM:HBC2	4:F:700:HEM:CMC	2.47	0.44
1:G:23:TYR:OH	5:G:944:HOH:O	2.21	0.44
1:G:36:TYR:HB3	1:G:37:PRO:HD2	1.99	0.44
1:I:76:ALA:N	1:I:77:PRO:CD	2.80	0.44
1:H:93:HIS:HE1	1:H:117:GLU:OE1	1.99	0.44
1:E:102:GLU:CD	5:E:897:HOH:O	2.56	0.44
1:L:16:ASP:HB2	5:L:837:HOH:O	2.18	0.43
1:C:38:GLU:CG	1:G:60:ARG:HH12	2.27	0.43
1:B:110:ARG:HD3	1:B:114:ASP:OD1	2.18	0.43
1:C:104:LEU:HG	1:C:109:ARG:HB2	2.01	0.43
1:F:23:TYR:CE1	1:F:36:TYR:CE2	3.05	0.43
1:F:76:ALA:N	1:F:77:PRO:CD	2.82	0.43
4:K:700:HEM:HMC1	4:K:700:HEM:CBC	2.39	0.43
1:A:30:GLU:CD	1:A:33:ARG:HH21	2.21	0.43
1:H:85:ARG:NH1	1:H:121:HIS:CD2	2.85	0.43
1:C:69:PRO:HB2	5:I:947:HOH:O	2.18	0.43
1:L:38:GLU:OE2	1:L:39:ASP:HB3	2.19	0.43
1:I:30:GLU:CD	1:I:30:GLU:H	2.21	0.42
1:F:68:HIS:HD2	1:K:31:VAL:HG13	1.84	0.42
1:L:16:ASP:HB2	1:L:49:ARG:HH11	1.84	0.42
1:A:118:MET:HE2	1:A:118:MET:HB2	1.88	0.42
1:E:19:VAL:HG21	1:E:49:ARG:HA	2.01	0.42
1:H:19:VAL:HG12	1:H:45:GLU:HG3	2.01	0.42
1:L:115:TYR:CE1	4:L:700:HEM:HBC1	2.54	0.42
1:L:16:ASP:HA	1:L:49:ARG:HD3	2.00	0.42

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:35:VAL:CG2	1:A:78:PHE:HB3	2.49	0.42
1:K:26:VAL:HG13	1:K:32:LEU:HB3	2.00	0.42
1:A:38:GLU:HB3	5:A:928:HOH:O	2.19	0.42
1:I:54:GLN:HE21	1:I:60:ARG:HA	1.85	0.42
1:L:76:ALA:N	1:L:77:PRO:CD	2.83	0.42
1:A:36:TYR:HA	1:A:37:PRO:HD3	1.86	0.42
1:E:21:ARG:O	1:E:25[B]:GLN:HG3	2.19	0.42
1:E:102:GLU:CG	5:E:897:HOH:O	2.56	0.42
1:F:73:MET:HG3	1:J:65:GLN:HE22	1.85	0.42
1:H:74:ARG:HD3	4:H:700:HEM:CGD	2.50	0.42
1:J:42:ALA:HA	1:K:33[B]:ARG:HH12	1.84	0.42
1:D:68:HIS:CE1	1:L:84:GLU:OE2	2.73	0.41
1:F:54:GLN:OE1	1:F:62:TYR:HB3	2.19	0.41
1:G:33:ARG:NH2	5:G:925:HOH:O	2.53	0.41
1:B:3:LYS:HA	1:B:3:LYS:HD3	1.82	0.41
1:J:76:ALA:N	1:J:77:PRO:CD	2.83	0.41
1:A:25:GLN:OE1	1:A:91:CYS:HA	2.20	0.41
1:A:54:GLN:NE2	1:A:62:TYR:HB3	2.36	0.41
1:D:79:ARG:NH2	1:G:72:ARG:HD3	2.35	0.41
1:F:125:ASN:O	1:I:125:ASN:HA	2.20	0.41
1:D:26:VAL:HG13	1:D:32:LEU:HB3	2.03	0.41
1:B:27:ALA:HA	1:B:33:ARG:HD3	2.02	0.41
1:D:126:SER:HA	1:D:127:PRO:HD3	1.92	0.41
1:G:76:ALA:N	1:G:77:PRO:CD	2.83	0.41
1:I:54:GLN:HE22	1:I:63:SER:N	2.05	0.41
1:C:38:GLU:CG	1:G:60:ARG:NH1	2.84	0.41
1:E:88:PHE:HE1	2:E:800:CYN:C	2.34	0.41
1:E:121:HIS:HD2	5:E:880:HOH:O	2.01	0.41
1:K:12:ALA:HA	1:K:53:GLU:OE1	2.20	0.41
1:L:126[A]:SER:OG	1:L:127:PRO:HD2	2.21	0.41
1:A:68:HIS:HE1	1:G:84:GLU:OE1	2.03	0.41
1:A:72:ARG:NH2	1:L:79:ARG:NE	2.69	0.41
1:B:35:VAL:CG2	1:B:78:PHE:HB3	2.51	0.41
1:D:92:MET:HG3	1:D:116:LEU:HD13	1.99	0.41
1:E:71:LEU:HD23	1:E:71:LEU:HA	1.87	0.40
1:A:65:GLN:O	3:H:801:SO4:O3	2.39	0.40
1:B:76:ALA:N	1:B:77:PRO:CD	2.83	0.40
1:F:68:HIS:CD2	1:F:69:PRO:HD2	2.56	0.40
1:I:33:ARG:NE	5:I:957:HOH:O	2.48	0.40
1:F:49[A]:ARG:HD2	5:F:849:HOH:O	2.22	0.40
1:F:59:PRO:CB	1:F:61:THR:HG22	2.51	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	126/128 (98%)	125 (99%)	1 (1%)	0	100	100
1	B	125/128 (98%)	122 (98%)	3 (2%)	0	100	100
1	C	126/128 (98%)	125 (99%)	1 (1%)	0	100	100
1	D	125/128 (98%)	122 (98%)	2 (2%)	1 (1%)	19	9
1	E	125/128 (98%)	124 (99%)	1 (1%)	0	100	100
1	F	126/128 (98%)	124 (98%)	2 (2%)	0	100	100
1	G	128/128 (100%)	127 (99%)	1 (1%)	0	100	100
1	H	126/128 (98%)	125 (99%)	1 (1%)	0	100	100
1	I	128/128 (100%)	128 (100%)	0	0	100	100
1	J	127/128 (99%)	126 (99%)	1 (1%)	0	100	100
1	K	129/128 (101%)	128 (99%)	1 (1%)	0	100	100
1	L	127/128 (99%)	125 (98%)	1 (1%)	1 (1%)	19	9
All	All	1518/1536 (99%)	1501 (99%)	15 (1%)	2 (0%)	51	43

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	L	105	ASP
1	D	37	PRO

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	109/109 (100%)	105 (96%)	4 (4%)	34	19
1	B	108/109 (99%)	105 (97%)	3 (3%)	43	29
1	C	109/109 (100%)	106 (97%)	3 (3%)	43	29
1	D	108/109 (99%)	103 (95%)	5 (5%)	27	12
1	E	108/109 (99%)	105 (97%)	3 (3%)	43	29
1	F	109/109 (100%)	103 (94%)	6 (6%)	21	8
1	G	111/109 (102%)	107 (96%)	4 (4%)	35	20
1	H	109/109 (100%)	106 (97%)	3 (3%)	43	29
1	I	111/109 (102%)	109 (98%)	2 (2%)	59	47
1	J	110/109 (101%)	108 (98%)	2 (2%)	59	47
1	K	112/109 (103%)	108 (96%)	4 (4%)	35	20
1	L	110/109 (101%)	105 (96%)	5 (4%)	27	12
All	All	1314/1308 (100%)	1270 (97%)	44 (3%)	37	24

All (44) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	23	TYR
1	A	39	ASP
1	A	104	LEU
1	A	125	ASN
1	B	23	TYR
1	B	110	ARG
1	B	117	GLU
1	C	23	TYR
1	C	102	GLU
1	C	125	ASN
1	D	3	LYS
1	D	23	TYR
1	D	32	LEU
1	D	71	LEU
1	D	117	GLU
1	E	23	TYR
1	E	74	ARG
1	E	110	ARG
1	F	23	TYR

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Mol	Chain	Res	Type
1	F	61	THR
1	F	82	LEU
1	F	89	LEU
1	F	113	LEU
1	F	118	MET
1	G	3	LYS
1	G	34	ARG
1	G	60	ARG
1	G	117	GLU
1	H	92	MET
1	H	113	LEU
1	H	125	ASN
1	I	28	GLU
1	I	125	ASN
1	J	104	LEU
1	J	113	LEU
1	K	30	GLU
1	K	32	LEU
1	K	82	LEU
1	K	104	LEU
1	L	3	LYS
1	L	104	LEU
1	L	110	ARG
1	L	113	LEU
1	L	118	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (27) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	54	GLN
1	A	68	HIS
1	A	125	ASN
1	B	54	GLN
1	B	121	HIS
1	C	25	GLN
1	C	93	HIS
1	C	125	ASN
1	D	54	GLN
1	D	65	GLN
1	E	65	GLN
1	E	68	HIS
1	E	93	HIS

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Mol	Chain	Res	Type
1	F	65	GLN
1	F	121	HIS
1	G	68	HIS
1	G	93	HIS
1	H	54	GLN
1	H	121	HIS
1	H	125	ASN
1	I	54	GLN
1	I	65	GLN
1	I	125	ASN
1	J	65	GLN
1	J	93	HIS
1	L	54	GLN
1	L	93	HIS

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

47 ligands are modelled in this entry.

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

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
4	HEM	I	700	1,2	41,50,50	1.89	5 (12%)	45,82,82	1.77	10 (22%)
2	CYN	B	800	4	0,1,1	-	-	-		
2	CYN	L	800	4	0,1,1	-	-	-		
3	SO4	D	802	-	4,4,4	0.29	0	6,6,6	0.21	0
3	SO4	L	802	-	4,4,4	0.16	0	6,6,6	0.17	0
3	SO4	B	801	-	4,4,4	0.29	0	6,6,6	0.46	0
3	SO4	L	801	-	4,4,4	0.14	0	6,6,6	0.16	0
2	CYN	E	800	4	0,1,1	-	-	-		
3	SO4	A	804	-	4,4,4	0.15	0	6,6,6	0.10	0
2	CYN	G	800	4	0,1,1	-	-	-		
3	SO4	E	802	-	4,4,4	0.51	0	6,6,6	1.14	1 (16%)
3	SO4	A	802	-	4,4,4	0.14	0	6,6,6	0.34	0
2	CYN	C	800	4	0,1,1	-	-	-		
3	SO4	F	803	-	4,4,4	0.35	0	6,6,6	0.37	0
4	HEM	G	700	1,2	41,50,50	1.85	6 (14%)	45,82,82	1.75	10 (22%)
3	SO4	A	803	-	4,4,4	0.11	0	6,6,6	0.24	0
4	HEM	E	700	1,2	41,50,50	1.86	6 (14%)	45,82,82	1.87	6 (13%)
2	CYN	H	800	4	0,1,1	-	-	-		
2	CYN	J	800	4	0,1,1	-	-	-		
3	SO4	J	801	-	4,4,4	0.16	0	6,6,6	0.18	0
3	SO4	F	802	-	4,4,4	0.25	0	6,6,6	0.34	0
4	HEM	F	700	1,2	41,50,50	1.84	6 (14%)	45,82,82	1.83	8 (17%)
3	SO4	G	802	-	4,4,4	0.18	0	6,6,6	0.33	0
4	HEM	A	700	1,2	41,50,50	1.82	4 (9%)	45,82,82	1.84	10 (22%)
4	HEM	L	700	1,2	41,50,50	1.75	5 (12%)	45,82,82	1.79	11 (24%)
2	CYN	D	800	4	0,1,1	-	-	-		
3	SO4	D	801	-	4,4,4	0.36	0	6,6,6	0.20	0
3	SO4	B	803	-	4,4,4	0.20	0	6,6,6	0.21	0
3	SO4	B	802	-	4,4,4	0.15	0	6,6,6	0.49	0
3	SO4	L	803	-	4,4,4	0.22	0	6,6,6	0.21	0
4	HEM	K	700	1,2	41,50,50	1.92	9 (21%)	45,82,82	1.79	7 (15%)
2	CYN	K	800	4	0,1,1	-	-	-		
4	HEM	C	700	1,2	41,50,50	1.78	7 (17%)	45,82,82	1.90	10 (22%)
2	CYN	I	800	4	0,1,1	-	-	-		
3	SO4	K	801	-	4,4,4	0.22	0	6,6,6	0.14	0
4	HEM	H	700	1,2	41,50,50	1.89	7 (17%)	45,82,82	1.96	10 (22%)
3	SO4	E	801	-	4,4,4	0.09	0	6,6,6	0.21	0
3	SO4	H	801	-	4,4,4	0.14	0	6,6,6	0.20	0
4	HEM	B	700	1,2	41,50,50	1.83	5 (12%)	45,82,82	1.79	9 (20%)
2	CYN	F	800	4	0,1,1	-	-	-		

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\| > 2$	Counts	RMSZ	$\#\|Z\| > 2$
3	SO4	F	801	-	4,4,4	0.15	0	6,6,6	0.11	0
2	CYN	A	800	4	0,1,1	-	-	-		
4	HEM	J	700	1,2	41,50,50	1.87	6 (14%)	45,82,82	1.82	11 (24%)
3	SO4	A	801	-	4,4,4	0.13	0	6,6,6	0.26	0
3	SO4	G	801	-	4,4,4	0.16	0	6,6,6	0.25	0
3	SO4	I	801	-	4,4,4	0.06	0	6,6,6	0.37	0
4	HEM	D	700	1,2	41,50,50	1.90	5 (12%)	45,82,82	1.79	10 (22%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	HEM	I	700	1,2	-	5/12/54/54	-
4	HEM	H	700	1,2	-	5/12/54/54	-
4	HEM	L	700	1,2	-	7/12/54/54	-
4	HEM	G	700	1,2	-	5/12/54/54	-
4	HEM	B	700	1,2	-	5/12/54/54	-
4	HEM	E	700	1,2	-	4/12/54/54	-
4	HEM	K	700	1,2	-	5/12/54/54	-
4	HEM	J	700	1,2	-	6/12/54/54	-
4	HEM	F	700	1,2	-	6/12/54/54	-
4	HEM	C	700	1,2	-	6/12/54/54	-
4	HEM	A	700	1,2	-	6/12/54/54	-
4	HEM	D	700	1,2	-	5/12/54/54	-

All (71) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	I	700	HEM	C3D-C2D	7.74	1.53	1.36
4	J	700	HEM	C3D-C2D	7.63	1.53	1.36
4	A	700	HEM	C3D-C2D	7.42	1.52	1.36
4	G	700	HEM	C3D-C2D	7.38	1.52	1.36
4	D	700	HEM	C3D-C2D	7.37	1.52	1.36
4	F	700	HEM	C3D-C2D	7.31	1.52	1.36
4	K	700	HEM	C3D-C2D	7.30	1.52	1.36
4	H	700	HEM	C3D-C2D	7.25	1.52	1.36
4	E	700	HEM	C3D-C2D	7.24	1.52	1.36

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	C	700	HEM	C3D-C2D	7.18	1.52	1.36
4	B	700	HEM	C3D-C2D	7.01	1.51	1.36
4	L	700	HEM	C3D-C2D	6.82	1.51	1.36
4	K	700	HEM	C3C-C2C	-4.62	1.34	1.40
4	E	700	HEM	C3C-C2C	-4.60	1.34	1.40
4	I	700	HEM	C3C-C2C	-4.43	1.34	1.40
4	D	700	HEM	C3C-C2C	-4.40	1.34	1.40
4	A	700	HEM	C3C-C2C	-4.31	1.34	1.40
4	B	700	HEM	C3C-C2C	-4.19	1.34	1.40
4	H	700	HEM	C3C-C2C	-4.19	1.34	1.40
4	L	700	HEM	C3C-C2C	-4.15	1.34	1.40
4	G	700	HEM	C3C-C2C	-4.04	1.34	1.40
4	J	700	HEM	C3C-C2C	-3.92	1.34	1.40
4	F	700	HEM	C3C-CAC	3.69	1.55	1.47
4	D	700	HEM	C3C-CAC	3.68	1.55	1.47
4	F	700	HEM	C3C-C2C	-3.68	1.35	1.40
4	H	700	HEM	C3C-CAC	3.61	1.55	1.47
4	K	700	HEM	C3C-CAC	3.50	1.55	1.47
4	B	700	HEM	C3C-CAC	3.49	1.55	1.47
4	C	700	HEM	C3C-C2C	-3.37	1.35	1.40
4	L	700	HEM	C3C-CAC	3.36	1.54	1.47
4	J	700	HEM	C3C-CAC	3.28	1.54	1.47
4	G	700	HEM	C3C-CAC	3.26	1.54	1.47
4	E	700	HEM	C3C-CAC	3.24	1.54	1.47
4	B	700	HEM	FE-NB	3.08	2.12	1.96
4	A	700	HEM	C3C-CAC	3.03	1.54	1.47
4	I	700	HEM	C3C-CAC	2.96	1.53	1.47
4	C	700	HEM	C3C-CAC	2.91	1.53	1.47
4	I	700	HEM	FE-ND	2.81	2.10	1.96
4	D	700	HEM	CAB-C3B	2.79	1.55	1.47
4	E	700	HEM	CAB-C3B	2.72	1.54	1.47
4	L	700	HEM	CAB-C3B	2.72	1.54	1.47
4	H	700	HEM	FE-ND	2.57	2.09	1.96
4	J	700	HEM	CAB-C3B	2.56	1.54	1.47
4	H	700	HEM	FE-NB	2.54	2.09	1.96
4	G	700	HEM	FE-ND	2.46	2.09	1.96
4	B	700	HEM	CAB-C3B	2.43	1.54	1.47
4	H	700	HEM	CAB-C3B	2.39	1.53	1.47
4	C	700	HEM	FE-ND	2.36	2.08	1.96
4	K	700	HEM	CAB-C3B	2.36	1.53	1.47
4	D	700	HEM	CAA-C2A	2.31	1.55	1.52
4	K	700	HEM	CMB-C2B	2.30	1.55	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	F	700	HEM	CAB-C3B	2.28	1.53	1.47
4	A	700	HEM	CAB-C3B	2.27	1.53	1.47
4	L	700	HEM	FE-NB	2.25	2.08	1.96
4	K	700	HEM	FE-ND	2.25	2.08	1.96
4	I	700	HEM	CAB-C3B	2.22	1.53	1.47
4	G	700	HEM	CMD-C2D	2.22	1.55	1.50
4	K	700	HEM	CMA-C3A	2.17	1.56	1.51
4	H	700	HEM	CAA-C2A	2.16	1.55	1.52
4	F	700	HEM	FE-ND	2.15	2.07	1.96
4	K	700	HEM	FE-NB	2.14	2.07	1.96
4	C	700	HEM	FE-NB	2.12	2.07	1.96
4	C	700	HEM	CMD-C2D	2.12	1.55	1.50
4	J	700	HEM	CMD-C2D	2.11	1.55	1.50
4	G	700	HEM	CAB-C3B	2.11	1.53	1.47
4	K	700	HEM	CMD-C2D	2.11	1.55	1.50
4	E	700	HEM	FE-ND	2.10	2.07	1.96
4	J	700	HEM	FE-ND	2.08	2.07	1.96
4	C	700	HEM	CAB-C3B	2.05	1.53	1.47
4	F	700	HEM	CMB-C2B	2.02	1.55	1.50
4	E	700	HEM	CMB-C2B	2.02	1.55	1.50

All (113) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	H	700	HEM	C4D-ND-C1D	6.90	112.20	105.07
4	I	700	HEM	C4D-ND-C1D	6.48	111.76	105.07
4	E	700	HEM	CBA-CAA-C2A	-6.48	101.57	112.62
4	C	700	HEM	C4D-ND-C1D	6.22	111.49	105.07
4	J	700	HEM	C4D-ND-C1D	6.16	111.43	105.07
4	E	700	HEM	C4D-ND-C1D	6.11	111.39	105.07
4	K	700	HEM	C4D-ND-C1D	6.03	111.30	105.07
4	B	700	HEM	C4D-ND-C1D	5.89	111.15	105.07
4	L	700	HEM	C4D-ND-C1D	5.85	111.11	105.07
4	A	700	HEM	C4D-ND-C1D	5.68	110.94	105.07
4	F	700	HEM	C4D-ND-C1D	5.66	110.92	105.07
4	C	700	HEM	CBA-CAA-C2A	-5.49	103.25	112.62
4	A	700	HEM	CBA-CAA-C2A	-5.47	103.29	112.62
4	K	700	HEM	CBA-CAA-C2A	-5.46	103.30	112.62
4	G	700	HEM	C4D-ND-C1D	5.44	110.70	105.07
4	D	700	HEM	C4D-ND-C1D	5.42	110.67	105.07
4	B	700	HEM	CBA-CAA-C2A	-5.35	103.49	112.62
4	D	700	HEM	CBA-CAA-C2A	-5.28	103.60	112.62

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	H	700	HEM	CBA-CAA-C2A	-5.24	103.68	112.62
4	F	700	HEM	CBA-CAA-C2A	-4.68	104.63	112.62
4	L	700	HEM	CBA-CAA-C2A	-4.01	105.78	112.62
4	J	700	HEM	CBA-CAA-C2A	-3.84	106.07	112.62
4	F	700	HEM	C4B-CHC-C1C	3.65	127.37	122.56
4	G	700	HEM	CBA-CAA-C2A	-3.56	106.55	112.62
4	D	700	HEM	CAD-CBD-CGD	-3.27	106.56	113.60
4	H	700	HEM	C1B-NB-C4B	3.25	108.43	105.07
4	L	700	HEM	CAD-CBD-CGD	-3.24	106.64	113.60
4	G	700	HEM	CMA-C3A-C4A	-3.21	123.53	128.46
4	K	700	HEM	C4C-CHD-C1D	3.16	126.73	122.56
4	J	700	HEM	CMA-C3A-C4A	-3.16	123.61	128.46
4	E	700	HEM	CAD-CBD-CGD	-3.14	106.85	113.60
4	I	700	HEM	CBA-CAA-C2A	-3.08	107.36	112.62
4	F	700	HEM	CMA-C3A-C4A	-2.98	123.88	128.46
4	C	700	HEM	C1B-NB-C4B	2.96	108.13	105.07
4	L	700	HEM	C4C-CHD-C1D	2.91	126.39	122.56
4	F	700	HEM	CAD-CBD-CGD	-2.91	107.35	113.60
4	I	700	HEM	C4C-CHD-C1D	2.90	126.38	122.56
4	G	700	HEM	O1A-CGA-CBA	-2.89	113.79	123.08
4	A	700	HEM	CAD-CBD-CGD	-2.87	107.43	113.60
4	A	700	HEM	C4B-CHC-C1C	2.85	126.32	122.56
4	B	700	HEM	C1B-NB-C4B	2.82	107.98	105.07
4	I	700	HEM	C1B-NB-C4B	2.81	107.98	105.07
4	H	700	HEM	C4B-CHC-C1C	2.80	126.26	122.56
4	J	700	HEM	C1B-NB-C4B	2.74	107.91	105.07
4	C	700	HEM	C4C-CHD-C1D	2.73	126.17	122.56
4	F	700	HEM	C1B-NB-C4B	2.70	107.86	105.07
4	H	700	HEM	C2C-C3C-C4C	2.67	108.77	106.90
4	A	700	HEM	C1B-NB-C4B	2.67	107.83	105.07
4	I	700	HEM	CAD-CBD-CGD	-2.65	107.89	113.60
4	K	700	HEM	C1D-C2D-C3D	-2.64	104.18	106.96
4	G	700	HEM	CBD-CAD-C3D	-2.64	105.29	112.63
4	L	700	HEM	CBD-CAD-C3D	-2.62	105.34	112.63
4	E	700	HEM	C4C-CHD-C1D	2.61	126.00	122.56
4	B	700	HEM	CAD-CBD-CGD	-2.58	108.05	113.60
4	L	700	HEM	O1A-CGA-CBA	-2.57	114.82	123.08
4	C	700	HEM	CAD-CBD-CGD	-2.57	108.07	113.60
4	D	700	HEM	CMA-C3A-C4A	-2.53	124.57	128.46
4	A	700	HEM	C4C-CHD-C1D	2.51	125.87	122.56
4	K	700	HEM	CAD-CBD-CGD	-2.51	108.21	113.60
4	I	700	HEM	CAA-CBA-CGA	-2.51	106.73	113.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	K	700	HEM	C1B-NB-C4B	2.49	107.65	105.07
4	B	700	HEM	C4B-CHC-C1C	2.49	125.85	122.56
4	L	700	HEM	CMD-C2D-C1D	2.47	128.81	125.04
4	J	700	HEM	O1A-CGA-CBA	-2.46	115.19	123.08
4	D	700	HEM	C4C-CHD-C1D	2.41	125.75	122.56
4	B	700	HEM	C4C-CHD-C1D	2.39	125.72	122.56
4	A	700	HEM	C1D-C2D-C3D	-2.38	104.46	106.96
4	F	700	HEM	CBD-CAD-C3D	-2.38	106.03	112.63
4	G	700	HEM	C1B-NB-C4B	2.37	107.52	105.07
4	E	700	HEM	C3B-C2B-C1B	2.36	108.24	106.49
4	D	700	HEM	CBB-CAB-C3B	-2.35	115.92	127.62
4	A	700	HEM	CMA-C3A-C4A	-2.33	124.88	128.46
4	B	700	HEM	CMD-C2D-C1D	2.28	128.52	125.04
4	B	700	HEM	CMA-C3A-C4A	-2.28	124.96	128.46
4	H	700	HEM	C4C-CHD-C1D	2.26	125.54	122.56
3	E	802	SO4	O4-S-O2	-2.25	97.55	109.31
4	E	700	HEM	C1B-NB-C4B	2.25	107.40	105.07
4	D	700	HEM	C1B-NB-C4B	2.25	107.39	105.07
4	H	700	HEM	CAD-CBD-CGD	-2.24	108.78	113.60
4	D	700	HEM	C4A-C3A-C2A	2.24	108.55	107.00
4	C	700	HEM	CMD-C2D-C1D	2.24	128.44	125.04
4	H	700	HEM	CHA-C4D-ND	2.23	127.14	124.38
4	F	700	HEM	C1D-C2D-C3D	-2.21	104.64	106.96
4	H	700	HEM	O1A-CGA-CBA	-2.20	116.01	123.08
4	L	700	HEM	CAA-CBA-CGA	-2.20	107.59	113.76
4	J	700	HEM	C1D-C2D-C3D	-2.20	104.64	106.96
4	J	700	HEM	CAD-CBD-CGD	-2.20	108.87	113.60
4	L	700	HEM	C1D-C2D-C3D	-2.18	104.67	106.96
4	C	700	HEM	CAD-C3D-C4D	2.17	128.45	124.66
4	I	700	HEM	O1A-CGA-CBA	-2.17	116.12	123.08
4	G	700	HEM	C1D-C2D-C3D	-2.17	104.68	106.96
4	G	700	HEM	C2C-C3C-C4C	2.16	108.41	106.90
4	C	700	HEM	CBB-CAB-C3B	-2.16	116.87	127.62
4	J	700	HEM	C4B-CHC-C1C	2.13	125.37	122.56
4	I	700	HEM	CBD-CAD-C3D	-2.13	106.72	112.63
4	D	700	HEM	C4B-CHC-C1C	2.12	125.36	122.56
4	C	700	HEM	C4B-CHC-C1C	2.11	125.35	122.56
4	K	700	HEM	CAA-CBA-CGA	-2.11	107.85	113.76
4	I	700	HEM	C1D-C2D-C3D	-2.09	104.76	106.96
4	H	700	HEM	CMA-C3A-C4A	-2.09	125.25	128.46
4	L	700	HEM	CMA-C3A-C4A	-2.07	125.28	128.46
4	D	700	HEM	C1D-C2D-C3D	-2.07	104.78	106.96

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
4	J	700	HEM	CMA-C3A-C2A	2.06	128.82	124.94
4	I	700	HEM	C2C-C3C-C4C	2.05	108.33	106.90
4	G	700	HEM	CMA-C3A-C2A	2.05	128.81	124.94
4	A	700	HEM	CBD-CAD-C3D	-2.04	106.95	112.63
4	B	700	HEM	CBB-CAB-C3B	-2.04	117.48	127.62
4	L	700	HEM	C1B-NB-C4B	2.03	107.17	105.07
4	J	700	HEM	CMC-C2C-C3C	2.03	128.48	124.68
4	J	700	HEM	CAA-CBA-CGA	-2.03	108.07	113.76
4	C	700	HEM	O2A-CGA-CBA	2.02	120.52	114.03
4	G	700	HEM	CAA-CBA-CGA	-2.00	108.14	113.76
4	A	700	HEM	CBB-CAB-C3B	-2.00	117.66	127.62

There are no chirality outliers.

All (65) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
4	J	700	HEM	C2A-CAA-CBA-CGA
4	A	700	HEM	C2B-C3B-CAB-CBB
4	C	700	HEM	C2B-C3B-CAB-CBB
4	F	700	HEM	C2B-C3B-CAB-CBB
4	L	700	HEM	C2B-C3B-CAB-CBB
4	C	700	HEM	C4B-C3B-CAB-CBB
4	F	700	HEM	C4B-C3B-CAB-CBB
4	G	700	HEM	C2A-CAA-CBA-CGA
4	H	700	HEM	C2A-CAA-CBA-CGA
4	I	700	HEM	C2A-CAA-CBA-CGA
4	L	700	HEM	C2A-CAA-CBA-CGA
4	D	700	HEM	C2B-C3B-CAB-CBB
4	J	700	HEM	C2B-C3B-CAB-CBB
4	K	700	HEM	C2A-CAA-CBA-CGA
4	J	700	HEM	CAD-CBD-CGD-O2D
4	L	700	HEM	C4B-C3B-CAB-CBB
4	I	700	HEM	CAD-CBD-CGD-O1D
4	L	700	HEM	CAD-CBD-CGD-O2D
4	D	700	HEM	CAD-CBD-CGD-O2D
4	E	700	HEM	CAA-CBA-CGA-O1A
4	E	700	HEM	CAA-CBA-CGA-O2A
4	C	700	HEM	CAD-CBD-CGD-O2D
4	F	700	HEM	CAA-CBA-CGA-O2A
4	B	700	HEM	CAA-CBA-CGA-O1A
4	F	700	HEM	CAA-CBA-CGA-O1A
4	H	700	HEM	CAD-CBD-CGD-O1D

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Mol	Chain	Res	Type	Atoms
4	B	700	HEM	CAA-CBA-CGA-O2A
4	D	700	HEM	CAA-CBA-CGA-O1A
4	I	700	HEM	CAD-CBD-CGD-O2D
4	A	700	HEM	CAA-CBA-CGA-O1A
4	G	700	HEM	CAD-CBD-CGD-O1D
4	A	700	HEM	CAD-CBD-CGD-O2D
4	H	700	HEM	CAD-CBD-CGD-O2D
4	K	700	HEM	CAD-CBD-CGD-O1D
4	E	700	HEM	CAD-CBD-CGD-O2D
4	G	700	HEM	CAD-CBD-CGD-O2D
4	H	700	HEM	CAA-CBA-CGA-O1A
4	K	700	HEM	CAA-CBA-CGA-O2A
4	K	700	HEM	CAD-CBD-CGD-O2D
4	A	700	HEM	CAD-CBD-CGD-O1D
4	C	700	HEM	CAD-CBD-CGD-O1D
4	A	700	HEM	CAA-CBA-CGA-O2A
4	K	700	HEM	CAA-CBA-CGA-O1A
4	J	700	HEM	CAA-CBA-CGA-O1A
4	L	700	HEM	CAD-CBD-CGD-O1D
4	H	700	HEM	CAA-CBA-CGA-O2A
4	J	700	HEM	CAD-CBD-CGD-O1D
4	J	700	HEM	CAA-CBA-CGA-O2A
4	E	700	HEM	CAD-CBD-CGD-O1D
4	D	700	HEM	CAA-CBA-CGA-O2A
4	D	700	HEM	CAD-CBD-CGD-O1D
4	L	700	HEM	CAA-CBA-CGA-O2A
4	G	700	HEM	CAA-CBA-CGA-O2A
4	C	700	HEM	CAA-CBA-CGA-O2A
4	F	700	HEM	CAD-CBD-CGD-O1D
4	I	700	HEM	CAA-CBA-CGA-O2A
4	L	700	HEM	CAA-CBA-CGA-O1A
4	F	700	HEM	CAD-CBD-CGD-O2D
4	G	700	HEM	CAA-CBA-CGA-O1A
4	I	700	HEM	CAA-CBA-CGA-O1A
4	C	700	HEM	CAA-CBA-CGA-O1A
4	B	700	HEM	C2B-C3B-CAB-CBB
4	B	700	HEM	CAD-CBD-CGD-O1D
4	A	700	HEM	C4B-C3B-CAB-CBB
4	B	700	HEM	CAD-CBD-CGD-O2D

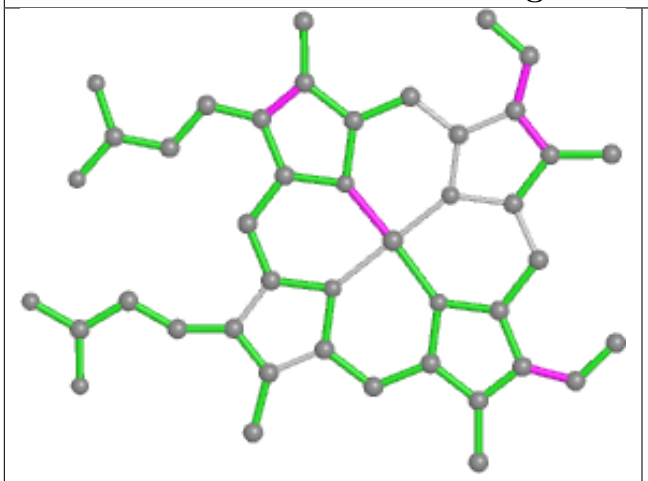
There are no ring outliers.

16 monomers are involved in 34 short contacts:

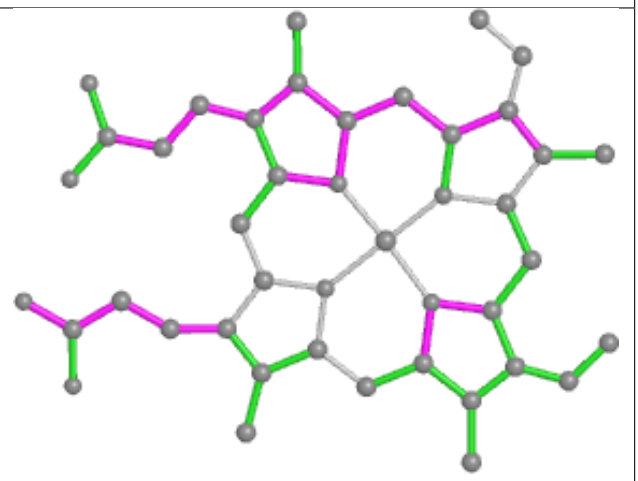
Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	I	700	HEM	2	0
3	B	801	SO4	1	0
2	E	800	CYN	1	0
3	A	802	SO4	1	0
4	G	700	HEM	2	0
4	E	700	HEM	2	0
4	F	700	HEM	3	0
3	G	802	SO4	1	0
4	A	700	HEM	2	0
4	L	700	HEM	3	0
4	K	700	HEM	3	0
4	C	700	HEM	2	0
4	H	700	HEM	1	0
3	H	801	SO4	6	0
4	J	700	HEM	3	0
3	G	801	SO4	1	0

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

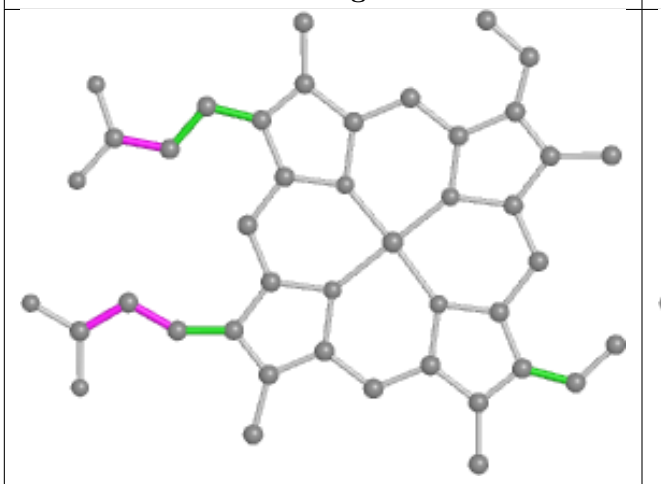
Ligand HEM I 700



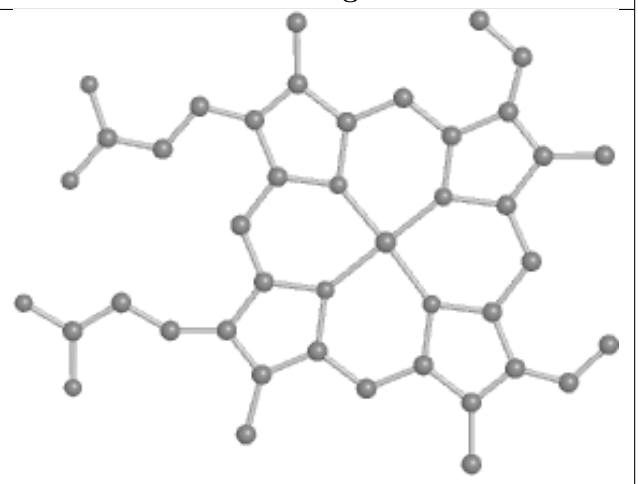
Bond lengths



Bond angles

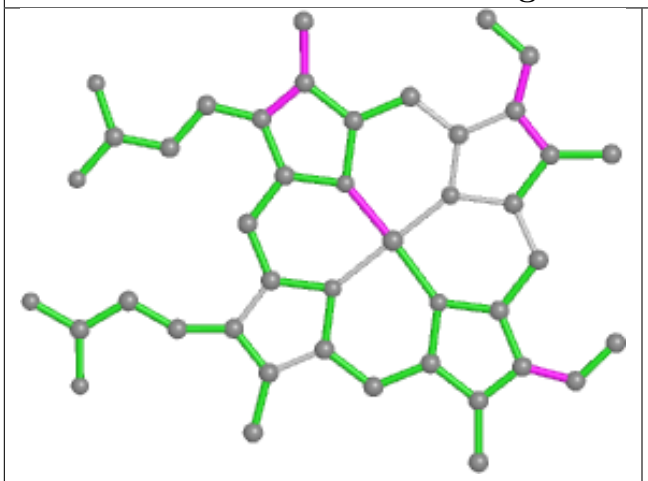


Torsions



Rings

Ligand HEM G 700



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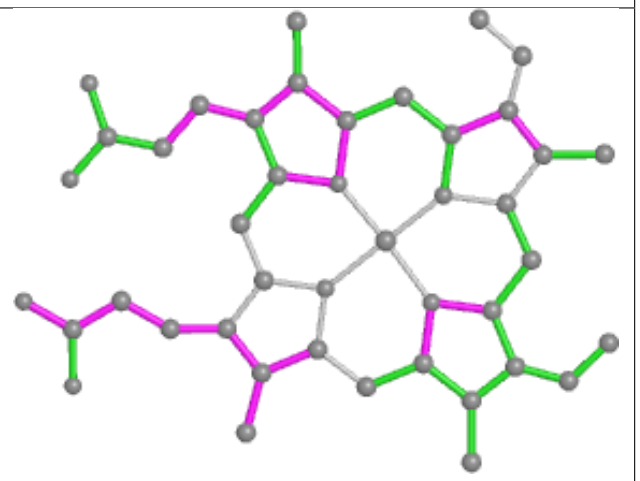
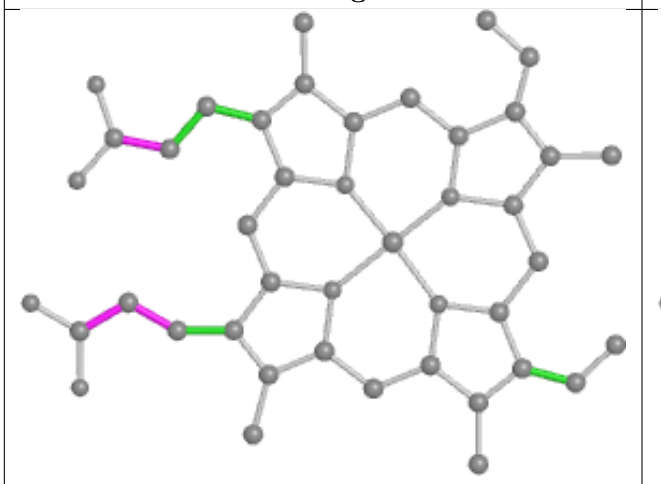


Bond angles



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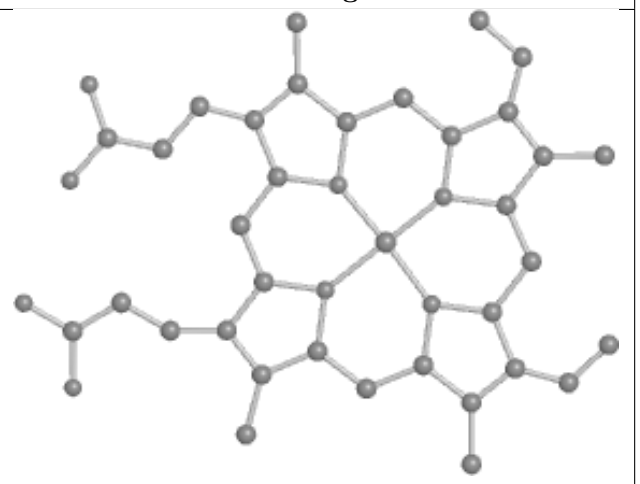

Rings

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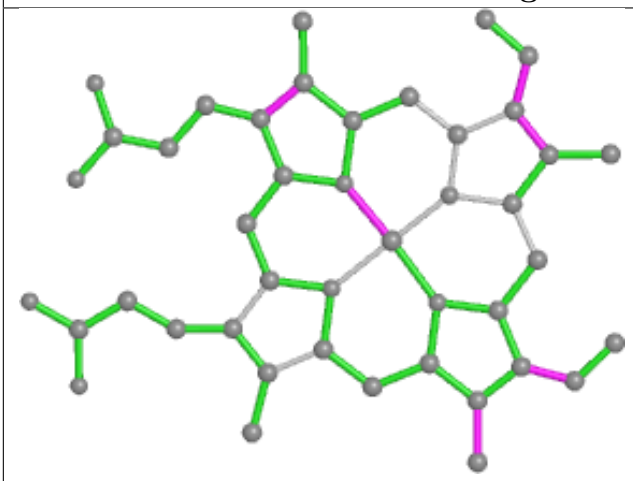

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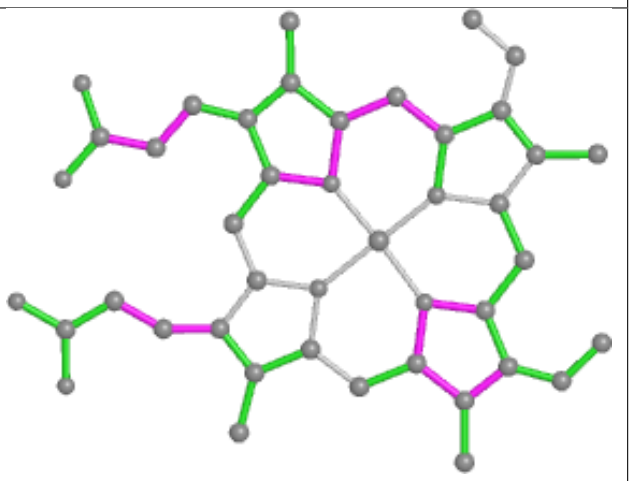

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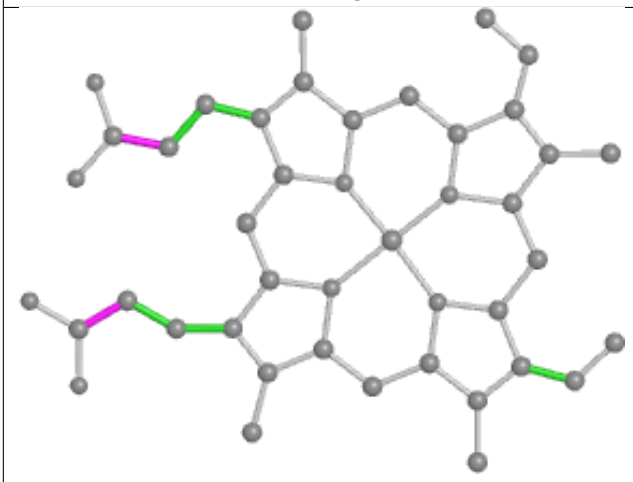
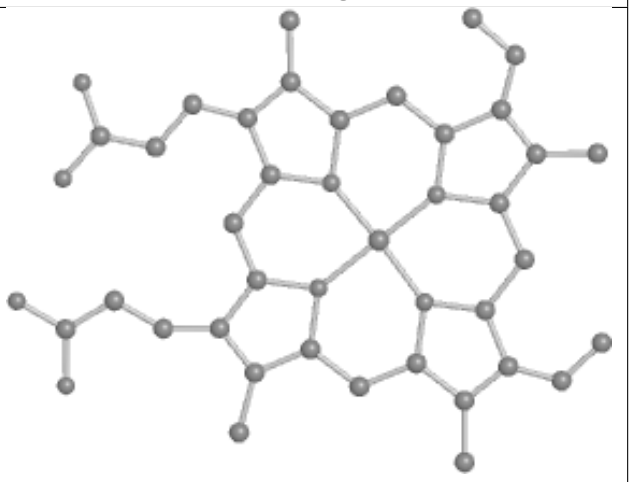
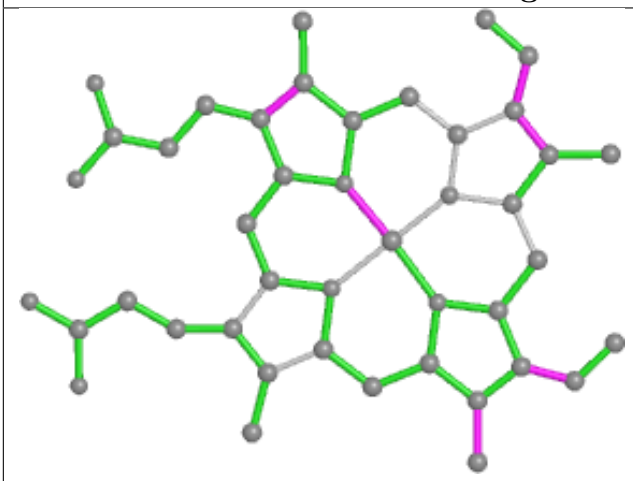


Torsions



Rings

Ligand HEM F 700



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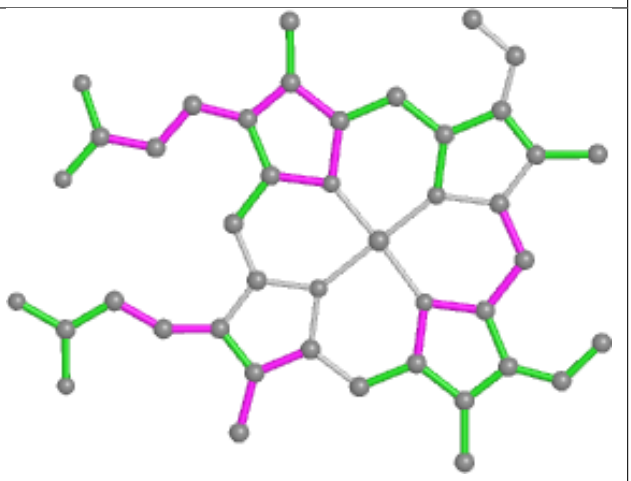
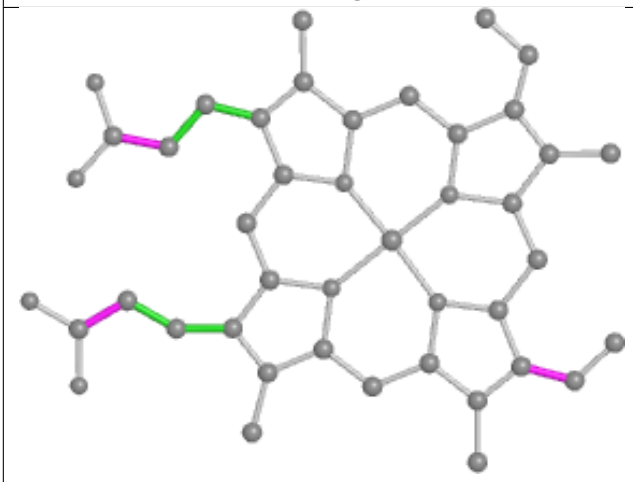
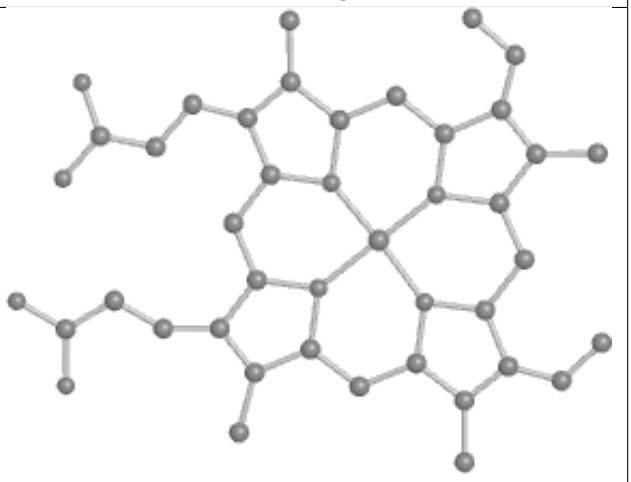


Bond angles

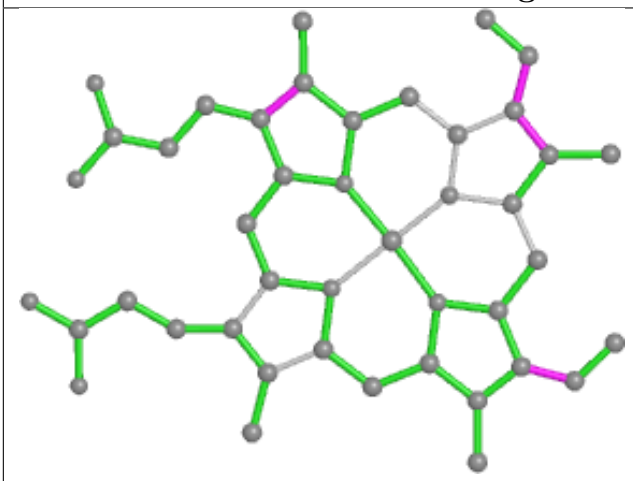


Torsions



Rings

Ligand HEM A 700



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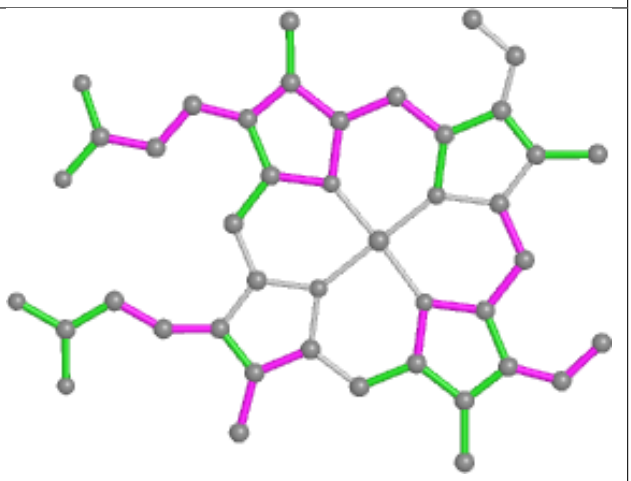
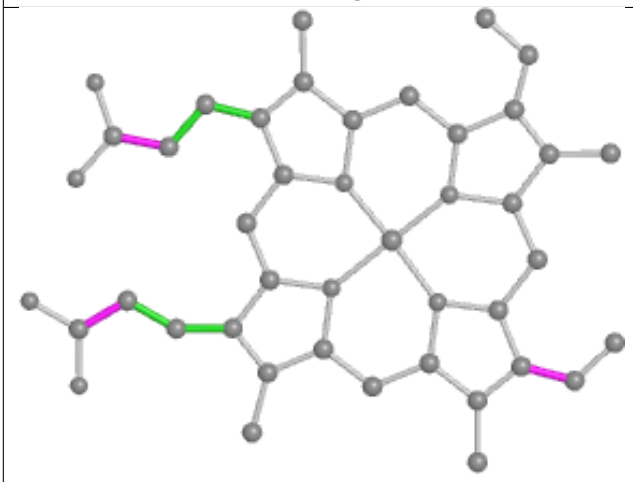
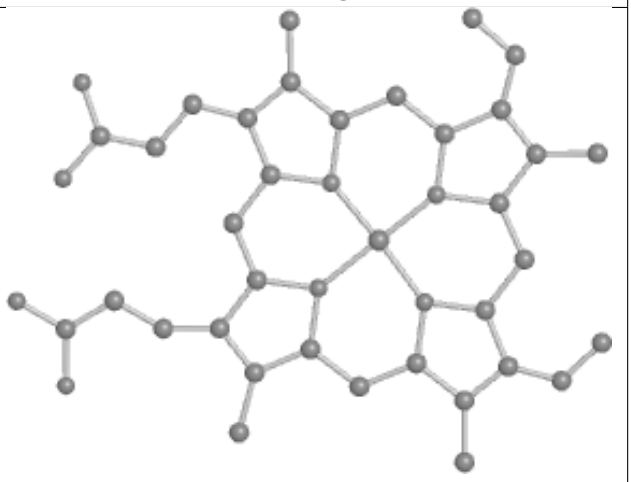


Bond angles

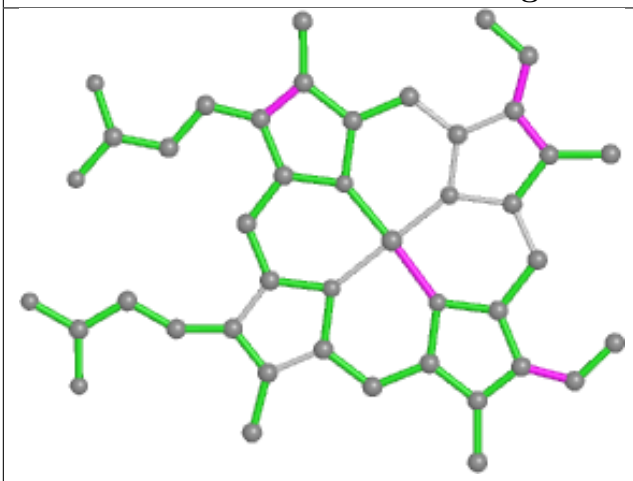


Torsions



Rings

Ligand HEM L 700



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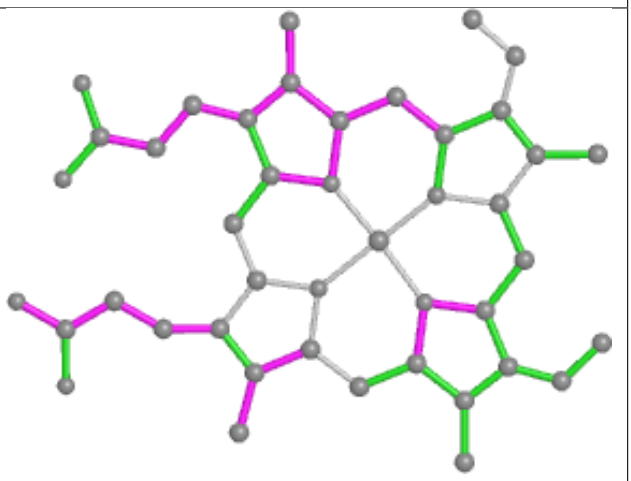

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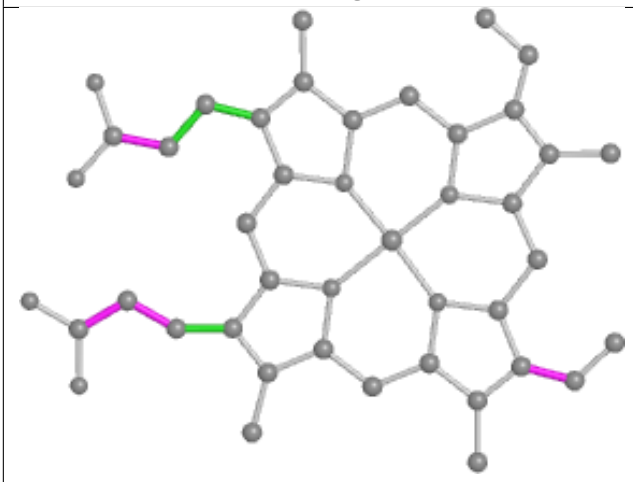
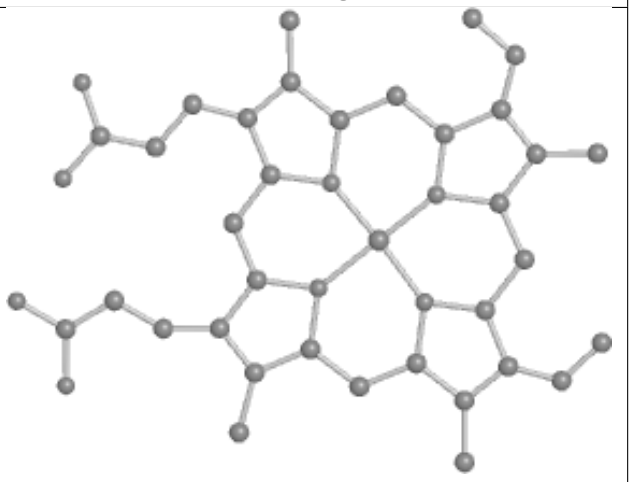


Torsions



Rings

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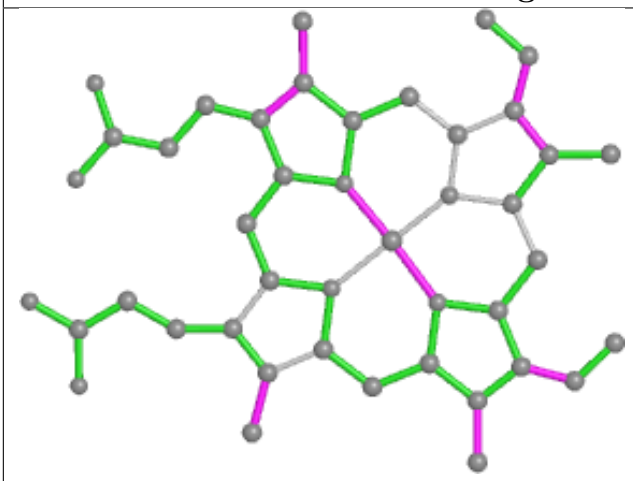

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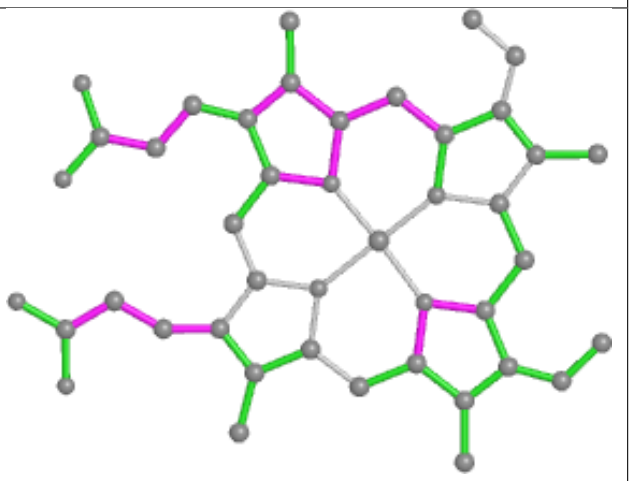
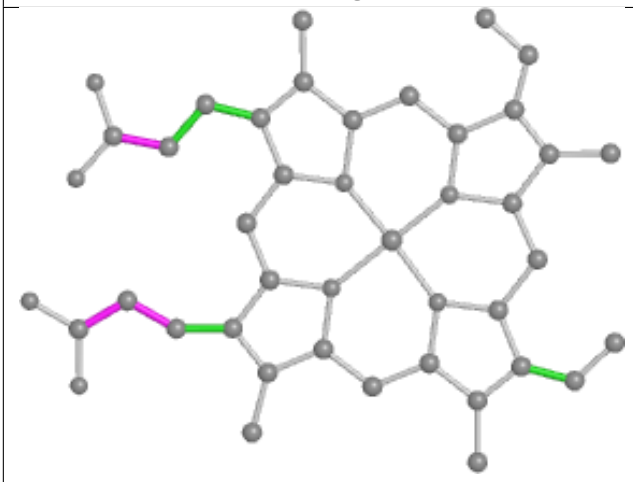


Bond angles



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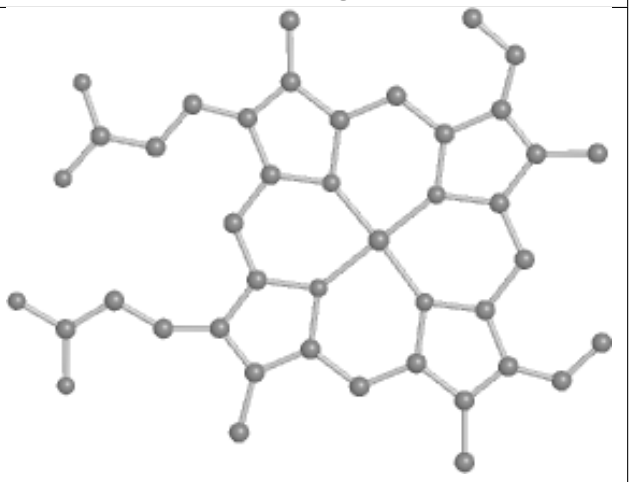
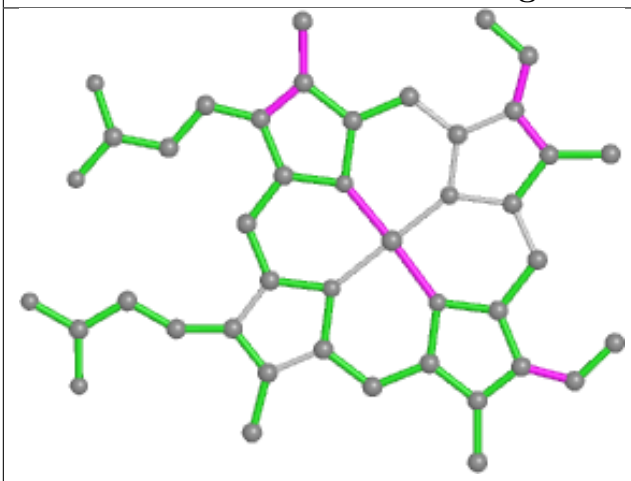


Rings

Ligand HEM C 700



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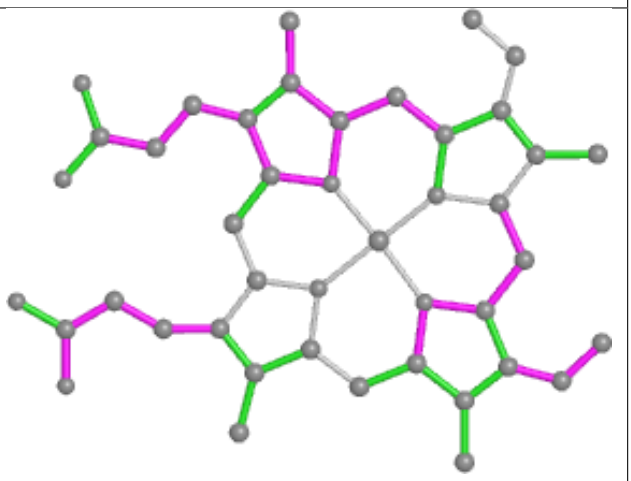

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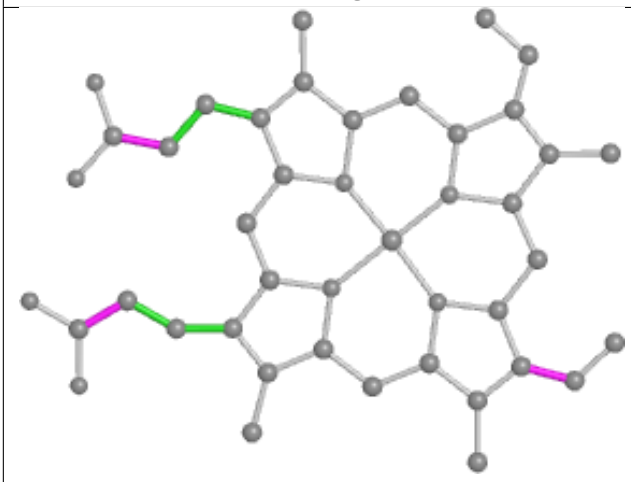

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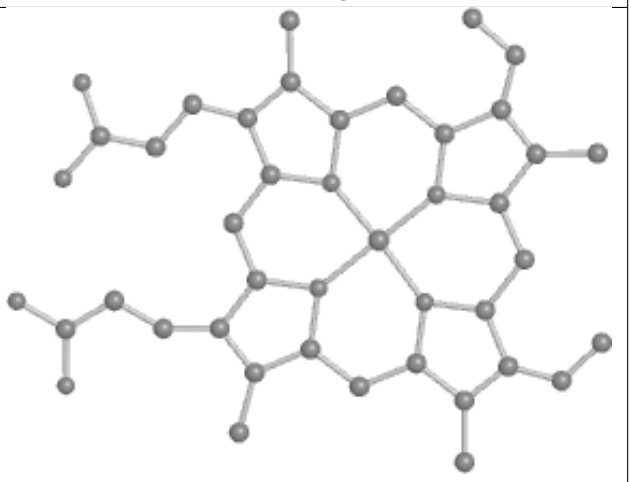

Rings

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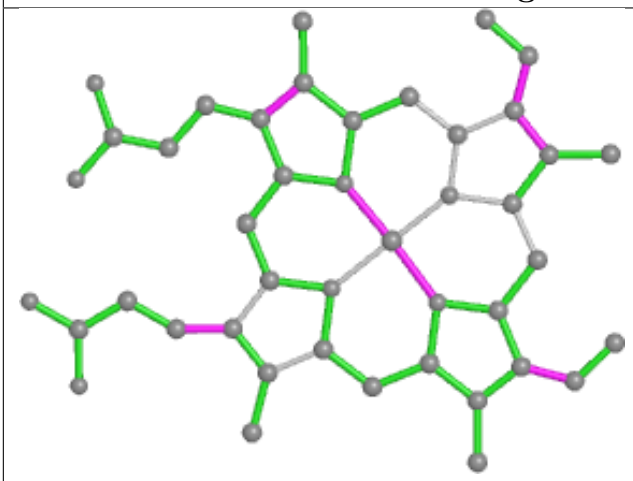

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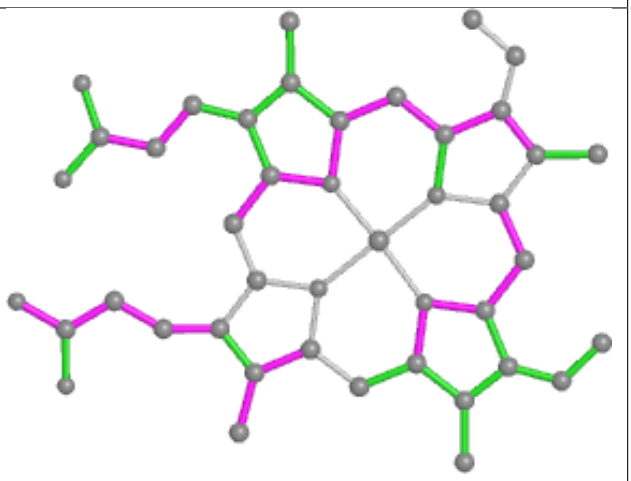
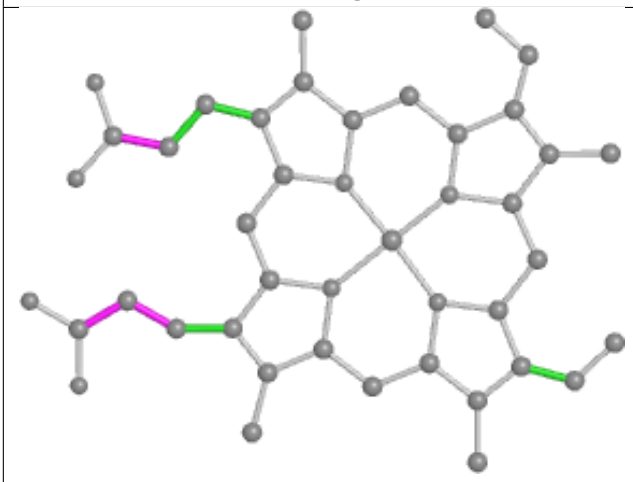
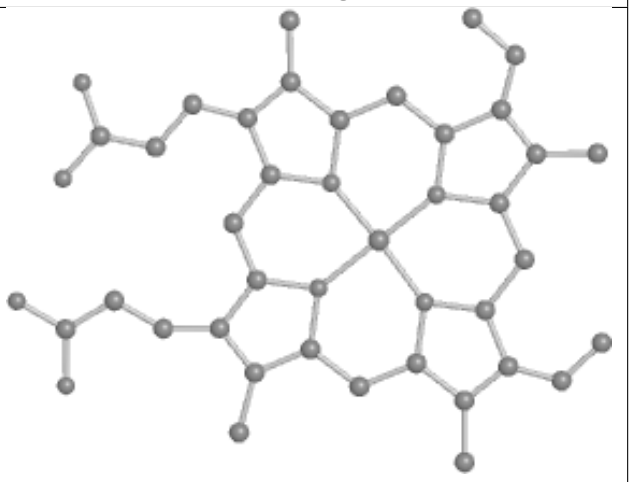


Bond angles



Torsions



Rings

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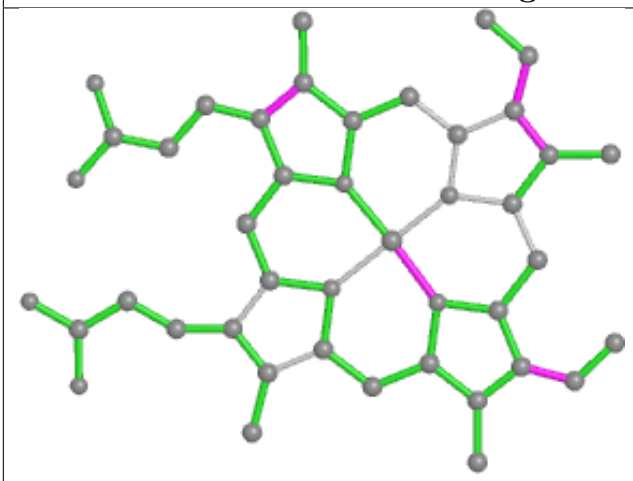

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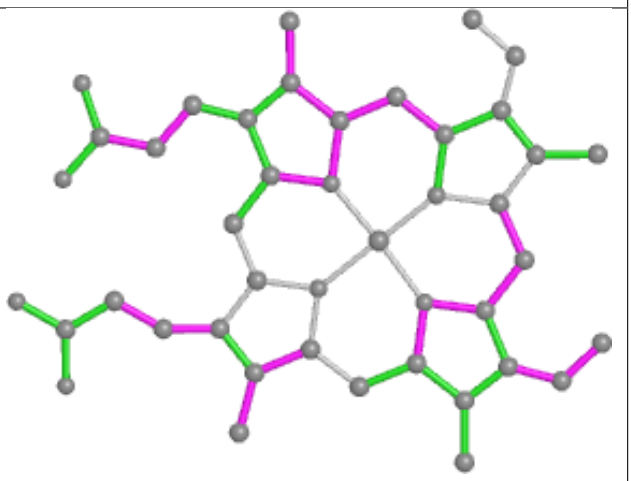
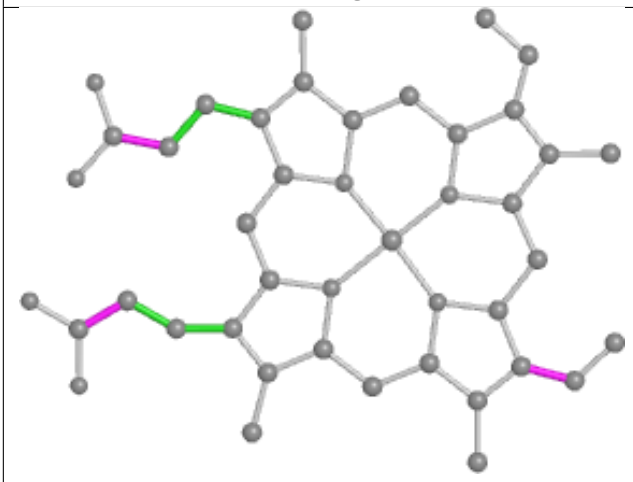
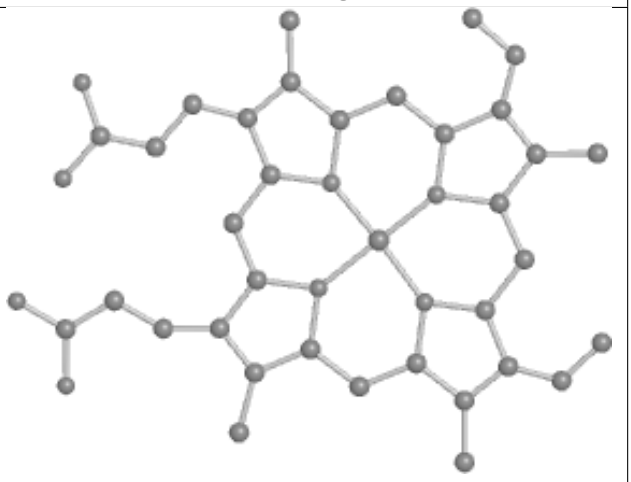


Bond angles

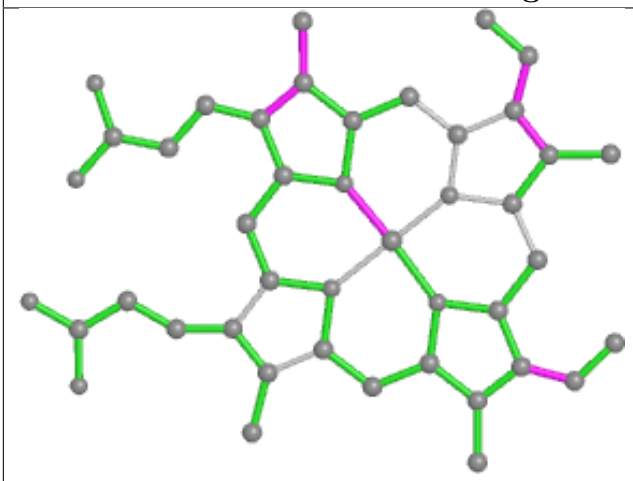


Torsions

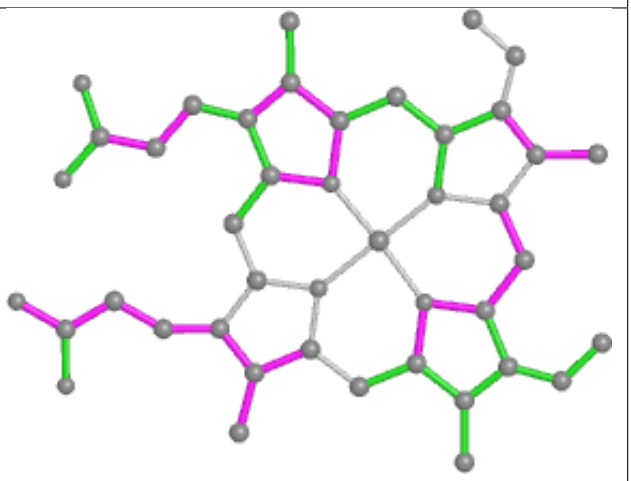


Rings

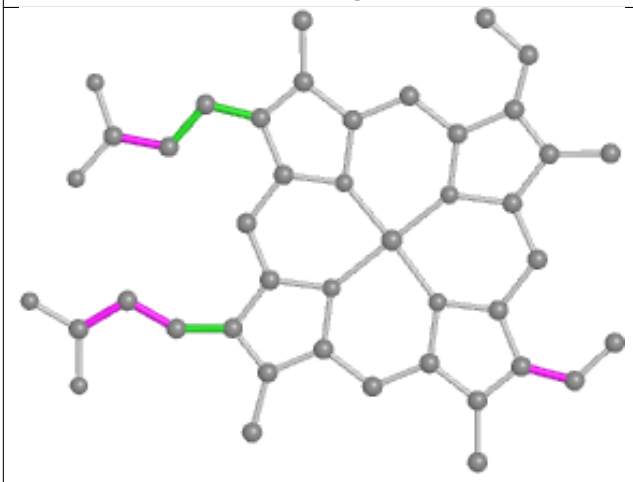
Ligand HEM J 700



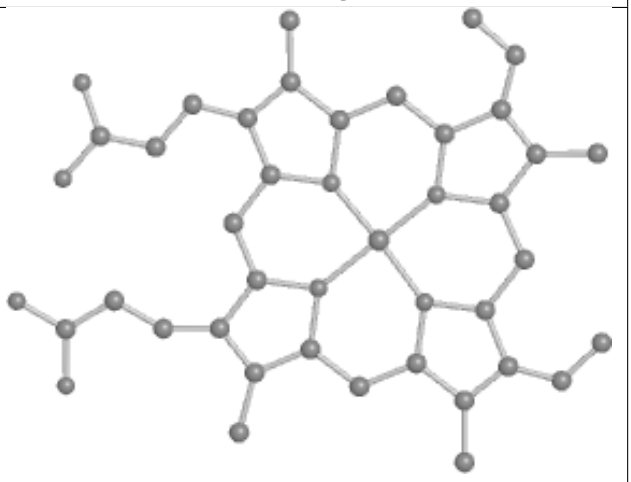
Bond lengths



Bond angles

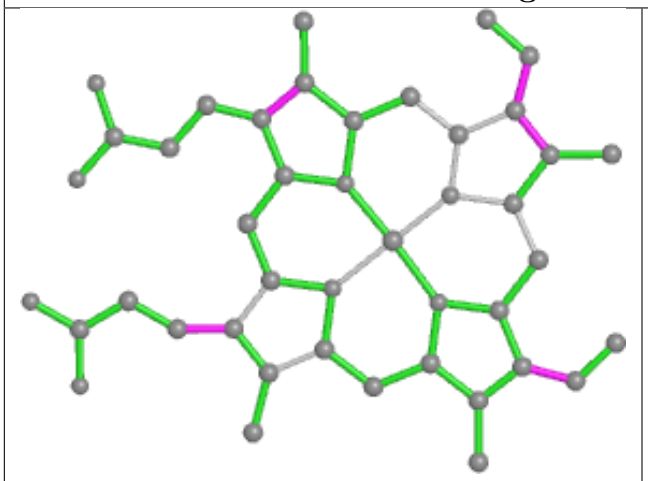


Torsions



Rings

Ligand HEM D 700



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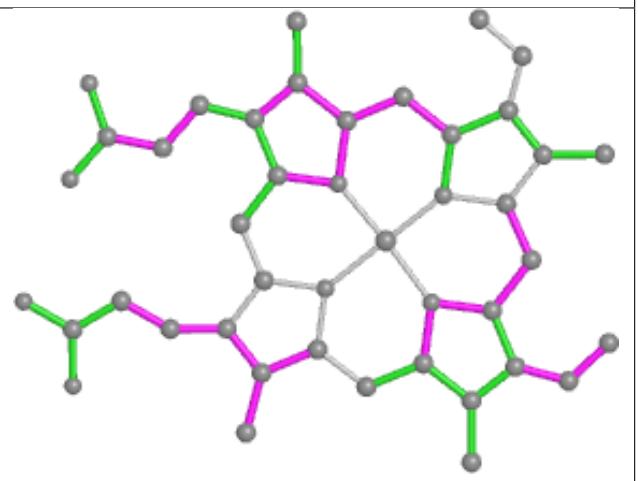
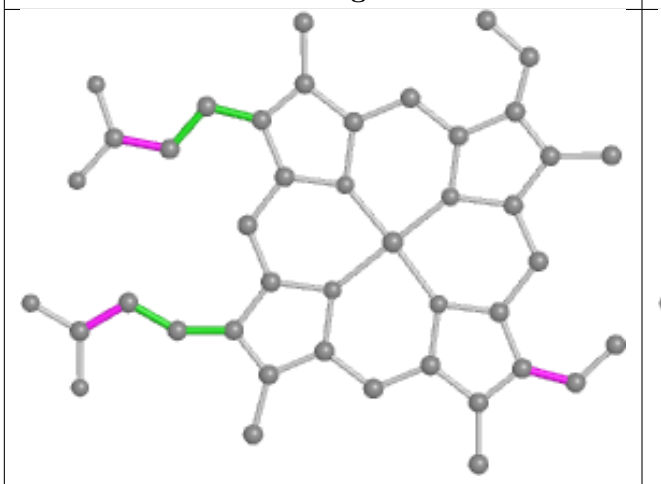
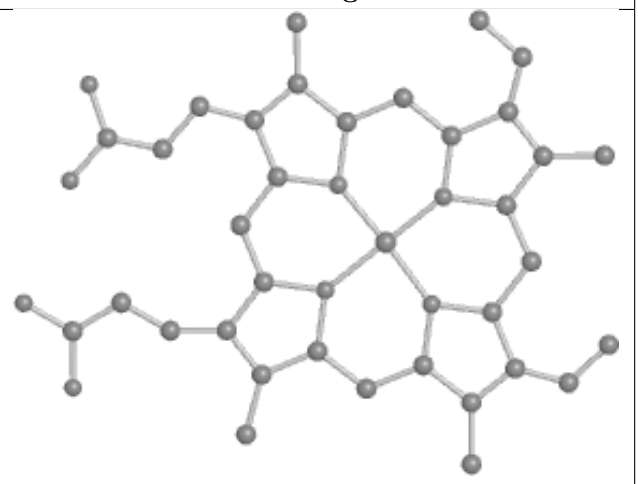


Bond angles



Torsions



Rings

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	126/128 (98%)	0.10	4 (3%)	47	55	14, 22, 49, 67	1 (0%)
1	B	127/128 (99%)	0.12	2 (1%)	72	77	16, 24, 44, 74	2 (1%)
1	C	127/128 (99%)	-0.01	1 (0%)	86	89	13, 19, 40, 66	2 (1%)
1	D	126/128 (98%)	0.06	4 (3%)	47	55	15, 23, 49, 66	2 (1%)
1	E	126/128 (98%)	0.19	6 (4%)	30	38	19, 26, 55, 76	0
1	F	126/128 (98%)	0.08	6 (4%)	30	38	16, 24, 53, 71	1 (0%)
1	G	127/128 (99%)	-0.18	1 (0%)	86	89	14, 21, 42, 56	0
1	H	127/128 (99%)	-0.13	2 (1%)	72	77	18, 25, 41, 56	0
1	I	127/128 (99%)	-0.25	0	100	100	14, 22, 39, 54	0
1	J	126/128 (98%)	0.37	5 (3%)	38	45	19, 30, 49, 62	2 (1%)
1	K	127/128 (99%)	0.37	8 (6%)	20	26	17, 27, 45, 74	1 (0%)
1	L	126/128 (98%)	0.28	10 (7%)	12	18	17, 30, 49, 67	0
All	All	1518/1536 (98%)	0.08	49 (3%)	47	55	13, 25, 49, 76	11 (0%)

All (49) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	2	PRO	8.5
1	K	2	PRO	6.4
1	L	3	LYS	5.4
1	C	2	PRO	5.1
1	F	39	ASP	5.0
1	D	39	ASP	4.5
1	E	40	ASP	4.2
1	D	40	ASP	4.1
1	K	3	LYS	4.1
1	E	39	ASP	3.9
1	L	102	GLU	3.8

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Mol	Chain	Res	Type	RSRZ
1	E	37	PRO	3.7
1	E	128	PHE	3.6
1	A	39	ASP	3.6
1	L	107	GLU	3.6
1	J	102	GLU	3.6
1	D	38	GLU	3.4
1	L	101	SER	3.2
1	L	110	ARG	3.1
1	K	106	ASP	3.0
1	K	102	GLU	3.0
1	F	40	ASP	3.0
1	H	2	PRO	3.0
1	F	38	GLU	3.0
1	A	3	LYS	3.0
1	L	105	ASP	3.0
1	E	38	GLU	2.8
1	K	110	ARG	2.8
1	J	104	LEU	2.8
1	F	33	ARG	2.8
1	K	33[A]	ARG	2.8
1	L	104	LEU	2.7
1	A	40	ASP	2.5
1	H	110	ARG	2.5
1	F	37	PRO	2.4
1	F	34	ARG	2.3
1	J	4	SER	2.3
1	L	108	HIS	2.3
1	E	70	ARG	2.3
1	B	3	LYS	2.2
1	L	9	VAL	2.2
1	J	107	GLU	2.2
1	D	33	ARG	2.2
1	K	107	GLU	2.2
1	G	102	GLU	2.2
1	K	101	SER	2.1
1	L	106	ASP	2.1
1	A	38	GLU	2.1
1	J	58	GLY	2.1

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

6.4 Ligands (i)

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	SO4	B	801	5/5	0.73	0.31	14,26,40,52	5
3	SO4	E	802	5/5	0.73	0.56	20,20,20,20	0
3	SO4	L	803	5/5	0.73	0.17	58,64,71,85	0
3	SO4	L	801	5/5	0.75	0.20	35,42,44,57	5
3	SO4	K	801	5/5	0.77	0.16	59,62,71,82	0
3	SO4	J	801	5/5	0.78	0.18	48,62,71,78	0
3	SO4	G	801	5/5	0.79	0.25	33,40,48,55	5
3	SO4	L	802	5/5	0.80	0.21	32,40,44,51	5
3	SO4	I	801	5/5	0.82	0.31	22,24,47,49	5
3	SO4	F	803	5/5	0.85	0.57	20,20,20,20	0
3	SO4	A	801	5/5	0.85	0.25	29,30,40,49	5
3	SO4	H	801	5/5	0.85	0.25	29,31,36,47	5
3	SO4	D	802	5/5	0.86	0.17	46,52,60,69	0
3	SO4	A	803	5/5	0.88	0.35	18,27,33,34	5
3	SO4	F	801	5/5	0.90	0.38	24,31,32,38	5
3	SO4	E	801	5/5	0.90	0.15	21,23,28,29	5
3	SO4	F	802	5/5	0.92	0.23	20,25,26,34	5
3	SO4	B	803	5/5	0.94	0.54	21,35,40,43	5
3	SO4	D	801	5/5	0.94	0.23	38,38,45,55	0
3	SO4	A	804	5/5	0.95	0.12	42,53,68,68	0
3	SO4	A	802	5/5	0.95	0.21	20,22,27,28	5
2	CYN	F	800	2/2	0.97	0.12	22,22,22,24	0
3	SO4	B	802	5/5	0.97	0.10	25,32,34,37	0
3	SO4	G	802	5/5	0.98	0.10	22,27,32,33	0
2	CYN	G	800	2/2	0.98	0.12	19,19,19,22	0
4	HEM	B	700	43/43	0.98	0.13	13,17,25,29	0
4	HEM	E	700	43/43	0.98	0.11	14,20,29,39	0
4	HEM	F	700	43/43	0.98	0.11	10,17,30,33	0
4	HEM	G	700	43/43	0.98	0.12	10,18,29,44	0
4	HEM	I	700	43/43	0.98	0.11	13,17,31,36	0
4	HEM	J	700	43/43	0.98	0.12	16,21,32,46	0
4	HEM	K	700	43/43	0.98	0.12	11,19,34,49	0

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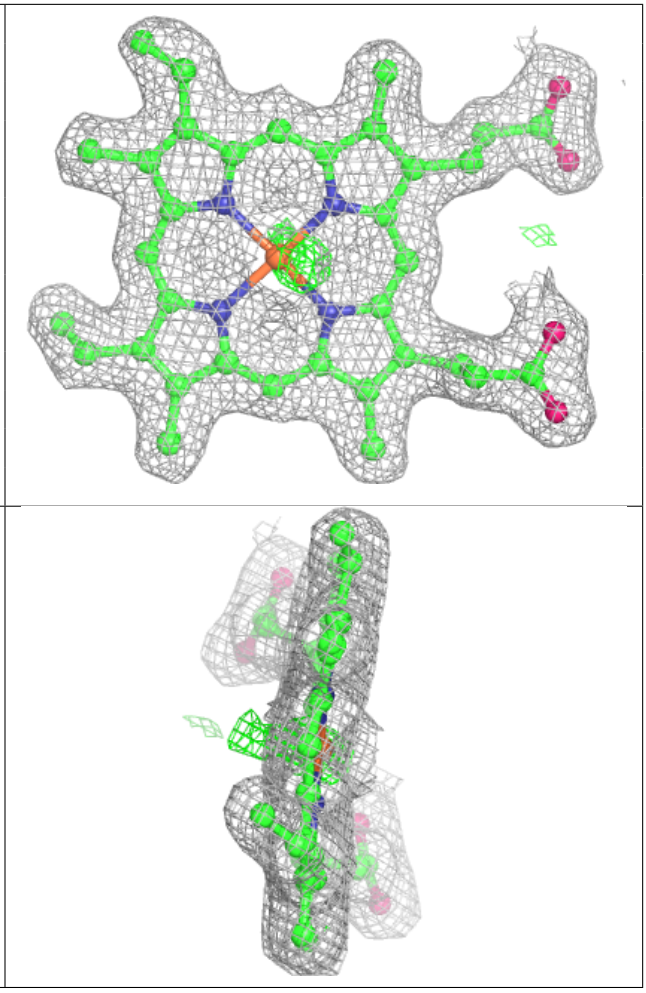
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
4	HEM	L	700	43/43	0.98	0.11	14,21,30,37	0
4	HEM	A	700	43/43	0.99	0.12	10,16,28,32	0
2	CYN	L	800	2/2	0.99	0.10	21,21,21,24	0
4	HEM	C	700	43/43	0.99	0.12	8,15,22,25	0
4	HEM	D	700	43/43	0.99	0.11	11,17,28,32	0
2	CYN	E	800	2/2	0.99	0.10	18,18,18,30	0
2	CYN	C	800	2/2	0.99	0.11	14,14,14,16	0
2	CYN	D	800	2/2	0.99	0.12	18,18,18,27	0
4	HEM	H	700	43/43	0.99	0.10	13,19,32,41	0
2	CYN	H	800	2/2	0.99	0.13	15,15,15,22	0
2	CYN	I	800	2/2	0.99	0.10	18,18,18,29	0
2	CYN	J	800	2/2	0.99	0.10	22,22,22,24	0
2	CYN	K	800	2/2	0.99	0.14	20,20,20,30	0
2	CYN	B	800	2/2	1.00	0.11	18,18,18,29	0
2	CYN	A	800	2/2	1.00	0.10	16,16,16,18	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.

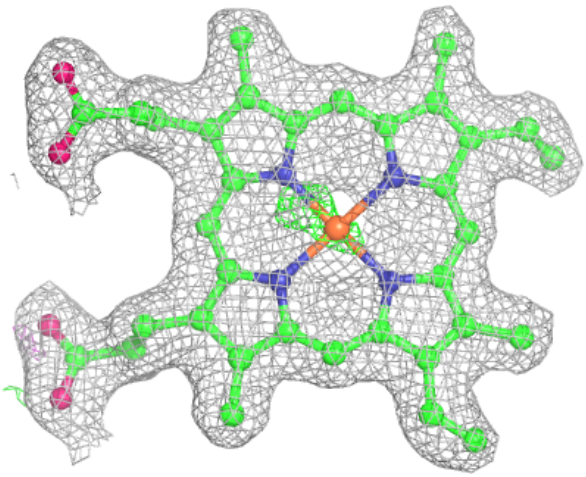
Electron density around HEM B 700:

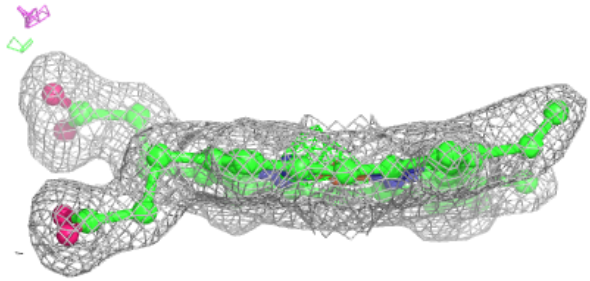
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

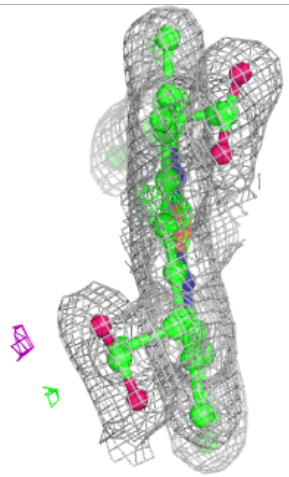


Electron density around HEM E 700:

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

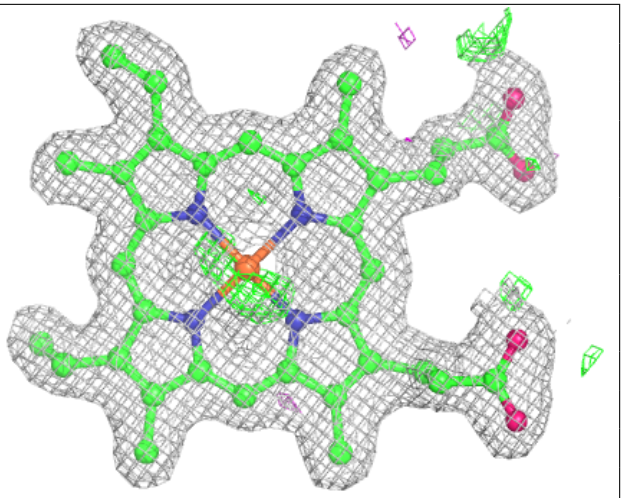


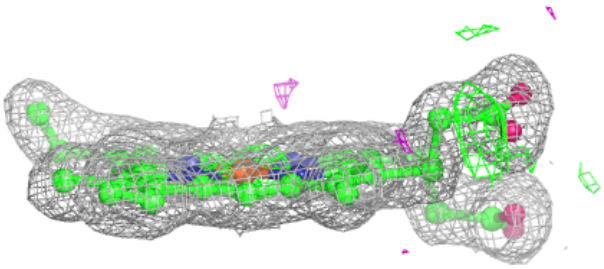


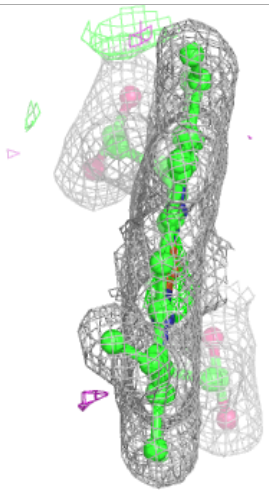


Electron density around HEM F 700:

2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

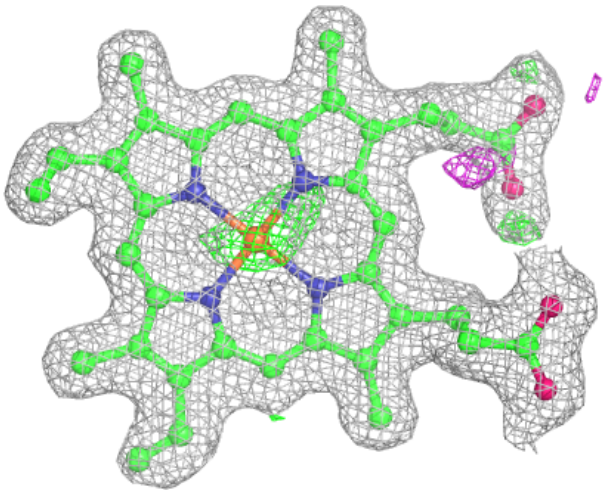


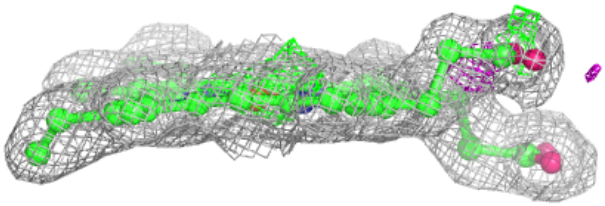


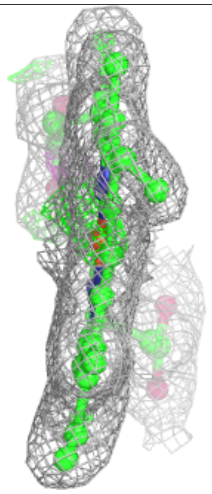


Electron density around HEM G 700:

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

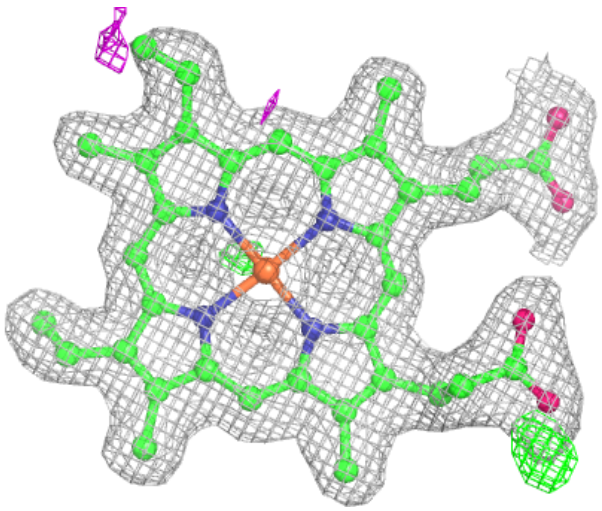


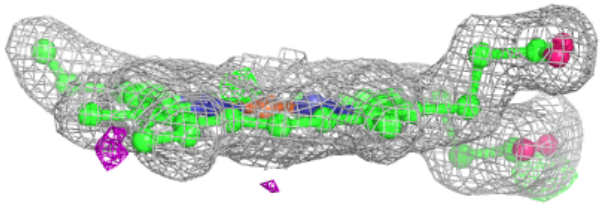


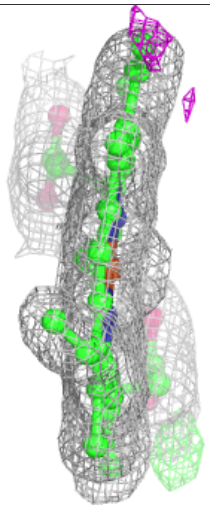


Electron density around HEM I 700:

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

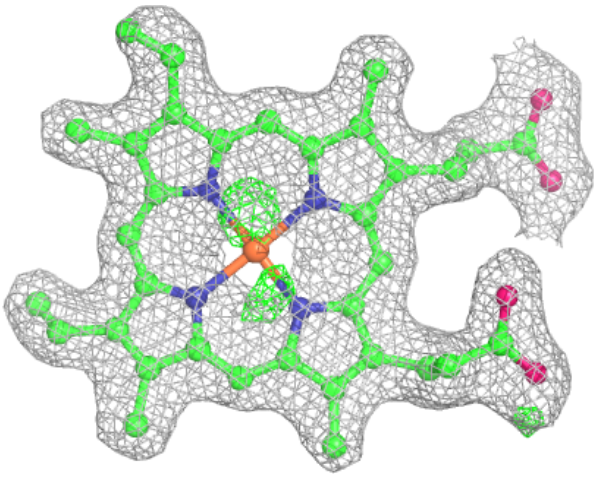


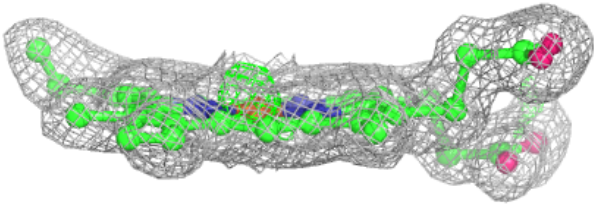


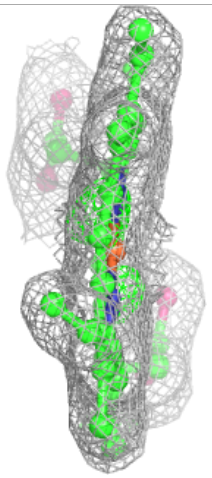


Electron density around HEM J 700:

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

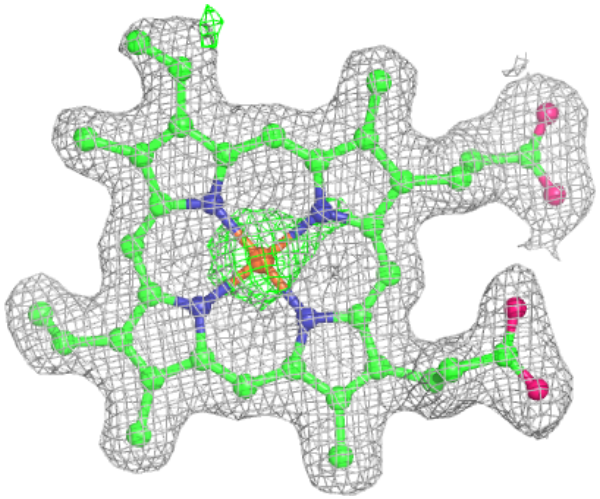


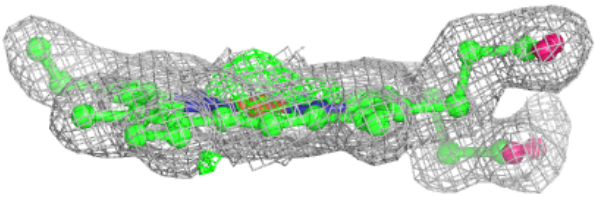


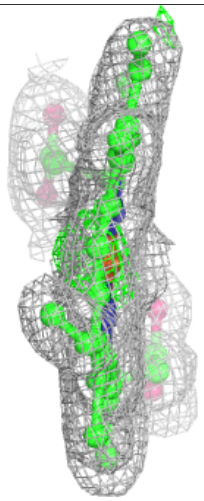


Electron density around HEM K 700:

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

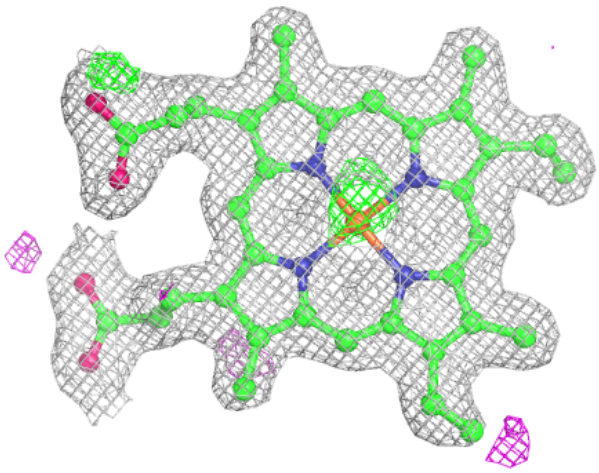


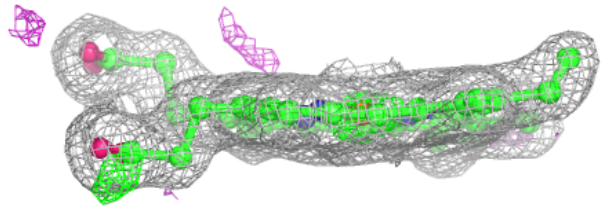




Electron density around HEM L 700:

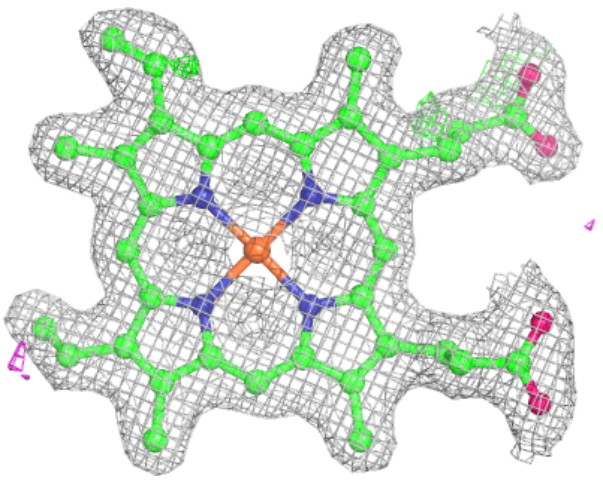
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

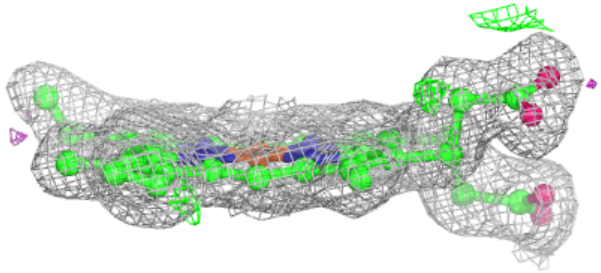


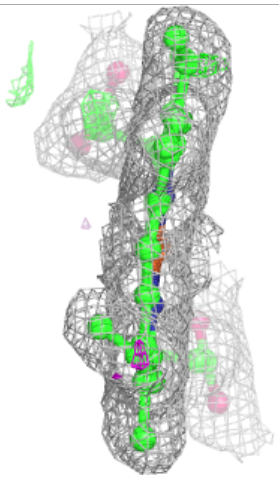


Electron density around HEM A 700:

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

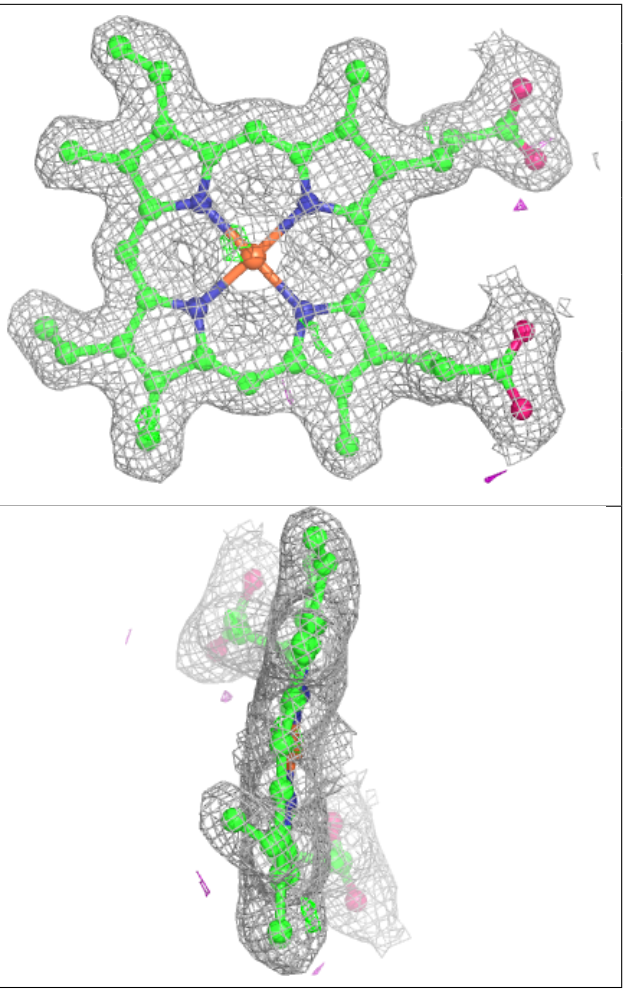






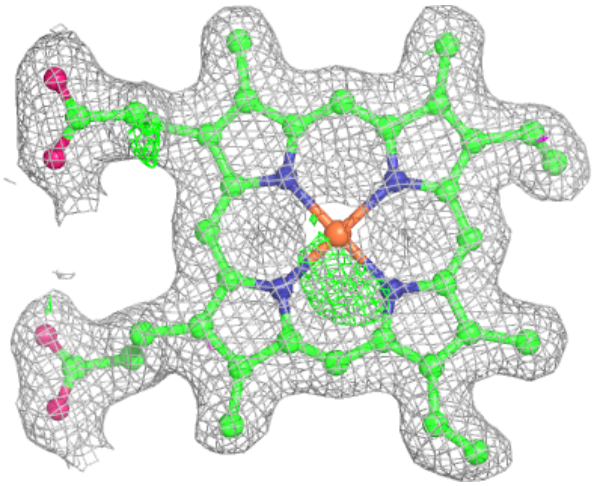
Electron density around HEM C 700:

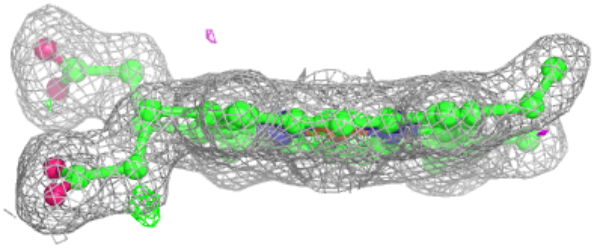
$2mF_o$-DF_c (at 0.7 rmsd) in gray
mF_o-DF_c (at 3 rmsd) in purple (negative)
and green (positive)

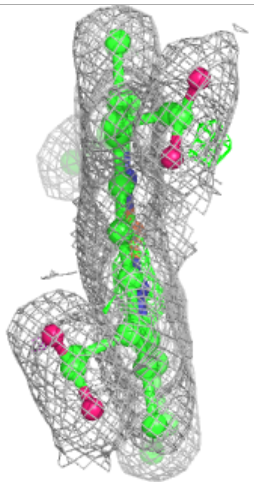


Electron density around HEM D 700:

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

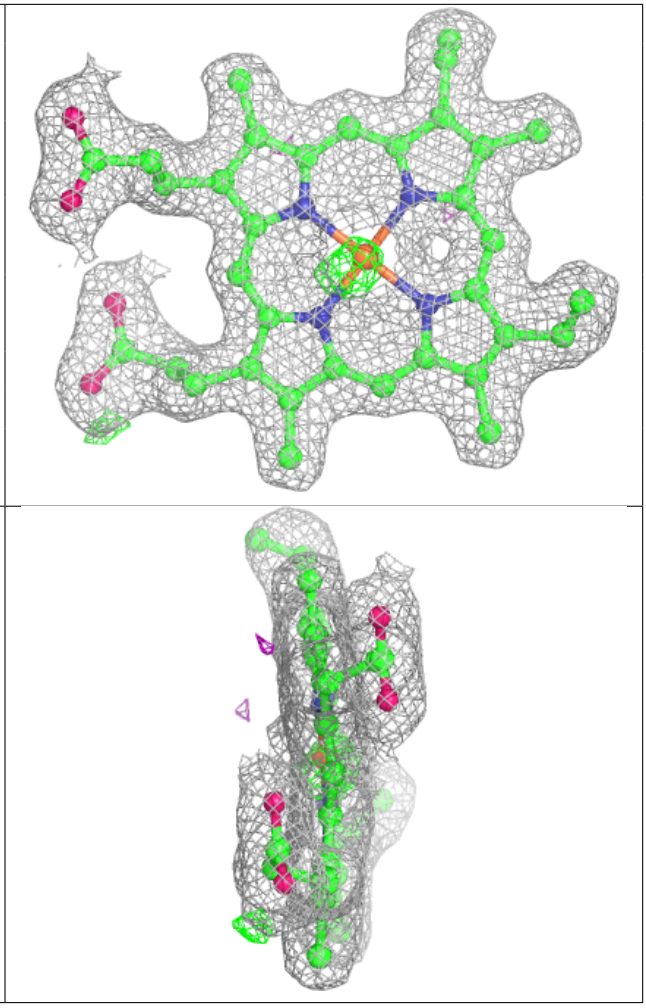






Electron density around HEM H 700:

$2mF_o$-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.