Full wwPDB X-ray Structure Validation Report ⓘ

Aug 24, 2026 – 03:00 am BST

PDB ID	:	2Y46 / pdb_00002y46
Title	:	Structure of the mixed-function P450 MycG in complex with mycinamicin IV in C 2 2 21 space group
Authors	:	Li, S.; Kells, P.M.; Sherman, D.H.; Podust, L.M.
Deposited on	:	2011-01-05
Resolution	:	1.83 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 ⓘ symbol.

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

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

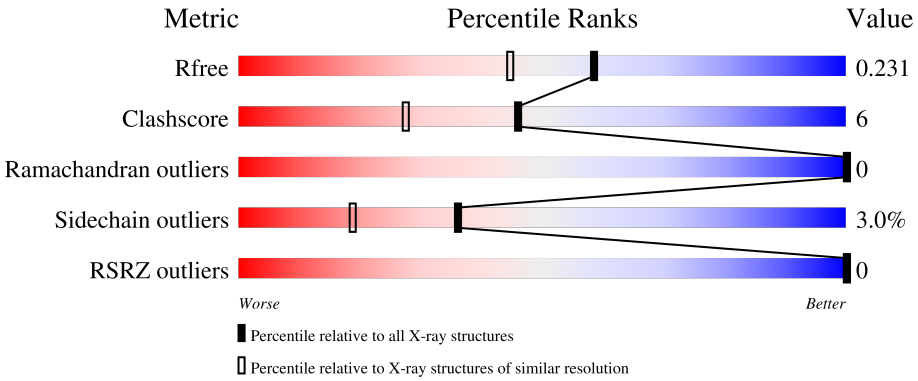
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
Ideal 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 1.83 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}	180053	1296 (1.84-1.84)
Clashscore	190562	1329 (1.84-1.84)
Ramachandran outliers	187476	1318 (1.84-1.84)
Sidechain outliers	187428	1318 (1.84-1.84)
RSRZ outliers	180081	1296 (1.84-1.84)

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	417	81% 12% • 6%
1	B	417	85% 9% • 6%
1	C	417	84% 9% • 6%

2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 10665 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 P-450-LIKE PROTEIN.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	393	Total	C	N	O	S	0	7	0
			3131	1964	572	584	11			
1	B	393	Total	C	N	O	S	0	7	0
			3143	1971	576	585	11			
1	C	393	Total	C	N	O	S	0	6	0
			3136	1967	576	582	11			

There are 60 discrepancies between the modelled and reference sequences:

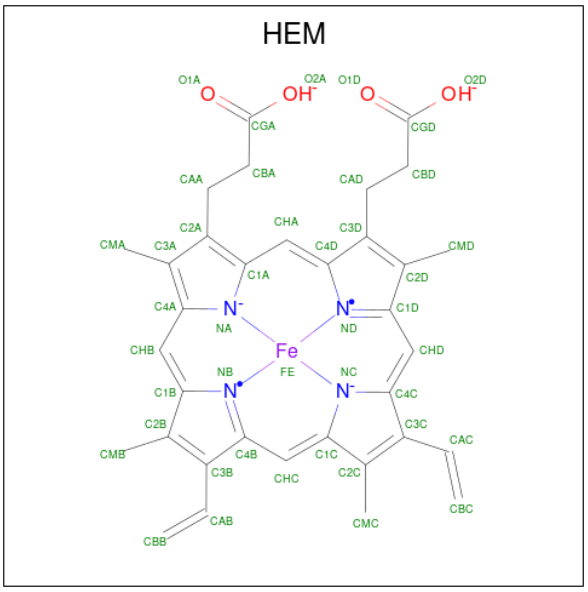
Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	expression tag	UNP Q59523
A	-18	GLY	-	expression tag	UNP Q59523
A	-17	SER	-	expression tag	UNP Q59523
A	-16	SER	-	expression tag	UNP Q59523
A	-15	HIS	-	expression tag	UNP Q59523
A	-14	HIS	-	expression tag	UNP Q59523
A	-13	HIS	-	expression tag	UNP Q59523
A	-12	HIS	-	expression tag	UNP Q59523
A	-11	HIS	-	expression tag	UNP Q59523
A	-10	HIS	-	expression tag	UNP Q59523
A	-9	SER	-	expression tag	UNP Q59523
A	-8	SER	-	expression tag	UNP Q59523
A	-7	GLY	-	expression tag	UNP Q59523
A	-6	LEU	-	expression tag	UNP Q59523
A	-5	VAL	-	expression tag	UNP Q59523
A	-4	PRO	-	expression tag	UNP Q59523
A	-3	ARG	-	expression tag	UNP Q59523
A	-2	GLY	-	expression tag	UNP Q59523
A	-1	SER	-	expression tag	UNP Q59523
A	0	HIS	-	expression tag	UNP Q59523
B	-19	MET	-	expression tag	UNP Q59523
B	-18	GLY	-	expression tag	UNP Q59523
B	-17	SER	-	expression tag	UNP Q59523

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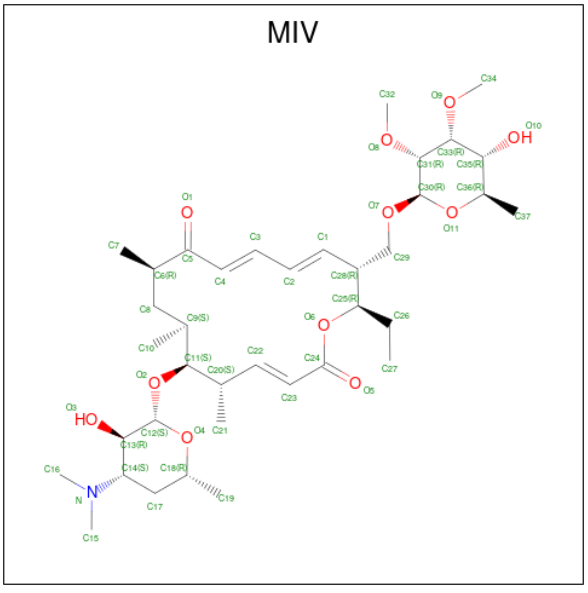
Chain	Residue	Modelled	Actual	Comment	Reference
B	-16	SER	-	expression tag	UNP Q59523
B	-15	HIS	-	expression tag	UNP Q59523
B	-14	HIS	-	expression tag	UNP Q59523
B	-13	HIS	-	expression tag	UNP Q59523
B	-12	HIS	-	expression tag	UNP Q59523
B	-11	HIS	-	expression tag	UNP Q59523
B	-10	HIS	-	expression tag	UNP Q59523
B	-9	SER	-	expression tag	UNP Q59523
B	-8	SER	-	expression tag	UNP Q59523
B	-7	GLY	-	expression tag	UNP Q59523
B	-6	LEU	-	expression tag	UNP Q59523
B	-5	VAL	-	expression tag	UNP Q59523
B	-4	PRO	-	expression tag	UNP Q59523
B	-3	ARG	-	expression tag	UNP Q59523
B	-2	GLY	-	expression tag	UNP Q59523
B	-1	SER	-	expression tag	UNP Q59523
B	0	HIS	-	expression tag	UNP Q59523
C	-19	MET	-	expression tag	UNP Q59523
C	-18	GLY	-	expression tag	UNP Q59523
C	-17	SER	-	expression tag	UNP Q59523
C	-16	SER	-	expression tag	UNP Q59523
C	-15	HIS	-	expression tag	UNP Q59523
C	-14	HIS	-	expression tag	UNP Q59523
C	-13	HIS	-	expression tag	UNP Q59523
C	-12	HIS	-	expression tag	UNP Q59523
C	-11	HIS	-	expression tag	UNP Q59523
C	-10	HIS	-	expression tag	UNP Q59523
C	-9	SER	-	expression tag	UNP Q59523
C	-8	SER	-	expression tag	UNP Q59523
C	-7	GLY	-	expression tag	UNP Q59523
C	-6	LEU	-	expression tag	UNP Q59523
C	-5	VAL	-	expression tag	UNP Q59523
C	-4	PRO	-	expression tag	UNP Q59523
C	-3	ARG	-	expression tag	UNP Q59523
C	-2	GLY	-	expression tag	UNP Q59523
C	-1	SER	-	expression tag	UNP Q59523
C	0	HIS	-	expression tag	UNP Q59523

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



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

- Molecule 3 is MYCINAMICIN IV (CCD ID: MIV) (formula: $C_{37}H_{61}NO_{11}$).



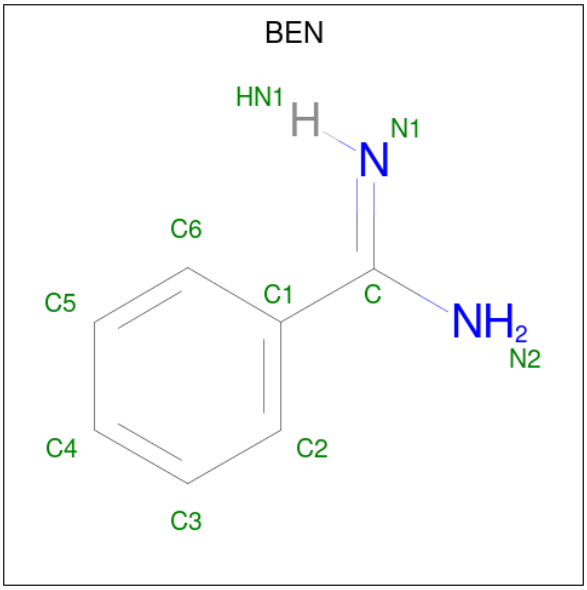
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
3	A	1	Total	C	N	O	0	0
			49	37	1	11		

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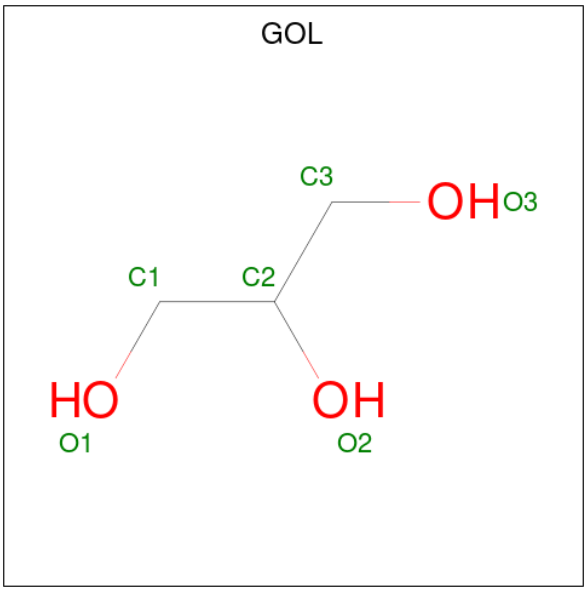
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	B	1	Total C N O 49 37 1 11	0	0
3	C	1	Total C N O 49 37 1 11	0	0

- Molecule 4 is BENZAMIDINE (CCD ID: BEN) (formula: $C_7H_8N_2$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C N 9 7 2	0	0
4	B	1	Total C N 9 7 2	0	0
4	C	1	Total C N 9 7 2	0	0

- Molecule 5 is GLYCEROL (CCD ID: GOL) (formula: $C_3H_8O_3$).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 6 3 3	0	0
5	A	1	Total C O 6 3 3	0	0
5	B	1	Total C O 6 3 3	0	0
5	B	1	Total C O 6 3 3	0	0
5	C	1	Total C O 6 3 3	0	0
5	C	1	Total C O 6 3 3	0	0

- Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	A	292	Total O 292 292	0	0
6	B	313	Total O 313 313	0	0
6	C	311	Total O 311 311	0	0

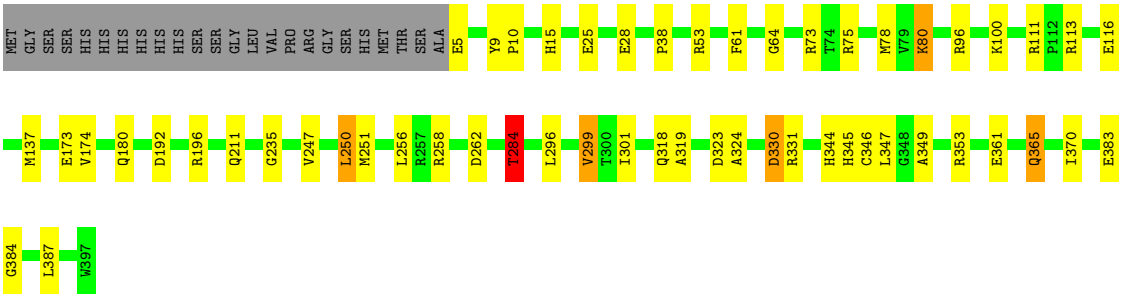
3 Residue-property plots (i)

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: P-450-LIKE PROTEIN

Chain A:

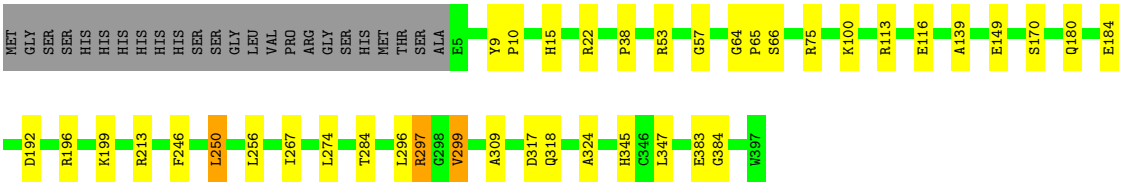




• Molecule 1: P-450-LIKE PROTEIN

Chain B:

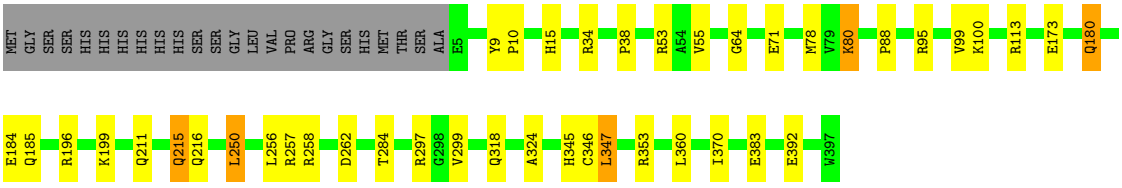




• Molecule 1: P-450-LIKE PROTEIN

Chain C:





4 Data and refinement statistics (i)

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α, β, γ	58.27Å 101.01Å 441.53Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	220.77 – 1.83 220.77 – 1.83	Depositor EDS
% Data completeness (in resolution range)	82.8 (220.77-1.83) 82.7 (220.77-1.83)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.52 (at 1.83Å)	Xtriage
Refinement program	REFMAC 5.5.0109	Depositor
R, R_{free}	0.157 , 0.233 0.157 , 0.231	Depositor DCC
R_{free} test set	4822 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å^2)	20.9	Xtriage
Anisotropy	0.167	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.34 , 21.5	EDS
L-test for twinning[2]	$< \|L\| > = 0.51$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	0.479 for 1/2*h-1/2*k,-3/2*h-1/2*k,-l 0.479 for 1/2*h+1/2*k,3/2*h-1/2*k,-l	Xtriage
F_o,F_c correlation	0.97	EDS
Total number of atoms	10665	wwPDB-VP
Average B, all atoms (Å^2)	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.48% 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

5.1 Standard geometry

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

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	1.29	7/3197 (0.2%)	1.13	4/4351 (0.1%)
1	B	1.28	7/3209 (0.2%)	1.12	1/4365 (0.0%)
1	C	1.26	2/3202 (0.1%)	1.11	1/4356 (0.0%)
All	All	1.27	16/9608 (0.2%)	1.12	6/13072 (0.0%)

All (16) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	64	GLY	N-CA	8.28	1.52	1.45
1	A	64	GLY	N-CA	6.75	1.53	1.44
1	B	309	ALA	CA-CB	6.32	1.61	1.53
1	C	64	GLY	N-CA	6.23	1.53	1.44
1	B	267	ILE	CA-CB	6.09	1.57	1.54
1	B	139	ALA	CA-CB	5.91	1.62	1.53
1	B	64	GLY	CA-C	5.90	1.56	1.51
1	B	57	GLY	N-CA	5.77	1.52	1.45
1	A	387	LEU	N-CA	5.76	1.52	1.45
1	A	319	ALA	CA-CB	5.38	1.62	1.53
1	A	301	ILE	C-O	-5.22	1.18	1.24
1	A	370	ILE	CA-CB	5.08	1.60	1.54
1	B	317	ASP	C-O	5.06	1.30	1.24
1	C	53	ARG	CG-CD	5.05	1.67	1.52
1	A	61	PHE	C-O	5.05	1.29	1.24
1	A	111	ARG	N-CA	5.04	1.51	1.46

All (6) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	C	55	VAL	N-CA-C	6.00	116.18	110.42
1	A	349	ALA	CA-C-N	-5.78	112.64	119.05

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	349	ALA	C-N-CA	-5.78	112.64	119.05
1	B	297	ARG	CA-CB-CG	-5.56	102.99	114.10
1	A	28	GLU	CB-CA-C	-5.24	101.98	110.79
1	A	284	THR	CB-CA-C	5.03	118.18	110.14

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	3131	0	3095	42	0
1	B	3143	0	3116	25	0
1	C	3136	0	3115	41	0
2	A	43	0	30	4	0
2	B	43	0	30	1	0
2	C	43	0	30	1	0
3	A	49	0	61	1	0
3	B	49	0	61	2	0
3	C	49	0	61	0	0
4	A	9	0	7	0	0
4	B	9	0	7	0	0
4	C	9	0	7	1	0
5	A	12	0	16	1	0
5	B	12	0	16	1	0
5	C	12	0	16	0	0
6	A	292	0	0	9	0
6	B	313	0	0	6	1
6	C	311	0	0	8	0
All	All	10665	0	9668	113	1

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 6.

All (113) 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:78:MET:HE2	1:A:80:LYS:CE	1.62	1.26
1:A:284:THR:HG22	6:A:2007:HOH:O	1.43	1.17
1:C:78:MET:HE2	1:C:80:LYS:CE	1.77	1.12
1:A:78:MET:HE2	1:A:80:LYS:HE3	1.20	1.11
1:C:78:MET:HE2	1:C:80:LYS:HE3	1.28	1.08
1:A:78:MET:CE	1:A:80:LYS:NZ	2.24	1.00
1:C:78:MET:CE	1:C:80:LYS:NZ	2.25	0.99
1:A:78:MET:HE2	1:A:80:LYS:NZ	1.78	0.98
1:C:78:MET:HE2	1:C:80:LYS:NZ	1.82	0.93
1:C:250:LEU:HD12	1:C:256:LEU:HG	1.51	0.92
1:A:78:MET:CE	1:A:80:LYS:CE	2.51	0.89
1:A:78:MET:CE	1:A:80:LYS:HZ2	1.83	0.89
1:C:78:MET:CE	1:C:80:LYS:HZ2	1.85	0.88
1:A:5:GLU:O	6:A:2003:HOH:O	1.96	0.84
1:C:9:TYR:OH	6:C:2007:HOH:O	1.98	0.81
1:C:78:MET:CE	1:C:80:LYS:CE	2.61	0.78
1:B:9:TYR:OH	6:B:2009:HOH:O	2.01	0.77
1:C:78:MET:HE1	1:C:80:LYS:NZ	1.99	0.77
1:A:78:MET:HE1	1:A:80:LYS:NZ	2.00	0.76
1:C:284:THR:OG1	6:C:2007:HOH:O	2.05	0.74
1:A:284:THR:CG2	6:A:2007:HOH:O	2.15	0.73
1:C:345[B]:HIS:HE1	6:C:2281:HOH:O	1.74	0.70
1:C:78:MET:HE1	1:C:80:LYS:HZ1	1.59	0.66
1:A:9:TYR:OH	6:A:2007:HOH:O	2.10	0.64
1:B:15:HIS:HE1	1:B:383:GLU:OE2	1.81	0.64
1:C:347:LEU:C	1:C:347:LEU:HD12	2.23	0.63
1:A:192[B]:ASP:O	1:A:196:ARG:HG2	1.99	0.62
1:A:15:HIS:HE1	1:A:383:GLU:OE2	1.82	0.62
1:A:250:LEU:HD22	1:A:256:LEU:HG	1.81	0.62
1:C:258:ARG:NH1	1:C:262:ASP:OD2	2.32	0.62
1:B:53:ARG:CD	6:B:2072:HOH:O	2.48	0.61
1:A:78:MET:HE1	1:A:80:LYS:HZ1	1.67	0.60
1:B:250:LEU:CD2	1:B:256:LEU:HG	2.32	0.60
5:B:1400:GOL:H12	6:B:2160:HOH:O	2.02	0.60
1:A:347:LEU:C	1:A:347:LEU:HD12	2.27	0.60
1:B:284:THR:OG1	6:B:2009:HOH:O	2.14	0.60
1:C:15:HIS:HE1	1:C:383:GLU:OE2	1.85	0.59
1:B:22:ARG:HH11	1:B:22:ARG:HG2	1.68	0.57
1:C:78:MET:HE2	1:C:80:LYS:HZ2	1.54	0.56
1:C:180:GLN:O	1:C:184:GLU:HG3	2.04	0.56
1:B:345[A]:HIS:HE1	6:B:2281:HOH:O	1.88	0.55
1:A:353:ARG:HD2	6:A:2264:HOH:O	2.06	0.55

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:38:PRO:HD2	1:A:73:ARG:O	2.06	0.55
5:A:1399:GOL:H12	6:A:2162:HOH:O	2.07	0.54
1:B:53:ARG:HD3	6:B:2072:HOH:O	2.08	0.53
1:B:347:LEU:C	1:B:347:LEU:HD23	2.33	0.53
1:C:250:LEU:CD1	1:C:256:LEU:HG	2.33	0.53
1:B:113:ARG:HD2	1:B:116:GLU:OE1	2.10	0.52
1:A:53:ARG:CD	6:A:2079:HOH:O	2.57	0.51
4:C:1398:BEN:N2	6:C:2311:HOH:O	2.33	0.51
1:A:247:VAL:HG12	1:A:251:MET:HE2	1.93	0.50
1:A:318:GLN:NE2	1:A:324:ALA:H	2.10	0.49
1:C:250:LEU:HD22	1:C:360:LEU:HD22	1.94	0.49
1:C:9:TYR:CG	1:C:10:PRO:HA	2.47	0.49
1:C:257:ARG:NH2	1:C:370:ILE:O	2.39	0.49
1:A:113:ARG:HD2	1:A:116:GLU:OE1	2.13	0.49
1:A:318:GLN:HE22	1:A:324:ALA:H	1.59	0.49
1:C:250:LEU:CD2	1:C:360:LEU:HD22	2.43	0.49
1:A:296:LEU:O	1:A:299:VAL:HG13	2.12	0.49
1:B:9:TYR:O	1:B:38:PRO:HD3	2.13	0.49
1:B:100:LYS:HE3	1:B:213:ARG:O	2.12	0.49
2:B:450:HEM:HBC2	2:B:450:HEM:HMC2	1.95	0.48
1:B:9:TYR:CG	1:B:10:PRO:HA	2.49	0.47
1:B:318:GLN:NE2	1:B:324:ALA:H	2.13	0.46
2:A:450:HEM:HMC2	2:A:450:HEM:HBC2	1.97	0.46
1:C:95:ARG:O	1:C:99:VAL:HG22	2.15	0.46
1:B:192[B]:ASP:O	1:B:196:ARG:HG3	2.15	0.46
1:C:297:ARG:HD3	1:C:297:ARG:HA	1.80	0.45
1:C:346:CYS:HB2	2:C:450:HEM:NA	2.32	0.45
1:A:250:LEU:CD2	1:A:256:LEU:HG	2.43	0.45
1:A:15:HIS:CE1	1:A:383:GLU:OE2	2.68	0.45
1:C:196:ARG:HA	1:C:199:LYS:HE2	1.98	0.45
1:C:318:GLN:NE2	1:C:324:ALA:H	2.15	0.44
1:B:65:PRO:O	1:B:66:SER:C	2.59	0.44
1:C:113:ARG:HA	1:C:113:ARG:HD2	1.80	0.44
1:B:296:LEU:O	1:B:299:VAL:HG13	2.17	0.44
1:A:53:ARG:HD3	6:A:2079:HOH:O	2.17	0.44
1:A:346:CYS:HB2	2:A:450:HEM:NA	2.33	0.44
1:A:192[A]:ASP:O	1:A:196:ARG:HG2	2.17	0.44
1:A:174:VAL:O	3:A:460:MIV:H161	2.18	0.43
1:C:353:ARG:HD2	6:C:2280:HOH:O	2.18	0.43
1:B:9:TYR:CD1	1:B:10:PRO:HA	2.54	0.43
1:A:9:TYR:CD1	1:A:10:PRO:HA	2.54	0.43

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:22:ARG:HH11	1:B:22:ARG:CG	2.31	0.43
1:C:297:ARG:O	6:C:2227:HOH:O	2.22	0.43
1:C:318:GLN:HE22	1:C:324:ALA:H	1.66	0.43
1:C:71:GLU:OE1	1:C:78:MET:HE3	2.19	0.42
1:A:330[A]:ASP:O	1:A:331:ARG:C	2.62	0.42
1:B:246:PHE:CE1	1:B:274:LEU:HB3	2.55	0.42
1:B:318:GLN:HE22	1:B:324:ALA:H	1.66	0.42
1:A:235:GLY:HA2	2:A:450:HEM:C2C	2.55	0.42
1:B:250:LEU:HD23	1:B:256:LEU:HG	2.01	0.42
1:A:361:GLU:O	1:A:365:GLN:HB2	2.19	0.42
1:B:75:ARG:HD2	1:B:384:GLY:O	2.21	0.41
1:C:9:TYR:O	1:C:38:PRO:HD3	2.21	0.41
1:C:34:ARG:HA	1:C:34:ARG:HD2	1.94	0.41
1:C:215:GLN:HG3	6:C:2093:HOH:O	2.19	0.41
1:A:344:HIS:O	1:A:345[B]:HIS:C	2.63	0.41
1:A:9:TYR:CG	1:A:10:PRO:HA	2.55	0.41
1:A:137:MET:HE2	1:A:137:MET:HB3	1.98	0.41
1:B:170:SER:HA	3:B:460:MIV:H212	2.03	0.41
1:A:323:ASP:HA	1:C:88:PRO:HB2	2.03	0.41
1:C:347:LEU:C	1:C:347:LEU:CD1	2.93	0.41
1:C:9:TYR:CD1	1:C:10:PRO:HA	2.55	0.40
1:A:258:ARG:NH1	1:A:262:ASP:OD2	2.54	0.40
3:B:460:MIV:C24	3:B:460:MIV:C1	2.99	0.40
1:C:199:LYS:H	1:C:199:LYS:HG2	1.73	0.40
1:A:346:CYS:HA	2:A:450:HEM:CHA	2.51	0.40
1:B:113:ARG:HD2	1:B:113:ARG:HA	1.93	0.40
1:C:185[A]:GLN:NE2	6:C:2171:HOH:O	2.53	0.40
1:C:215:GLN:O	1:C:216:GLN:HB2	2.22	0.40
1:A:53:ARG:NE	6:A:2079:HOH:O	2.53	0.40
1:A:75:ARG:HD2	1:A:384:GLY:O	2.22	0.40

All (1) 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)
6:B:2041:HOH:O	6:B:2228:HOH:O[4_555]	2.19	0.01

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	398/417 (95%)	385 (97%)	13 (3%)	0	100	100
1	B	398/417 (95%)	388 (98%)	10 (2%)	0	100	100
1	C	397/417 (95%)	386 (97%)	11 (3%)	0	100	100
All	All	1193/1251 (95%)	1159 (97%)	34 (3%)	0	100	100

There are no Ramachandran outliers to report.

5.3.2 Protein sidechains (i)

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

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

Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
1	A	327/346 (94%)	314 (96%)	13 (4%)	28	11
1	B	329/346 (95%)	322 (98%)	7 (2%)	47	30
1	C	329/346 (95%)	319 (97%)	10 (3%)	36	19
All	All	985/1038 (95%)	955 (97%)	30 (3%)	36	19

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	25	GLU
1	A	80	LYS
1	A	96	ARG
1	A	100	LYS
1	A	173	GLU

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Mol	Chain	Res	Type
1	A	180	GLN
1	A	211	GLN
1	A	250	LEU
1	A	284	THR
1	A	299	VAL
1	A	330[A]	ASP
1	A	330[B]	ASP
1	A	365	GLN
1	B	149	GLU
1	B	180	GLN
1	B	184	GLU
1	B	199	LYS
1	B	250	LEU
1	B	297	ARG
1	B	299	VAL
1	C	80	LYS
1	C	100	LYS
1	C	173	GLU
1	C	180	GLN
1	C	211	GLN
1	C	215	GLN
1	C	250	LEU
1	C	299	VAL
1	C	347	LEU
1	C	392	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (18) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	HIS
1	A	124	GLN
1	A	130	GLN
1	A	180	GLN
1	A	211	GLN
1	A	259	GLN
1	A	318	GLN
1	B	15	HIS
1	B	124	GLN
1	B	130	GLN
1	B	259	GLN
1	B	318	GLN
1	C	15	HIS

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Mol	Chain	Res	Type
1	C	211	GLN
1	C	215	GLN
1	C	318	GLN
1	C	335	GLN
1	C	365	GLN

5.3.3 RNA

There are no RNA molecules in this entry.

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

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

5.5 Carbohydrates

There are no oligosaccharides in this entry.

5.6 Ligand geometry

15 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	$\#\|Z\| > 2$
2	HEM	C	450	1,6	50,50,50	1.87	12 (24%)	66,82,82	1.42	7 (10%)
3	MIV	A	460	-	51,51,51	1.11	2 (3%)	62,71,71	1.45	10 (16%)
4	BEN	C	1398	-	9,9,9	0.63	0	7,11,11	0.67	0
4	BEN	B	1398	4	9,9,9	0.71	0	7,11,11	1.01	0
4	BEN	A	1398	4	9,9,9	0.70	0	7,11,11	1.00	0
2	HEM	B	450	1,6	50,50,50	1.89	9 (18%)	66,82,82	1.35	7 (10%)
5	GOL	A	1399	-	5,5,5	0.46	0	5,5,5	2.43	1 (20%)
3	MIV	C	460	-	51,51,51	1.06	3 (5%)	62,71,71	1.59	12 (19%)

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	$\#\|Z\|>2$	Counts	RMSZ	$\#\|Z\|>2$
5	GOL	A	1400	-	5,5,5	1.15	0	5,5,5	1.74	1 (20%)
3	MIV	B	460	-	51,51,51	1.01	2 (3%)	62,71,71	1.56	12 (19%)
5	GOL	C	1400	-	5,5,5	0.41	0	5,5,5	1.13	1 (20%)
5	GOL	B	1399	-	5,5,5	1.03	0	5,5,5	1.78	1 (20%)
5	GOL	B	1400	-	5,5,5	0.46	0	5,5,5	1.65	1 (20%)
5	GOL	C	1399	-	5,5,5	0.92	0	5,5,5	1.49	1 (20%)
2	HEM	A	450	1,6	50,50,50	1.83	9 (18%)	66,82,82	1.25	6 (9%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	C	450	1,6	-	0/14/54/54	-
3	MIV	A	460	-	-	10/55/91/91	0/2/3/3
4	BEN	C	1398	-	-	0/4/4/4	0/1/1/1
4	BEN	B	1398	4	-	4/4/4/4	0/1/1/1
4	BEN	A	1398	4	-	4/4/4/4	0/1/1/1
2	HEM	B	450	1,6	-	0/14/54/54	-
5	GOL	A	1399	-	-	4/4/4/4	-
3	MIV	C	460	-	-	9/55/91/91	0/2/3/3
5	GOL	A	1400	-	-	3/4/4/4	-
3	MIV	B	460	-	-	7/55/91/91	0/2/3/3
5	GOL	C	1400	-	-	2/4/4/4	-
5	GOL	B	1399	-	-	3/4/4/4	-
5	GOL	B	1400	-	-	2/4/4/4	-
5	GOL	C	1399	-	-	3/4/4/4	-
2	HEM	A	450	1,6	-	0/14/54/54	-

All (37) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	450	HEM	C3D-C2D	7.06	1.51	1.36
2	A	450	HEM	C3D-C2D	6.41	1.50	1.36
2	C	450	HEM	C3D-C2D	6.22	1.50	1.36
3	A	460	MIV	O6-C24	5.22	1.45	1.34
2	B	450	HEM	FE-NB	5.20	2.11	1.94

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	C	450	HEM	FE-NB	4.96	2.10	1.94
2	A	450	HEM	FE-ND	4.92	2.10	1.94
2	C	450	HEM	FE-ND	4.78	2.09	1.94
2	B	450	HEM	FE-ND	4.61	2.09	1.94
2	A	450	HEM	FE-NB	4.40	2.08	1.94
3	B	460	MIV	O6-C24	4.34	1.43	1.34
3	C	460	MIV	O6-C24	4.27	1.43	1.34
2	B	450	HEM	CAC-C3C	3.19	1.56	1.47
2	C	450	HEM	CAC-C3C	3.16	1.56	1.47
2	A	450	HEM	CAB-C3B	3.02	1.55	1.47
2	C	450	HEM	CMA-C3A	2.93	1.57	1.50
2	A	450	HEM	CMC-C2C	2.86	1.56	1.50
2	A	450	HEM	CAC-C3C	2.85	1.55	1.47
2	A	450	HEM	CMA-C3A	2.85	1.56	1.50
2	B	450	HEM	CMA-C3A	2.83	1.56	1.50
2	C	450	HEM	CAB-C3B	2.74	1.54	1.47
3	C	460	MIV	O9-C33	2.73	1.49	1.42
2	C	450	HEM	FE-NC	2.53	2.03	1.95
2	B	450	HEM	CMC-C2C	2.42	1.55	1.50
2	C	450	HEM	CMC-C2C	2.42	1.55	1.50
2	B	450	HEM	CAB-C3B	2.41	1.54	1.47
2	C	450	HEM	C1B-NB	-2.32	1.36	1.40
2	A	450	HEM	CMB-C2B	2.29	1.55	1.50
2	C	450	HEM	CMB-C2B	2.21	1.55	1.50
2	B	450	HEM	FE-NA	2.20	2.02	1.95
2	A	450	HEM	C4D-ND	-2.16	1.36	1.40
3	B	460	MIV	C29-C28	2.13	1.54	1.52
2	B	450	HEM	CMB-C2B	2.09	1.55	1.50
3	A	460	MIV	O9-C33	2.07	1.47	1.42
2	C	450	HEM	O1D-CGD	2.06	1.29	1.22
2	C	450	HEM	FE-NA	2.04	2.02	1.95
3	C	460	MIV	C29-C28	2.00	1.54	1.52

All (60) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	C	450	HEM	C4D-ND-C1D	5.93	111.20	105.07
2	B	450	HEM	C4D-ND-C1D	5.50	110.75	105.07
5	A	1399	GOL	C3-C2-C1	-5.30	91.10	111.70
2	A	450	HEM	C4D-ND-C1D	4.87	110.11	105.07
3	C	460	MIV	O9-C33-C31	4.27	118.98	108.94
3	B	460	MIV	C20-C22-C23	-3.85	115.94	126.44

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
3	B	460	MIV	C2-C3-C4	-3.75	115.32	124.67
3	A	460	MIV	C34-O9-C33	3.63	124.04	114.52
5	B	1400	GOL	C3-C2-C1	-3.62	97.61	111.70
3	C	460	MIV	C20-C22-C23	-3.58	116.67	126.44
5	B	1399	GOL	O2-C2-C3	3.52	124.64	109.12
2	B	450	HEM	C1B-NB-C4B	3.48	108.67	105.07
3	B	460	MIV	O6-C24-O5	-3.41	117.79	123.35
3	C	460	MIV	O4-C18-C17	3.39	114.34	109.14
3	C	460	MIV	O6-C24-O5	-3.38	117.85	123.35
3	B	460	MIV	C35-C33-C31	-3.33	104.05	111.66
5	A	1400	GOL	O2-C2-C1	3.27	123.53	109.12
3	A	460	MIV	O9-C33-C35	3.25	118.06	110.29
3	C	460	MIV	C34-O9-C33	3.17	122.84	114.52
2	B	450	HEM	CHD-C4C-NC	3.15	127.85	124.44
3	C	460	MIV	O9-C33-C35	3.03	117.53	110.29
2	C	450	HEM	CHD-C4C-NC	3.01	127.69	124.44
2	A	450	HEM	C1D-C2D-C3D	-2.98	103.82	106.96
5	C	1399	GOL	O2-C2-C1	2.92	121.98	109.12
3	B	460	MIV	O8-C31-C33	2.90	115.77	108.94
3	B	460	MIV	C21-C20-C22	-2.90	102.99	109.99
3	C	460	MIV	C2-C3-C4	-2.88	117.49	124.67
3	B	460	MIV	O4-C18-C17	2.86	113.52	109.14
2	C	450	HEM	CAC-C3C-C4C	-2.81	118.10	124.90
2	B	450	HEM	C1D-C2D-C3D	-2.78	104.03	106.96
3	A	460	MIV	C21-C20-C22	-2.78	103.27	109.99
2	C	450	HEM	C1B-NB-C4B	2.70	107.86	105.07
3	A	460	MIV	O8-C31-C33	2.66	115.20	108.94
2	A	450	HEM	CAC-C3C-C4C	-2.65	118.50	124.90
2	A	450	HEM	CHD-C4C-NC	2.63	127.28	124.44
3	C	460	MIV	C28-C1-C2	-2.63	118.82	125.45
3	C	460	MIV	O8-C31-C33	2.62	115.11	108.94
3	B	460	MIV	O6-C24-C23	2.52	117.09	111.38
2	A	450	HEM	C4B-C3B-C2B	2.48	109.09	107.11
2	C	450	HEM	C4C-NC-C1C	2.42	107.72	105.35
3	A	460	MIV	O6-C24-O5	-2.41	119.42	123.35
3	A	460	MIV	C15-N-C16	2.40	117.45	110.38
2	B	450	HEM	CHC-C1C-NC	2.38	127.01	124.44
3	A	460	MIV	C20-C22-C23	-2.37	119.97	126.44
3	B	460	MIV	C13-C14-N	-2.33	104.09	110.83
3	A	460	MIV	C2-C3-C4	-2.29	118.95	124.67
3	A	460	MIV	C28-C1-C2	-2.29	119.67	125.45
3	C	460	MIV	C35-C33-C31	-2.26	106.49	111.66

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	450	HEM	C4B-C3B-C2B	2.26	108.91	107.11
3	B	460	MIV	C28-C1-C2	-2.25	119.77	125.45
2	C	450	HEM	CHC-C4B-NB	2.25	126.86	124.42
3	B	460	MIV	O3-C13-C12	2.22	115.45	110.05
3	A	460	MIV	O4-C18-C17	2.22	112.53	109.14
2	B	450	HEM	CMD-C2D-C3D	2.10	131.81	126.12
3	C	460	MIV	O3-C13-C14	2.07	113.47	109.77
5	C	1400	GOL	C3-C2-C1	-2.04	103.79	111.70
2	A	450	HEM	C1B-NB-C4B	2.01	107.15	105.07
3	B	460	MIV	O9-C33-C31	2.01	113.67	108.94
2	C	450	HEM	CHB-C4A-NA	2.01	127.54	123.85
3	C	460	MIV	O6-C24-C23	2.00	115.92	111.38

There are no chirality outliers.

All (51) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	460	MIV	O6-C25-C26-C27
3	B	460	MIV	O6-C25-C26-C27
3	C	460	MIV	O6-C25-C26-C27
5	A	1399	GOL	C1-C2-C3-O3
5	A	1399	GOL	O2-C2-C3-O3
5	B	1399	GOL	O1-C1-C2-C3
5	B	1400	GOL	C1-C2-C3-O3
5	B	1400	GOL	O2-C2-C3-O3
3	A	460	MIV	C35-C33-O9-C34
3	C	460	MIV	C35-C33-O9-C34
3	A	460	MIV	C30-C31-O8-C32
5	C	1399	GOL	O1-C1-C2-O2
3	C	460	MIV	O11-C30-O7-C29
3	A	460	MIV	C31-C30-O7-C29
3	B	460	MIV	C31-C30-O7-C29
3	C	460	MIV	C31-C30-O7-C29
5	A	1400	GOL	O1-C1-C2-C3
5	B	1399	GOL	C1-C2-C3-O3
5	C	1399	GOL	O1-C1-C2-C3
5	C	1400	GOL	C1-C2-C3-O3
3	A	460	MIV	O11-C30-O7-C29
3	B	460	MIV	O11-C30-O7-C29
5	A	1400	GOL	O1-C1-C2-O2
5	B	1399	GOL	O2-C2-C3-O3
3	A	460	MIV	C1-C2-C3-C4

Continued on next page...

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Mol	Chain	Res	Type	Atoms
3	B	460	MIV	C1-C2-C3-C4
3	C	460	MIV	C1-C2-C3-C4
3	A	460	MIV	C33-C31-O8-C32
3	B	460	MIV	C17-C14-N-C15
5	A	1399	GOL	O1-C1-C2-O2
5	A	1400	GOL	C1-C2-C3-O3
5	C	1399	GOL	C1-C2-C3-O3
3	A	460	MIV	C28-C25-C26-C27
3	B	460	MIV	C28-C25-C26-C27
3	B	460	MIV	C28-C29-O7-C30
3	C	460	MIV	C28-C29-O7-C30
5	C	1400	GOL	O2-C2-C3-O3
4	A	1398	BEN	N2-C-C1-C2
4	A	1398	BEN	N2-C-C1-C6
4	B	1398	BEN	N2-C-C1-C2
4	B	1398	BEN	N2-C-C1-C6
3	C	460	MIV	C33-C31-O8-C32
5	A	1399	GOL	O1-C1-C2-C3
3	C	460	MIV	C30-C31-O8-C32
4	A	1398	BEN	N1-C-C1-C2
4	A	1398	BEN	N1-C-C1-C6
4	B	1398	BEN	N1-C-C1-C2
4	B	1398	BEN	N1-C-C1-C6
3	C	460	MIV	C28-C25-C26-C27
3	A	460	MIV	C22-C23-C24-O6
3	A	460	MIV	C28-C29-O7-C30

There are no ring outliers.

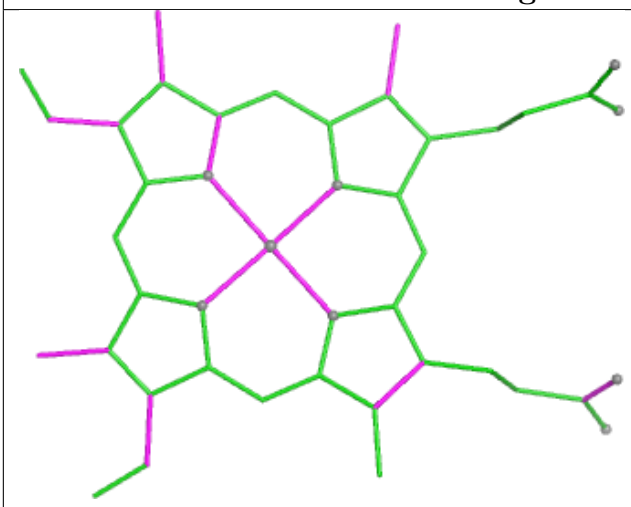
8 monomers are involved in 12 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	C	450	HEM	1	0
3	A	460	MIV	1	0
4	C	1398	BEN	1	0
2	B	450	HEM	1	0
5	A	1399	GOL	1	0
3	B	460	MIV	2	0
5	B	1400	GOL	1	0
2	A	450	HEM	4	0

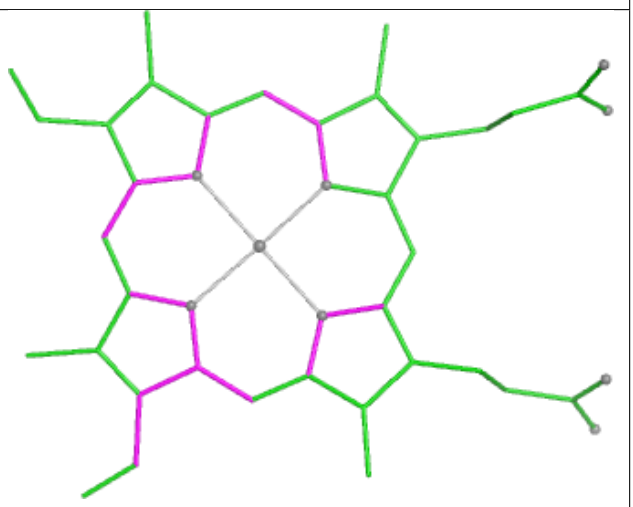
The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In

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.

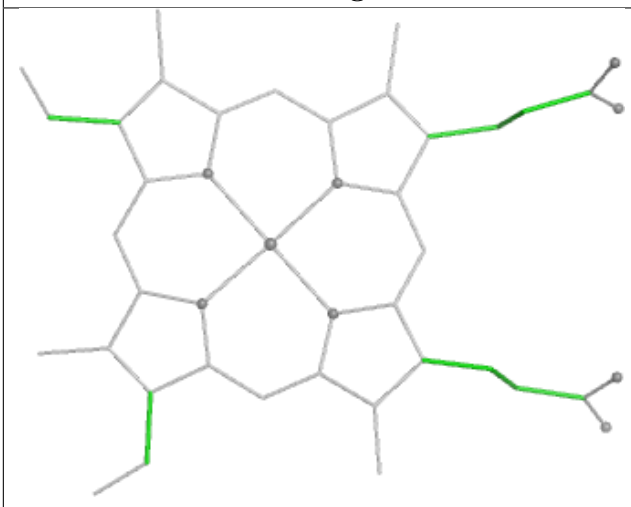
Ligand HEM C 450



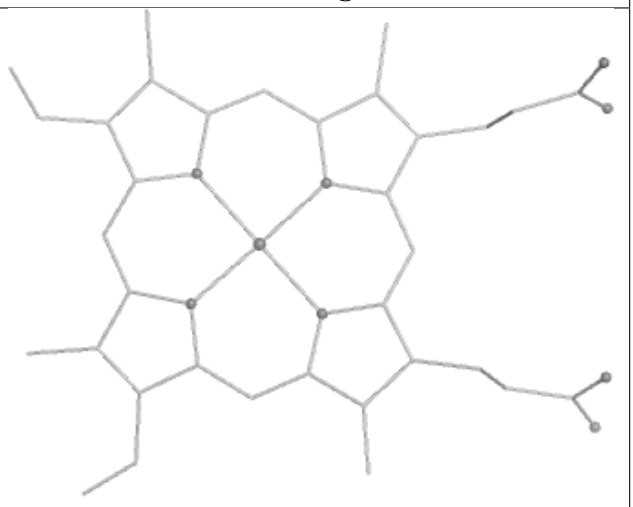
Bond lengths



Bond angles

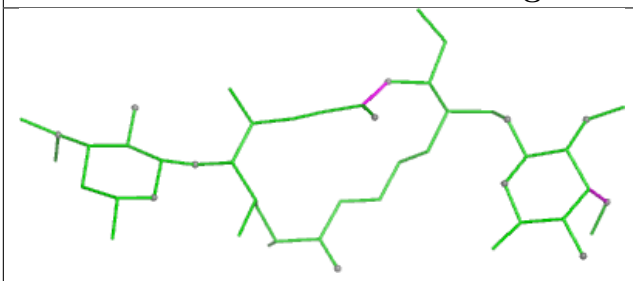


Torsions

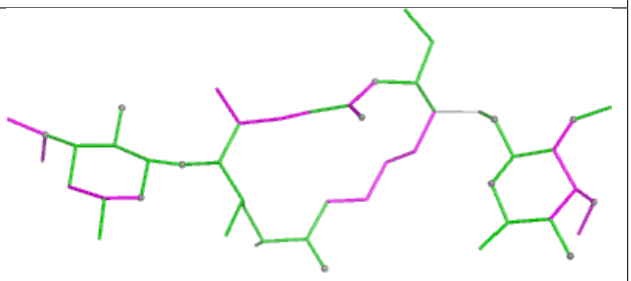


Rings

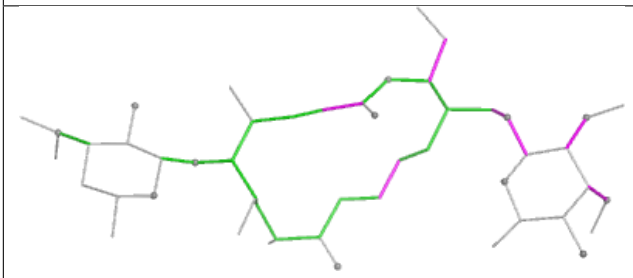
Ligand MIV A 460



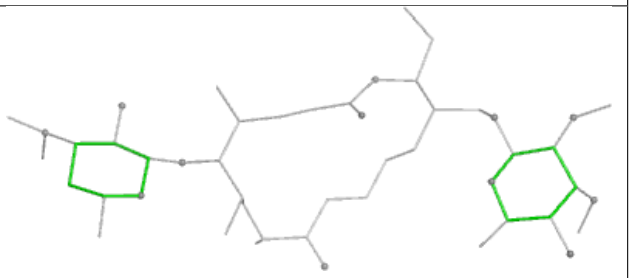
Bond lengths



Bond angles

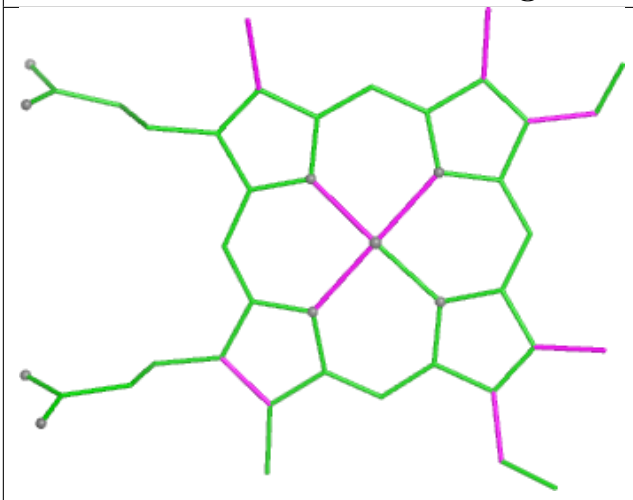


Torsions

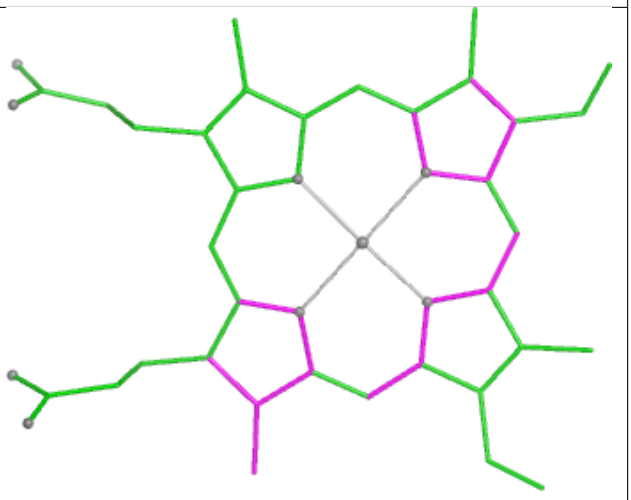


Rings

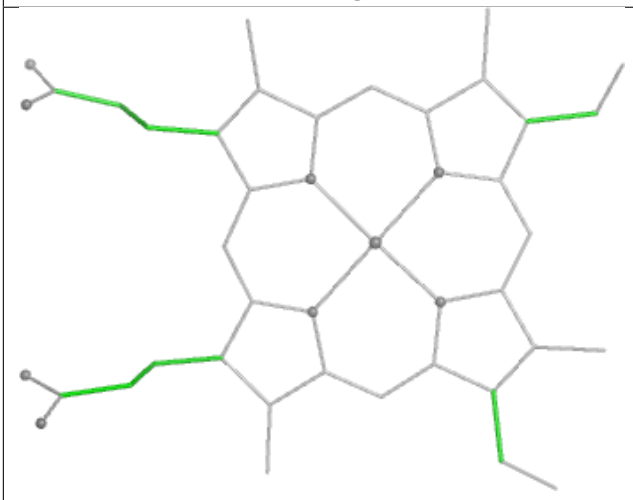
Ligand HEM B 450



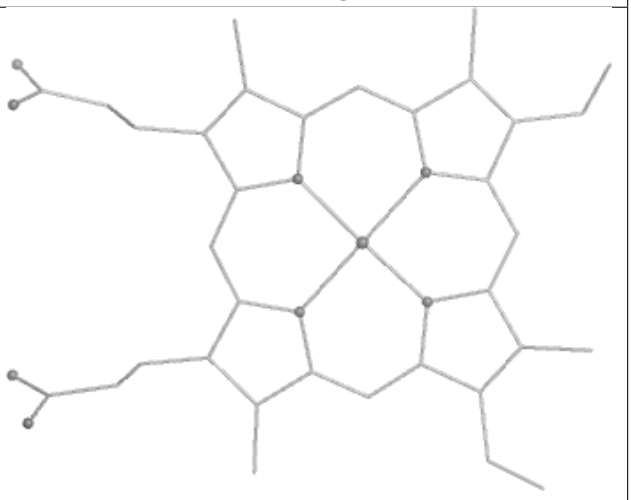
Bond lengths



Bond angles

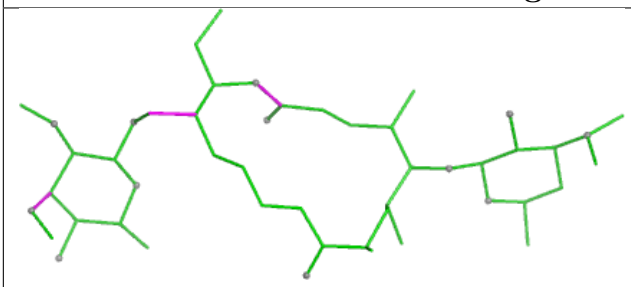


Torsions

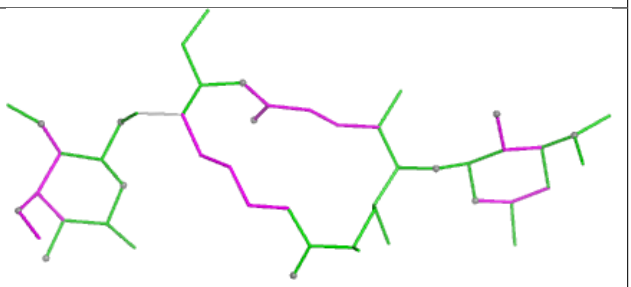


Rings

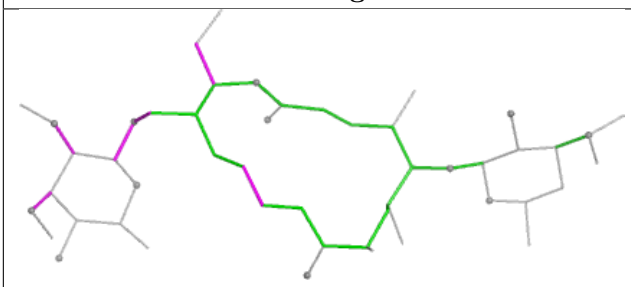
Ligand MIV C 460



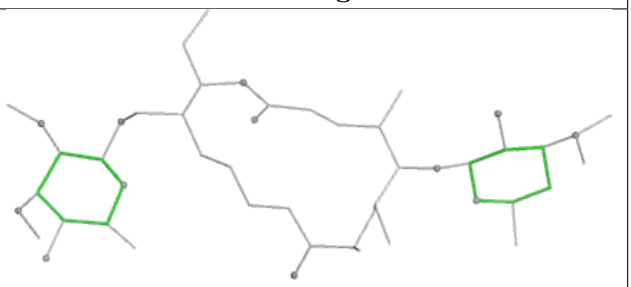
Bond lengths



Bond angles

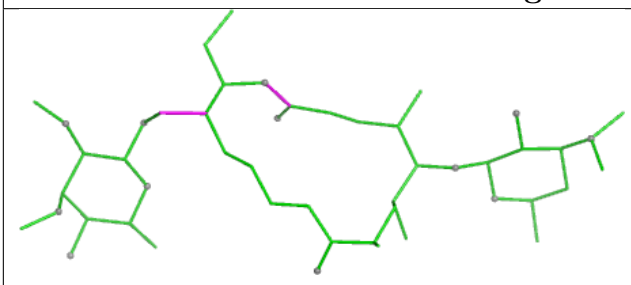


Torsions

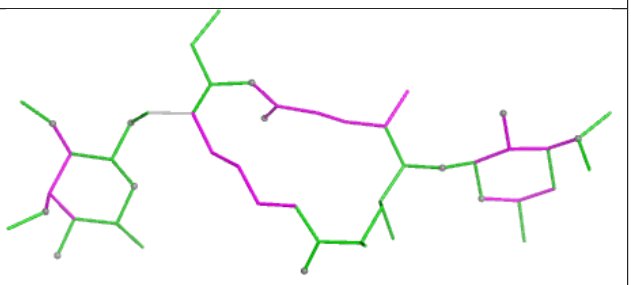


Rings

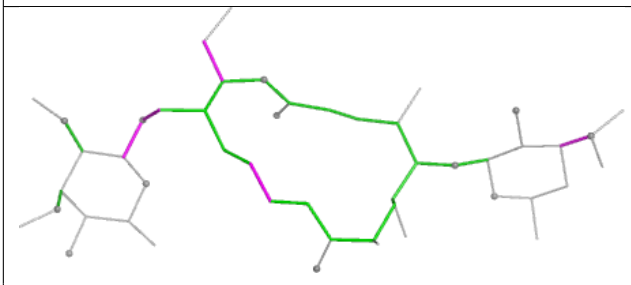
Ligand MIV B 460



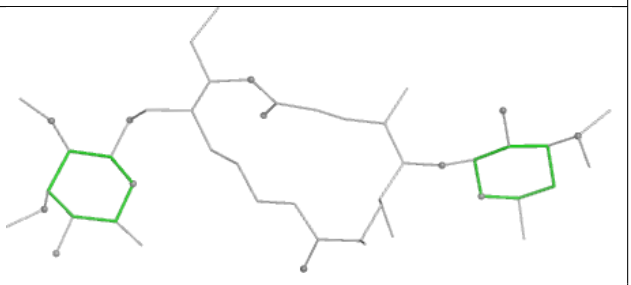
Bond lengths



Bond angles

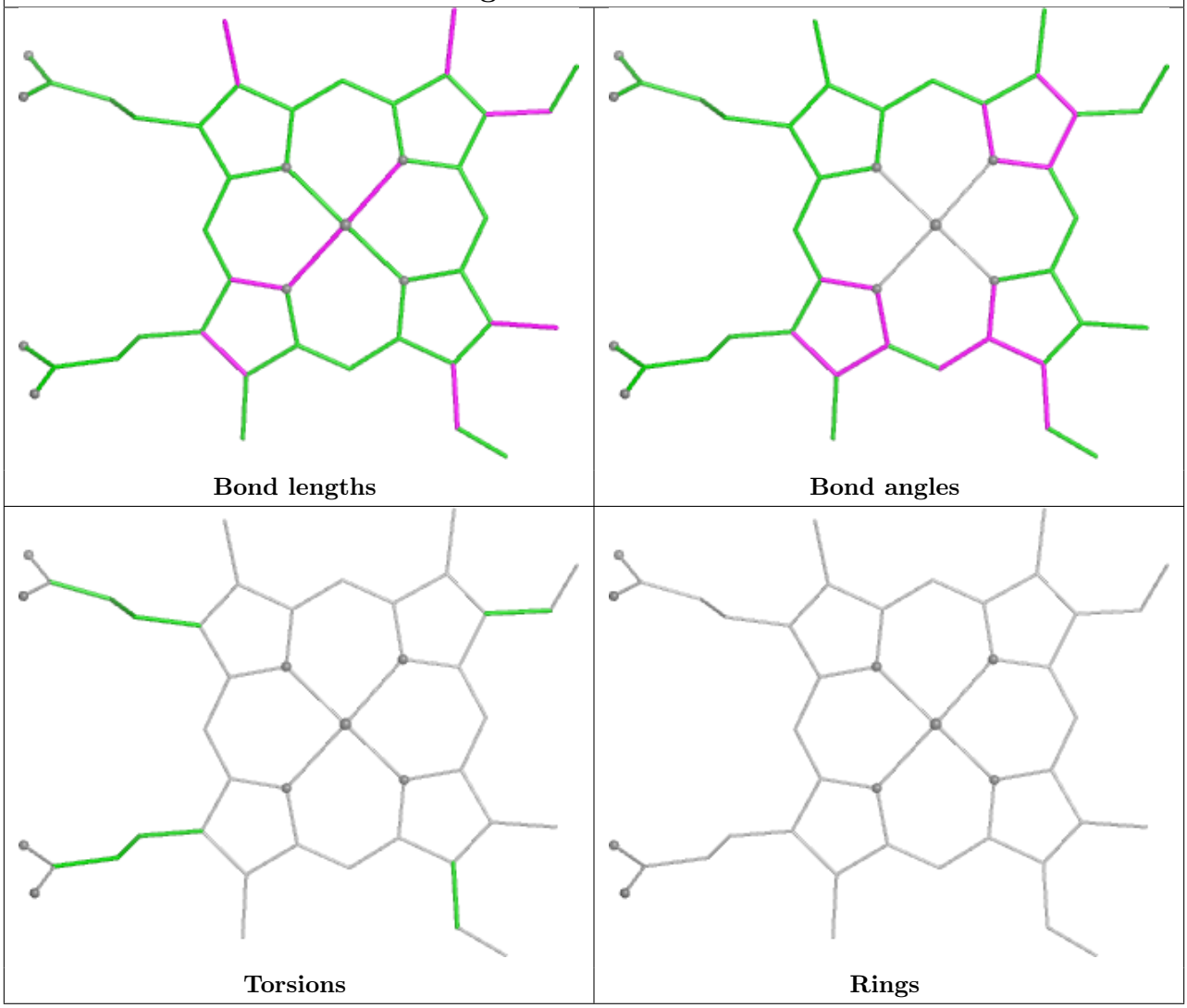


Torsions



Rings

Ligand HEM A 450



5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data

6.1 Protein, DNA and RNA chains

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	393/417 (94%)	-1.50	0	100	100	6, 20, 36, 44	7 (1%)
1	B	393/417 (94%)	-1.49	0	100	100	6, 20, 37, 44	7 (1%)
1	C	393/417 (94%)	-1.51	0	100	100	6, 20, 36, 49	6 (1%)
All	All	1179/1251 (94%)	-1.50	0	100	100	6, 20, 37, 49	20 (1%)

There are no RSRZ outliers to report.

6.2 Non-standard residues in protein, DNA, RNA chains

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

6.3 Carbohydrates

There are no oligosaccharides in this entry.

6.4 Ligands

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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	MIV	A	460	49/49	0.99	0.03	26,32,43,44	0
3	MIV	B	460	49/49	0.99	0.03	25,30,42,44	0
3	MIV	C	460	49/49	0.99	0.03	26,30,42,43	0
4	BEN	A	1398	9/9	0.99	0.03	37,37,37,39	9
4	BEN	B	1398	9/9	0.99	0.03	37,37,37,39	9

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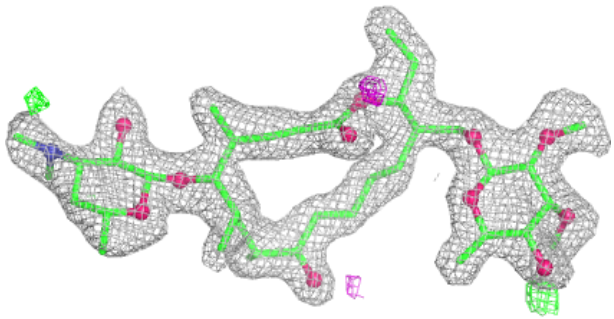
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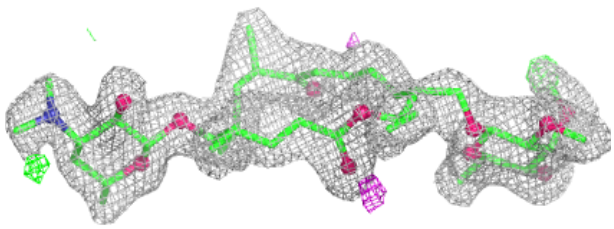
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	GOL	A	1399	6/6	0.99	0.02	24,25,26,26	0
5	GOL	A	1400	6/6	0.99	0.03	14,19,20,21	0
5	GOL	B	1399	6/6	0.99	0.04	13,21,24,24	0
5	GOL	C	1399	6/6	0.99	0.05	14,22,23,23	0
5	GOL	C	1400	6/6	0.99	0.03	22,24,24,26	0
2	HEM	A	450	43/43	1.00	0.02	9,11,14,16	0
2	HEM	B	450	43/43	1.00	0.02	9,11,14,16	0
5	GOL	B	1400	6/6	1.00	0.02	25,28,28,28	0
4	BEN	C	1398	9/9	1.00	0.02	27,28,29,30	9
2	HEM	C	450	43/43	1.00	0.02	9,10,14,17	0

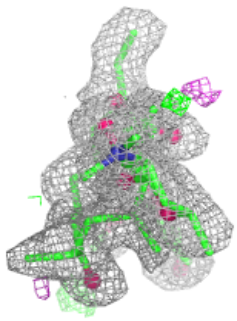
The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation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

Electron density around MIV A 460:

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

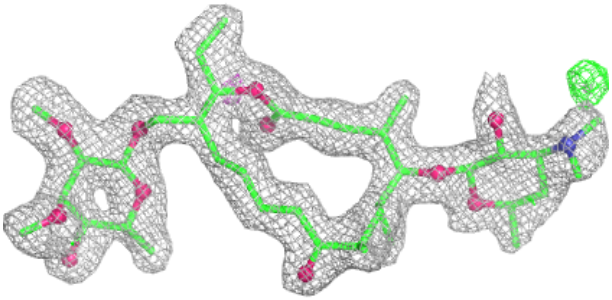


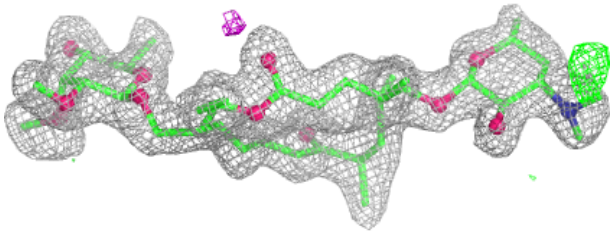


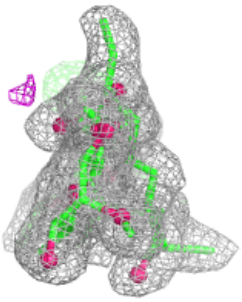


Electron density around MIV B 460:

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

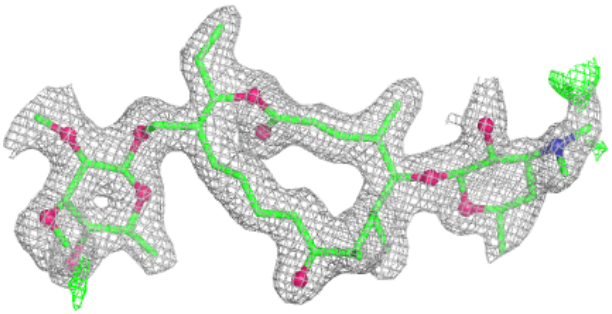


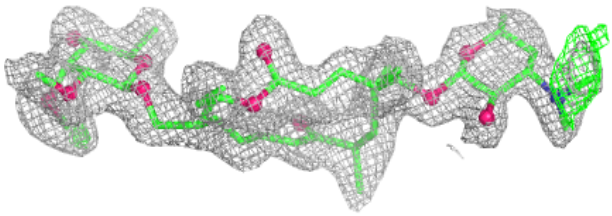


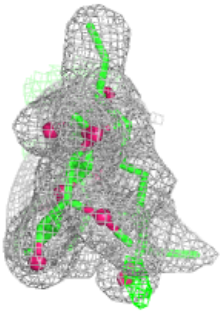


Electron density around MIV C 460:

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

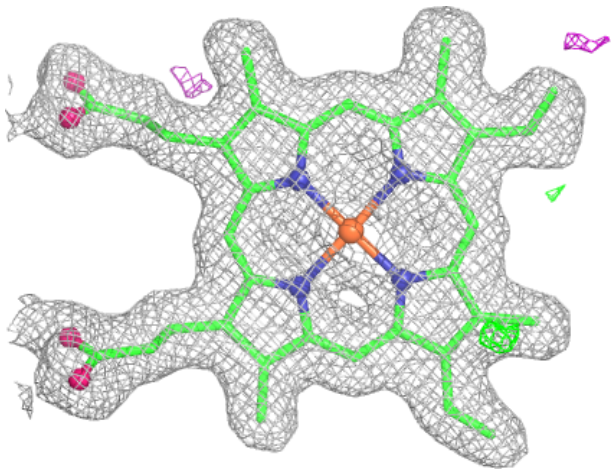


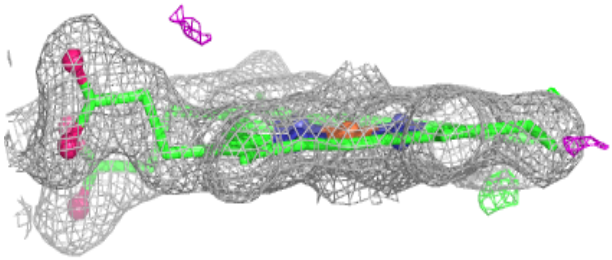


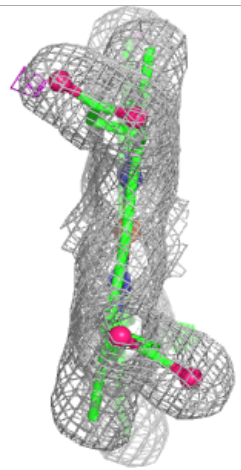


Electron density around HEM A 450:

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

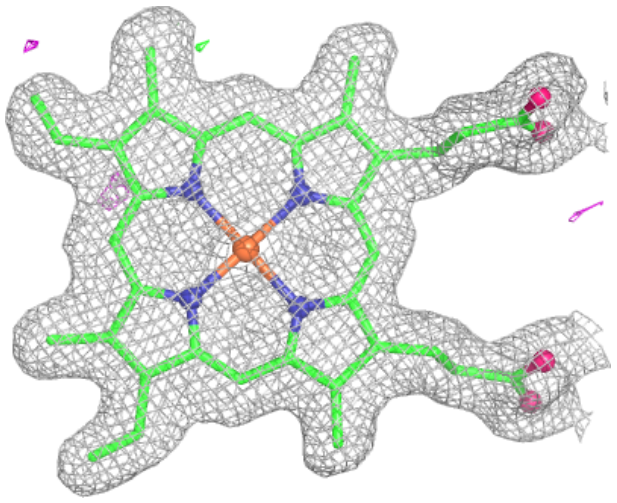


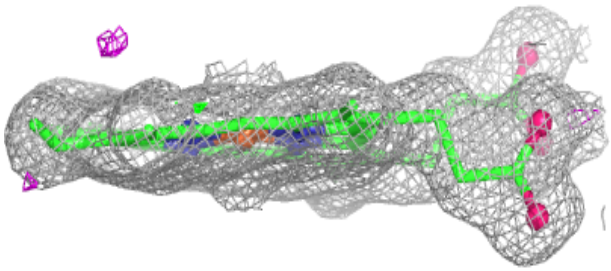


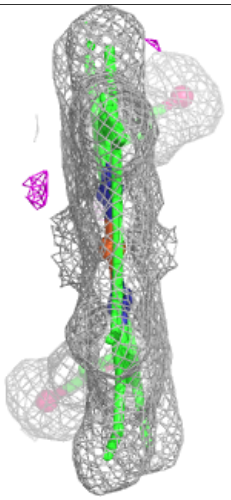


Electron density around HEM B 450:

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

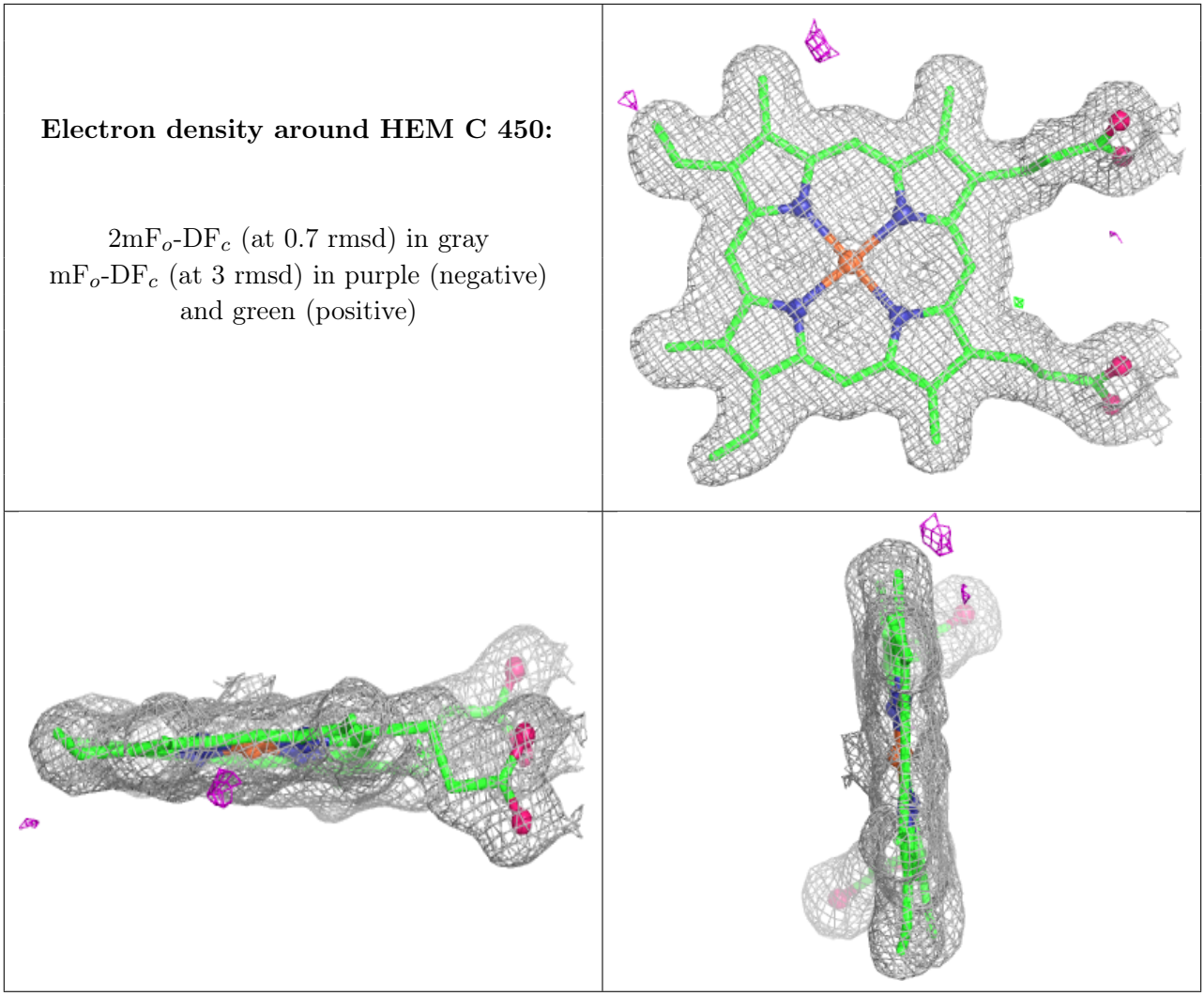






Electron density around HEM C 450:

$2mF_o$-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

There are no such residues in this entry.