Full wwPDB X-ray Structure Validation Report (i)

Aug 5, 2026 – 03:24 PM JST

PDB ID	:	6JO1 / pdb_00006jo1
Title	:	Structure of the CYP102A1 Haem Domain with N-(S)-Ibuprofenoyl-L-Phenylalanine
Authors	:	Stanfield, J.K.; Kasai, C.; Sugimoto, H.; Shiro, Y.; Watanabe, Y.; Shoji, O.
Deposited on	:	2019-03-19
Resolution	:	2.10 Å(reported)

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

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

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

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

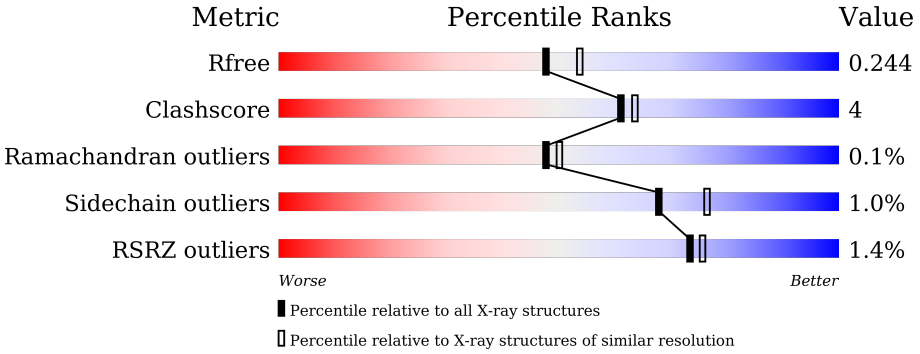
MolProbity	:	4-5-2 with Phenix2.0
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Buster-report	:	wwPDB partial adaption of 1.1.7 (2018)
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.015 (Gargrove)
Density-Fitness	:	1.0.12
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 2.10 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	6658 (2.10-2.10)
Clashscore	190562	7164 (2.10-2.10)
Ramachandran outliers	187476	7099 (2.10-2.10)
Sidechain outliers	187428	7100 (2.10-2.10)
RSRZ outliers	180081	6662 (2.10-2.10)

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	456	.% 91% 8% •
1	B	456	2% 88% 11% ••

2 Entry composition (i)

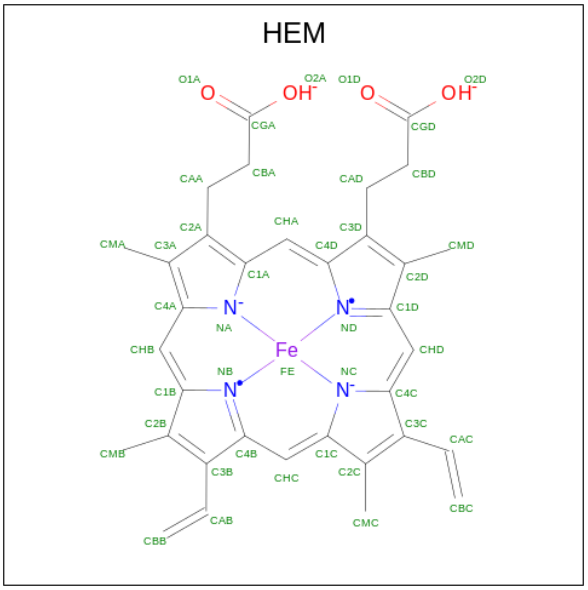
There are 6 unique types of molecules in this entry. The entry contains 7739 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Bifunctional cytochrome P450/NADPH–P450 reductase.

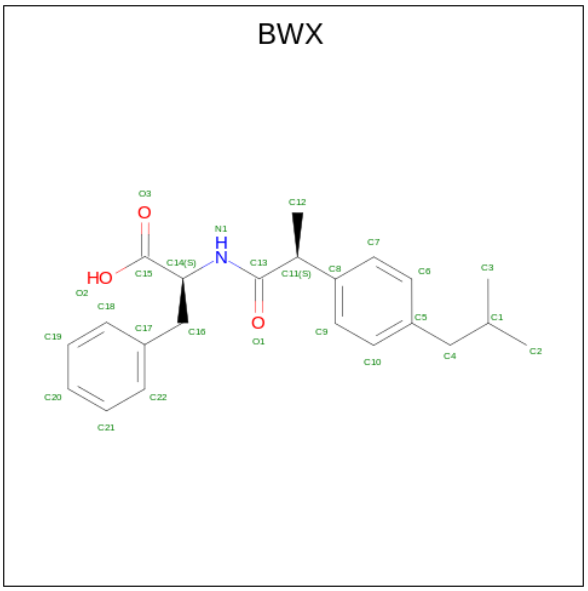
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	453	3682	2350	625	689	18	0	3	0
1	B	453	3690	2354	626	692	18	0	4	0

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



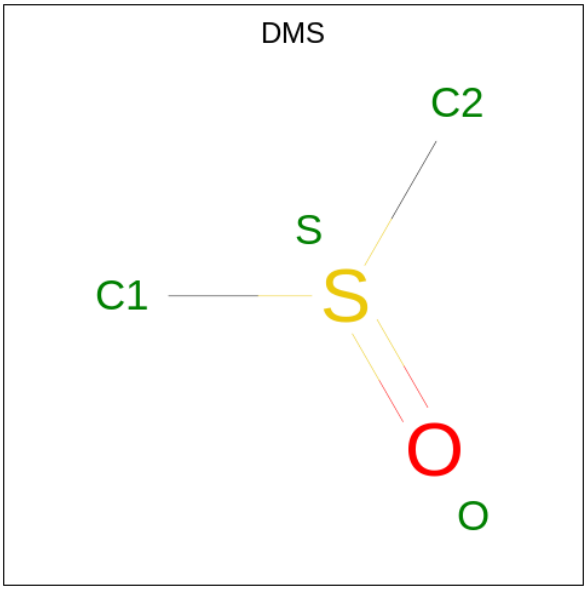
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			Total	C	Fe	N	O		
2	A	1	43	34	1	4	4	0	0
2	B	1	43	34	1	4	4	0	0

- Molecule 3 is (2S)-2-[[(2S)-2-[4-(2-methylpropyl)phenyl]propanoyl]amino]-3-phenyl-propanoic acid (CCD ID: BWX) (formula: $C_{22}H_{27}NO_3$) (labeled as "Ligand of Interest" by depositor).



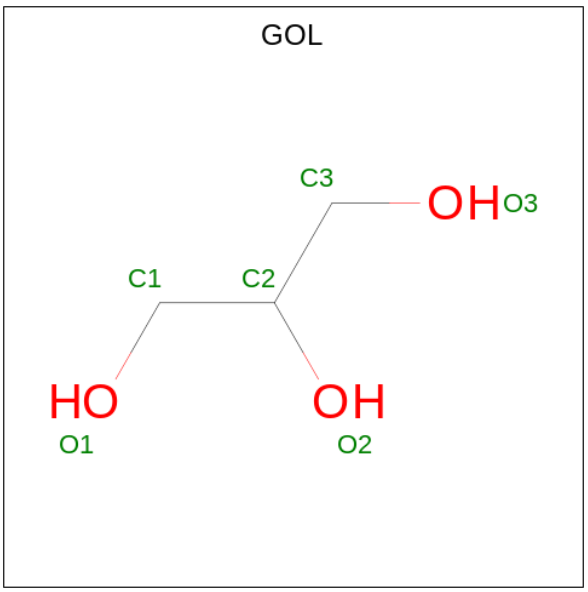
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	26	22	1	3	0	0
			Total	C	N	O		
3	B	1	26	22	1	3	0	0

- Molecule 4 is DIMETHYL SULFOXIDE (CCD ID: DMS) (formula: C_2H_6OS).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	O	S		
4	A	1	4	2	1	1	0	0
			Total	C	O	S		
4	B	1	4	2	1	1	0	0

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



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

- Molecule 6 is water.

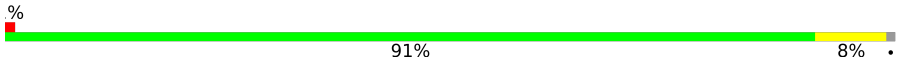
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
6	A	96	96	96	0	0
			Total	O		
6	B	107	107	107	0	0

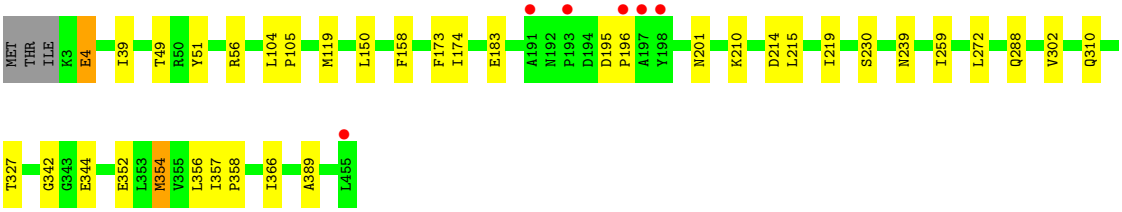
3 Residue-property plots (i)

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: Bifunctional cytochrome P450/NADPH–P450 reductase

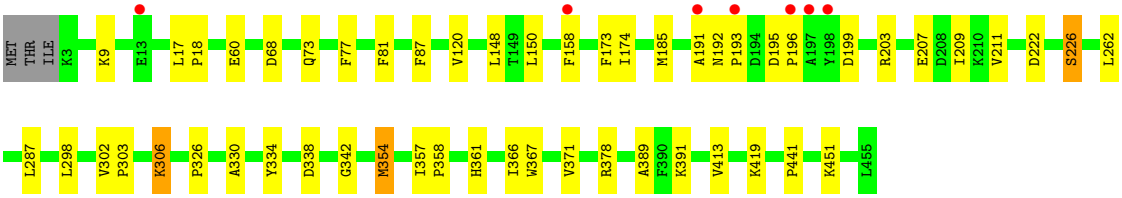
Chain A:





• Molecule 1: Bifunctional cytochrome P450/NADPH–P450 reductase

Chain B:



4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α, β, γ	58.90Å 148.81Å 64.69Å 90.00° 98.12° 90.00°	Depositor
Resolution (Å)	46.55 – 2.10 46.55 – 2.10	Depositor EDS
% Data completeness (in resolution range)	99.9 (46.55-2.10) 99.9 (46.55-2.10)	Depositor EDS
R_{merge}	0.14	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.52 (at 2.10Å)	Xtriage
Refinement program	REFMAC 5.8.0238	Depositor
R, R_{free}	0.192 , 0.241 0.199 , 0.244	Depositor DCC
R_{free} test set	3142 reflections (4.90%)	wwPDB-VP
Wilson B-factor (Å^2)	35.8	Xtriage
Anisotropy	0.122	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.32 , 23.1	EDS
L-test for twinning[2]	$< \|L\| > = 0.50$, $< L^2 > = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.96	EDS
Total number of atoms	7739	wwPDB-VP
Average B, all atoms (Å^2)	43.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.80% 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: GOL, DMS, HEM, BWX

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	1.04	2/3767 (0.1%)	1.37	6/5090 (0.1%)
1	B	1.03	1/3775 (0.0%)	1.35	1/5101 (0.0%)
All	All	1.03	3/7542 (0.0%)	1.36	7/10191 (0.1%)

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	441	PRO	C-O	-5.53	1.19	1.24
1	A	259	ILE	C-O	5.39	1.30	1.24
1	A	302	VAL	N-CA	5.35	1.50	1.46

All (7) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	239	ASN	CA-C-N	6.29	125.96	119.92
1	A	239	ASN	C-N-CA	6.29	125.96	119.92
1	A	310	GLN	CB-CA-C	5.79	120.34	110.56
1	A	352	GLU	CB-CA-C	-5.30	100.78	109.53
1	B	338	ASP	CA-CB-CG	5.27	117.87	112.60
1	A	119	MET	CA-C-N	5.13	127.03	120.56
1	A	119	MET	C-N-CA	5.13	127.03	120.56

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	3682	0	3644	21	0
1	B	3690	0	3647	37	0
2	A	43	0	30	2	0
2	B	43	0	30	3	0
3	A	26	0	0	0	0
3	B	26	0	0	0	0
4	A	4	0	6	0	0
4	B	4	0	6	1	0
5	A	6	0	8	0	0
5	B	12	0	16	2	0
6	A	96	0	0	0	0
6	B	107	0	0	2	0
All	All	7739	0	7387	63	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 4.

All (63) 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:4[B]:GLU:O	1:A:4[B]:GLU:HG2	1.63	0.95
1:B:306:LYS:H	1:B:306:LYS:CD	2.00	0.75
1:A:195:ASP:OD1	1:A:196:PRO:HD2	1.87	0.74
2:B:501:HEM:HBC2	2:B:501:HEM:HMC2	1.77	0.67
1:A:195:ASP:CG	1:A:196:PRO:CD	2.68	0.67
1:A:56:ARG:NH2	1:A:344:GLU:OE1	2.29	0.66
1:A:195:ASP:CG	1:A:196:PRO:HD2	2.22	0.64
1:A:4[B]:GLU:O	1:A:4[B]:GLU:CG	2.42	0.64
1:B:192:ASN:N	1:B:193:PRO:HD3	2.16	0.60
1:B:196:PRO:O	1:B:199:ASP:HB2	2.02	0.59
1:A:4[A]:GLU:O	1:A:4[A]:GLU:HG3	2.05	0.56
2:A:501:HEM:HBC2	2:A:501:HEM:HMC1	1.87	0.56
1:B:148:LEU:HD21	1:B:413:VAL:HG21	1.89	0.55
1:B:9:LYS:NZ	6:B:602:HOH:O	2.40	0.54
1:B:306:LYS:H	1:B:306:LYS:HD2	1.71	0.53
1:A:39:ILE:HA	1:A:51:TYR:O	2.09	0.52
1:B:378:ARG:HG2	6:B:677:HOH:O	2.09	0.51
1:A:158:PHE:CE2	1:A:219:ILE:HD13	2.45	0.51
1:B:306:LYS:H	1:B:306:LYS:HD3	1.75	0.51

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:195:ASP:OD1	1:B:196:PRO:HD2	2.11	0.50
1:A:173:PHE:HB2	1:A:215:LEU:HD22	1.94	0.49
1:A:183:GLU:OE1	1:A:201:ASN:HB3	2.13	0.48
1:B:306:LYS:CD	1:B:306:LYS:N	2.74	0.48
1:B:192:ASN:N	1:B:193:PRO:CD	2.77	0.48
1:B:120:VAL:HG11	1:B:302:VAL:HG13	1.95	0.47
1:B:357:ILE:N	1:B:358:PRO:CD	2.78	0.47
1:A:210:LYS:HE3	1:A:214:ASP:OD2	2.14	0.47
1:A:150:LEU:HD11	1:A:174:ILE:HD11	1.97	0.46
1:A:272:LEU:N	1:A:327:THR:HG21	2.30	0.46
1:B:367:TRP:HB2	1:B:371:VAL:HG12	1.98	0.46
1:B:185:MET:CB	5:B:504:GOL:H12	2.46	0.45
1:B:287:LEU:C	1:B:287:LEU:HD23	2.41	0.45
1:B:68:ASP:HB3	1:B:334:TYR:CE1	2.51	0.45
1:B:298:LEU:HD22	1:B:303:PRO:HB3	1.98	0.45
1:B:222:ASP:O	1:B:226:SER:OG	2.34	0.45
1:A:104:LEU:HB3	1:A:105:PRO:HD3	1.98	0.44
1:A:357:ILE:N	1:A:358:PRO:CD	2.79	0.44
1:B:173:PHE:CE2	1:B:262:LEU:HD13	2.52	0.44
1:B:185:MET:HB3	5:B:504:GOL:H31	1.99	0.44
1:B:158:PHE:CD1	1:B:158:PHE:N	2.85	0.44
1:A:288[B]:GLN:HA	1:A:288[B]:GLN:OE1	2.18	0.44
1:B:366:ILE:HG21	1:B:389:ALA:HB1	2.00	0.44
1:B:87:PHE:CD1	4:B:503:DMS:H11	2.52	0.44
2:B:501:HEM:HBC2	2:B:501:HEM:CMC	2.44	0.44
1:B:326:PRO:HG3	1:B:357:ILE:HG22	2.00	0.43
1:A:195:ASP:OD1	1:A:196:PRO:CD	2.60	0.43
1:B:17:LEU:N	1:B:18:PRO:CD	2.82	0.43
1:B:207:GLU:O	1:B:211:VAL:HG23	2.18	0.43
1:B:150:LEU:HD21	1:B:174:ILE:HG13	2.00	0.42
1:B:357:ILE:CG2	1:B:361:HIS:CE1	3.03	0.42
1:B:81:PHE:HB3	1:B:209:ILE:HG12	2.00	0.42
1:B:191:ALA:C	1:B:193:PRO:HD3	2.45	0.42
1:B:150:LEU:HD21	1:B:174:ILE:CD1	2.50	0.41
2:A:501:HEM:HMB2	2:A:501:HEM:HBB2	2.03	0.41
1:B:419:LYS:O	1:B:451:LYS:HD3	2.21	0.41
1:B:73:GLN:HG3	1:B:77:PHE:CE2	2.55	0.41
1:B:60:GLU:OE2	1:B:342:GLY:HA2	2.20	0.41
1:B:330:ALA:HB1	1:B:354:MET:CE	2.49	0.41
1:A:342:GLY:O	1:A:344:GLU:HG2	2.20	0.41
1:A:366:ILE:HG21	1:A:389:ALA:HB1	2.02	0.41

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:203:ARG:NH2	1:B:207:GLU:OE2	2.53	0.41
2:B:501:HEM:HMB2	2:B:501:HEM:HBB2	2.02	0.41
1:A:354:MET:HE3	1:A:354:MET:HA	2.04	0.40

There are no symmetry-related clashes.

5.3 Torsion angles (i)

5.3.1 Protein backbone (i)

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

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

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
1	A	454/456 (100%)	438 (96%)	14 (3%)	2 (0%)	30	28
1	B	455/456 (100%)	437 (96%)	18 (4%)	0	100	100
All	All	909/912 (100%)	875 (96%)	32 (4%)	2 (0%)	48	44

All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	4[A]	GLU
1	A	4[B]	GLU

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	400/400 (100%)	396 (99%)	4 (1%)	68	76

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles
1	B	401/400 (100%)	397 (99%)	4 (1%)	68 76
All	All	801/800 (100%)	793 (99%)	8 (1%)	68 76

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	49	THR
1	A	230	SER
1	A	354	MET
1	A	356	LEU
1	B	226	SER
1	B	306	LYS
1	B	354	MET
1	B	391	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (10) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	73	GLN
1	A	229	GLN
1	A	239	ASN
1	A	310	GLN
1	B	109	GLN
1	B	229	GLN
1	B	239	ASN
1	B	285	HIS
1	B	359	GLN
1	B	420	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 oligosaccharides in this entry.

5.6 Ligand geometry (i)

9 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	A	501	4,1	50,50,50	1.65	11 (22%)	66,82,82	1.57	11 (16%)
3	BWX	A	502	-	27,27,27	0.95	0	35,36,36	1.25	4 (11%)
2	HEM	B	501	4,1	50,50,50	1.57	7 (14%)	66,82,82	1.74	16 (24%)
4	DMS	A	503	2	3,3,3	0.44	0	3,3,3	0.24	0
5	GOL	A	504	-	5,5,5	0.13	0	5,5,5	0.40	0
5	GOL	B	505	-	5,5,5	0.11	0	5,5,5	0.26	0
5	GOL	B	504	-	5,5,5	0.15	0	5,5,5	0.37	0
3	BWX	B	502	-	27,27,27	0.95	1 (3%)	35,36,36	1.56	6 (17%)
4	DMS	B	503	2	3,3,3	0.39	0	3,3,3	0.28	0

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
5	GOL	A	504	-	-	2/4/4/4	-
2	HEM	A	501	4,1	-	3/14/54/54	-
3	BWX	A	502	-	-	5/24/24/24	0/2/2/2
2	HEM	B	501	4,1	-	2/14/54/54	-
5	GOL	B	505	-	-	2/4/4/4	-
3	BWX	B	502	-	-	3/24/24/24	0/2/2/2
5	GOL	B	504	-	-	1/4/4/4	-

All (19) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	501	HEM	FE-NB	5.13	2.10	1.94
2	A	501	HEM	FE-NB	4.49	2.08	1.94
2	A	501	HEM	C1B-NB	-3.82	1.33	1.40
2	A	501	HEM	FE-NC	3.81	2.08	1.95
2	B	501	HEM	FE-NC	3.71	2.07	1.95
2	B	501	HEM	C4B-NB	-3.47	1.31	1.38
2	B	501	HEM	C1B-NB	-3.28	1.34	1.40
2	B	501	HEM	C1C-NC	-2.76	1.34	1.39
2	A	501	HEM	C4D-ND	-2.70	1.35	1.40
2	A	501	HEM	C1D-ND	-2.64	1.33	1.38
2	A	501	HEM	C4B-NB	-2.51	1.33	1.38
3	B	502	BWX	C8-C11	-2.46	1.48	1.52
2	A	501	HEM	C1C-NC	-2.45	1.35	1.39
2	A	501	HEM	C1C-C2C	-2.36	1.40	1.45
2	B	501	HEM	C4D-C3D	2.35	1.49	1.45
2	A	501	HEM	C3C-C4C	-2.31	1.42	1.46
2	B	501	HEM	C4D-ND	-2.25	1.36	1.40
2	A	501	HEM	O1A-CGA	2.23	1.29	1.22
2	A	501	HEM	CMA-C3A	2.02	1.55	1.50

All (37) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	501	HEM	CHC-C4B-NB	5.26	130.14	124.42
2	A	501	HEM	CHC-C4B-NB	5.24	130.11	124.42
2	B	501	HEM	C1B-NB-C4B	4.63	109.86	105.07
3	B	502	BWX	C16-C14-N1	-4.28	101.77	110.79
2	A	501	HEM	C1B-NB-C4B	4.26	109.47	105.07
3	A	502	BWX	O1-C13-C11	-3.87	116.91	122.15
3	B	502	BWX	C12-C11-C8	-3.80	103.82	112.92
2	A	501	HEM	C4C-NC-C1C	3.61	108.88	105.35
2	B	501	HEM	CHD-C4C-NC	3.37	128.08	124.44
3	B	502	BWX	O1-C13-C11	-3.18	117.83	122.15
2	A	501	HEM	CMD-C2D-C1D	3.15	129.83	125.04
2	A	501	HEM	CHD-C4C-NC	3.10	127.79	124.44
2	B	501	HEM	O2A-CGA-CBA	3.07	123.90	114.03
2	B	501	HEM	C4C-NC-C1C	3.06	108.34	105.35
2	B	501	HEM	CHD-C1D-C2D	-2.98	120.32	124.98
2	B	501	HEM	CHD-C1D-ND	2.94	127.61	124.42
2	B	501	HEM	O1A-CGA-CBA	-2.87	113.87	123.08
3	B	502	BWX	C8-C11-C13	2.75	115.32	108.16
3	A	502	BWX	C8-C11-C13	2.72	115.26	108.16

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	A	501	HEM	CAD-C3D-C4D	2.57	129.16	124.66
2	B	501	HEM	CAD-C3D-C4D	2.53	129.08	124.66
2	A	501	HEM	CHA-C4D-C3D	-2.53	120.58	125.33
2	A	501	HEM	CHA-C4D-ND	2.44	127.38	124.37
2	A	501	HEM	O2A-CGA-CBA	2.44	121.86	114.03
2	B	501	HEM	C4A-C3A-C2A	2.39	109.63	106.83
3	B	502	BWX	C16-C14-C15	2.37	116.13	110.42
2	B	501	HEM	CMB-C2B-C1B	2.37	128.65	125.04
2	B	501	HEM	CHA-C4D-ND	2.29	127.20	124.37
2	B	501	HEM	CHA-C4D-C3D	-2.27	121.07	125.33
2	B	501	HEM	CAD-CBD-CGD	2.25	118.45	113.60
3	A	502	BWX	C14-N1-C13	2.18	126.34	121.67
2	A	501	HEM	CHD-C1D-C2D	-2.16	121.60	124.98
2	B	501	HEM	C3B-C2B-C1B	-2.15	104.89	106.49
3	A	502	BWX	O2-C15-C14	2.13	120.49	113.40
2	A	501	HEM	CHD-C1D-ND	2.11	126.71	124.42
2	B	501	HEM	O2D-CGD-O1D	-2.08	118.11	123.30
3	B	502	BWX	C10-C9-C8	-2.08	119.11	121.20

There are no chirality outliers.

All (18) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
3	A	502	BWX	C8-C11-C13-N1
3	A	502	BWX	C8-C11-C13-O1
3	B	502	BWX	C8-C11-C13-N1
3	B	502	BWX	C8-C11-C13-O1
5	A	504	GOL	C1-C2-C3-O3
5	A	504	GOL	O2-C2-C3-O3
5	B	505	GOL	C1-C2-C3-O3
5	B	504	GOL	C1-C2-C3-O3
3	B	502	BWX	C2-C1-C4-C5
5	B	505	GOL	O2-C2-C3-O3
3	A	502	BWX	C14-C16-C17-C18
3	A	502	BWX	C14-C16-C17-C22
2	A	501	HEM	CAD-CBD-CGD-O2D
2	B	501	HEM	CAD-CBD-CGD-O2D
3	A	502	BWX	C3-C1-C4-C5
2	A	501	HEM	CAD-CBD-CGD-O1D
2	B	501	HEM	CAD-CBD-CGD-O1D
2	A	501	HEM	CAA-CBA-CGA-O2A

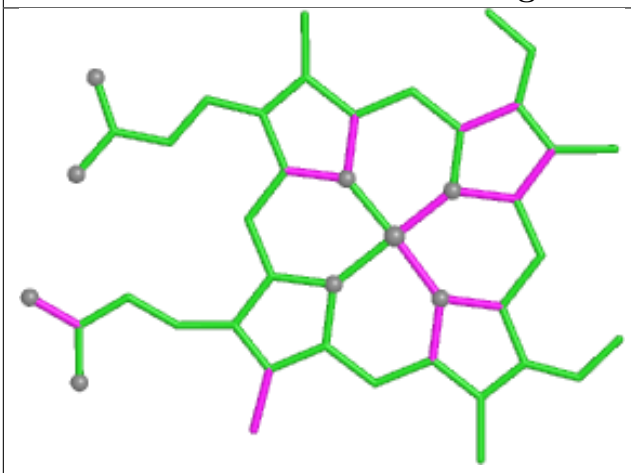
There are no ring outliers.

4 monomers are involved in 8 short contacts:

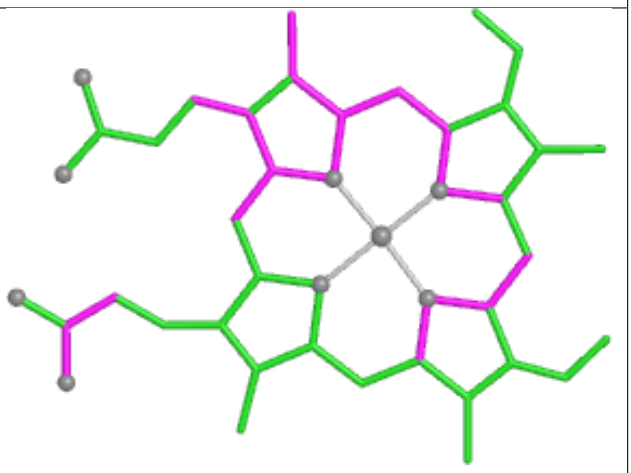
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	501	HEM	2	0
2	B	501	HEM	3	0
5	B	504	GOL	2	0
4	B	503	DMS	1	0

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.

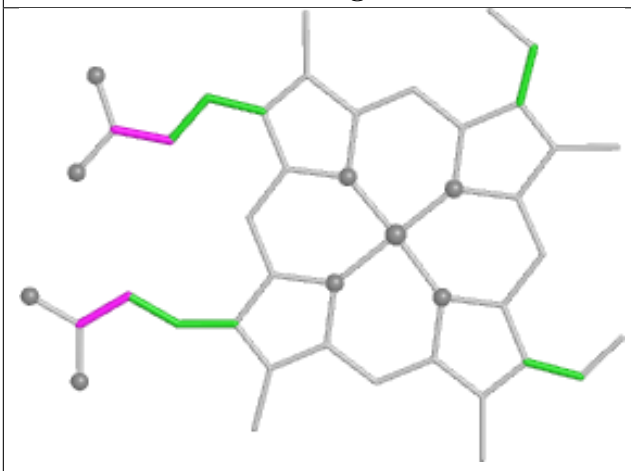
Ligand HEM A 501



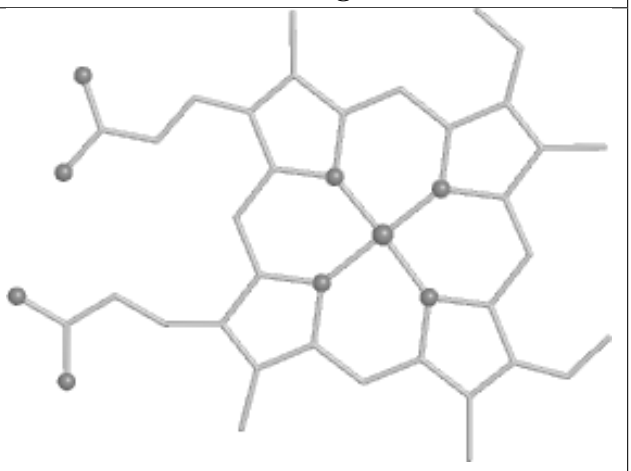
Bond lengths



Bond angles

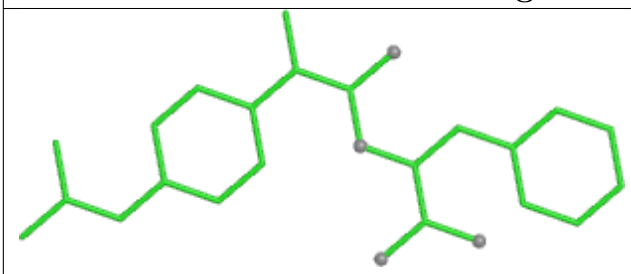


Torsions

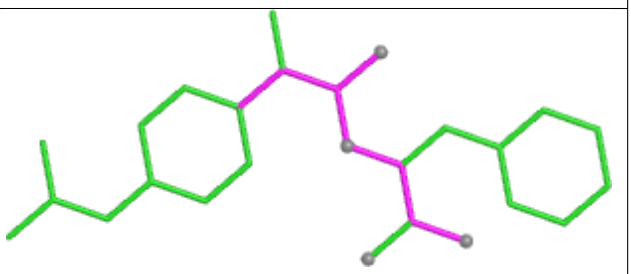


Rings

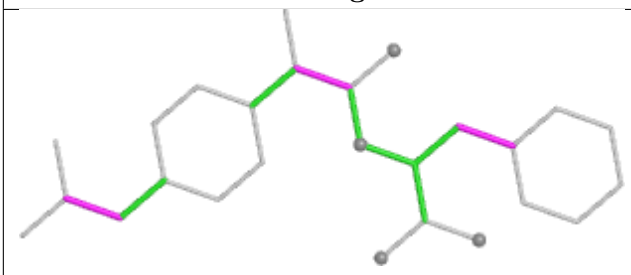
Ligand BWX A 502



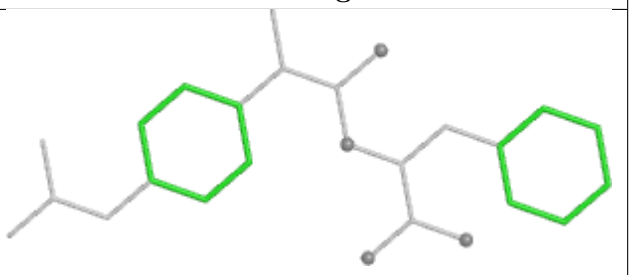
Bond lengths



Bond angles

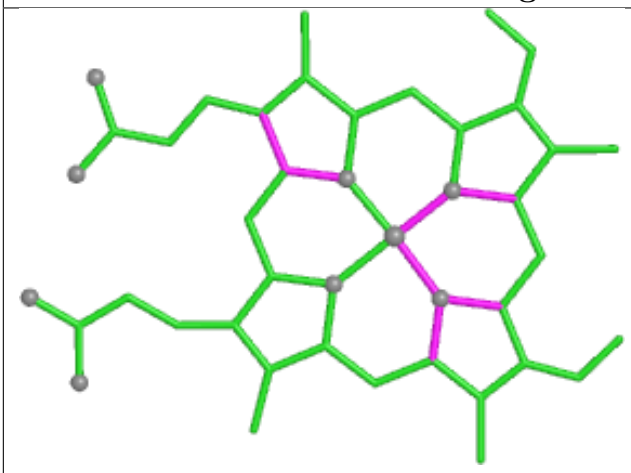


Torsions

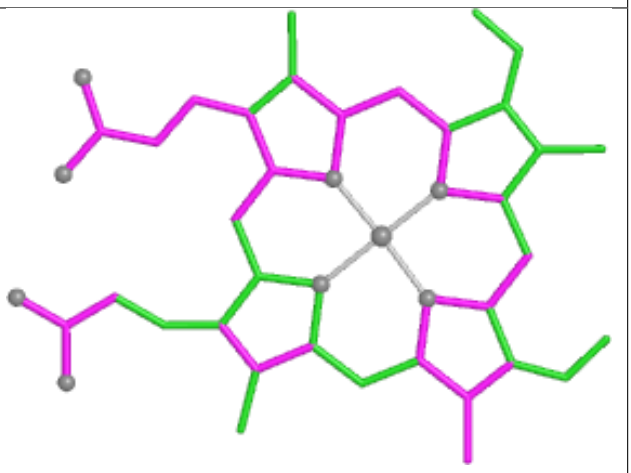


Rings

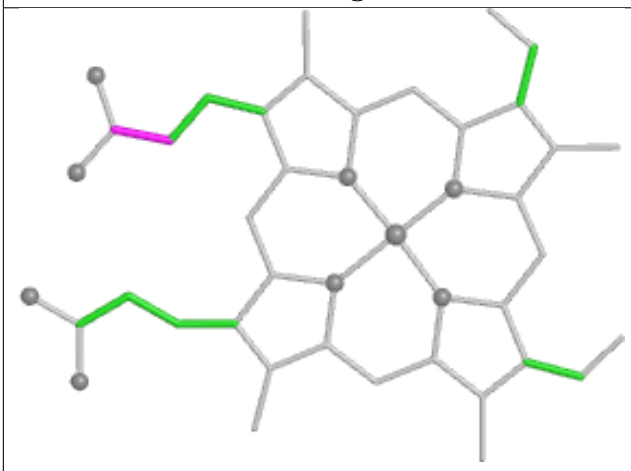
Ligand HEM B 501



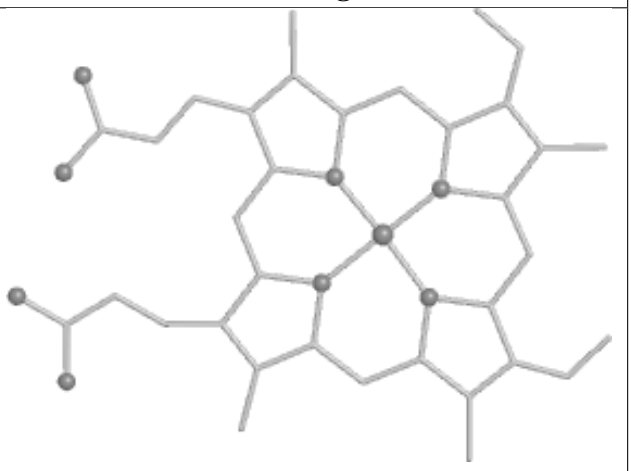
Bond lengths



Bond angles

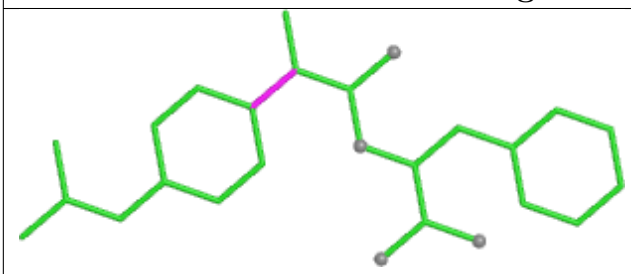


Torsions

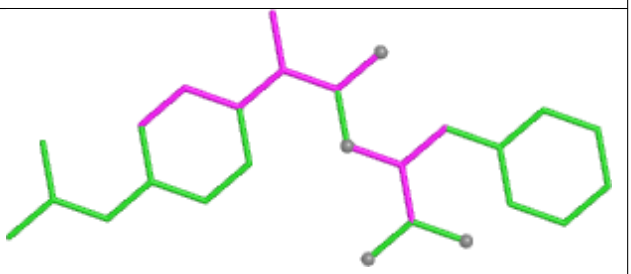


Rings

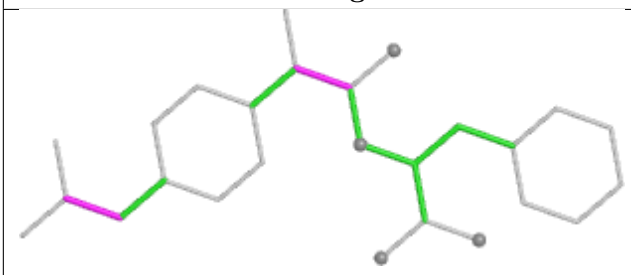
Ligand BWX B 502



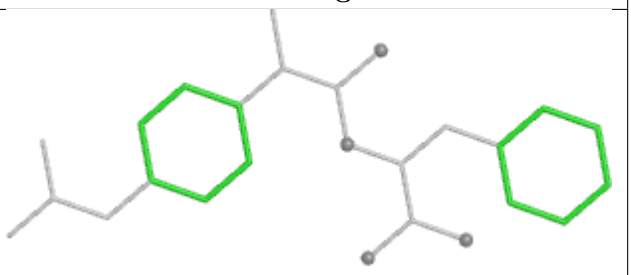
Bond lengths



Bond angles



Torsions



Rings

5.7 Other polymers (i)

There are no such residues in this entry.

5.8 Polymer linkage issues (i)

There are no chain breaks in this entry.

6 Fit of model and data (i)

6.1 Protein, DNA and RNA chains (i)

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

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2			OWAB(Å^2)	Q<0.9
1	A	453/456 (99%)	0.03	6 (1%)	75	77	20, 40, 66, 97	3 (0%)
1	B	453/456 (99%)	0.01	7 (1%)	72	74	18, 39, 74, 114	4 (0%)
All	All	906/912 (99%)	0.02	13 (1%)	73	75	18, 39, 69, 114	7 (0%)

All (13) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	197	ALA	4.5
1	B	191	ALA	4.2
1	A	191	ALA	4.0
1	B	197	ALA	4.0
1	B	13	GLU	3.1
1	B	196	PRO	2.9
1	B	158	PHE	2.6
1	B	198	TYR	2.4
1	A	196	PRO	2.4
1	B	193	PRO	2.4
1	A	198	TYR	2.3
1	A	193	PRO	2.1
1	A	455	LEU	2.0

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

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

6.3 Carbohydrates (i)

There are no oligosaccharides in this entry.

6.4 Ligands (i)

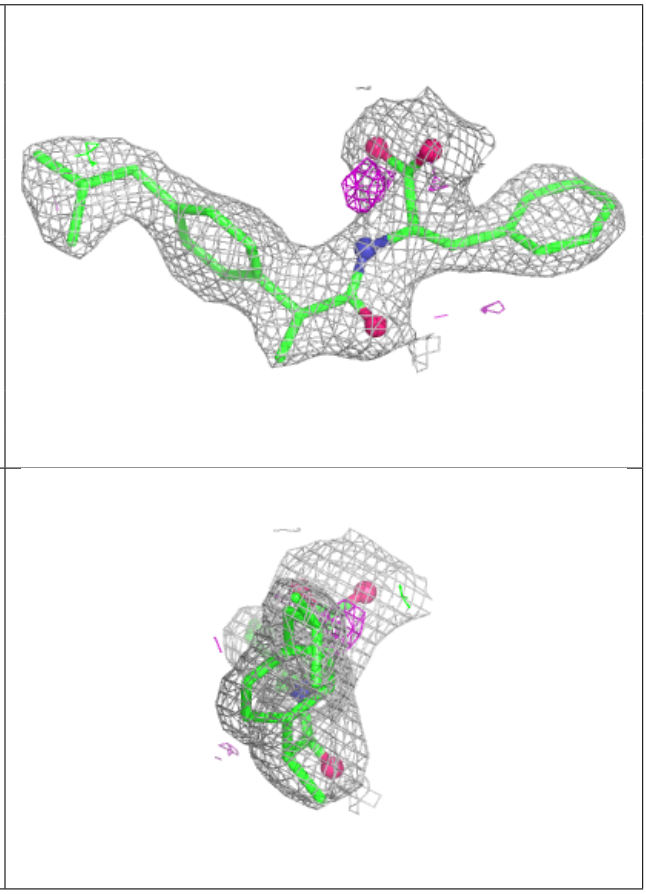
In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. The B-factors column lists the minimum, median, 95th percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists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

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
5	GOL	B	504	6/6	0.79	0.17	67,71,74,75	0
5	GOL	B	505	6/6	0.86	0.12	65,69,70,71	0
5	GOL	A	504	6/6	0.88	0.12	40,61,67,73	0
3	BWX	B	502	26/26	0.92	0.11	40,48,63,67	0
3	BWX	A	502	26/26	0.92	0.11	42,48,54,57	0
4	DMS	A	503	4/4	0.96	0.16	47,49,55,62	0
4	DMS	B	503	4/4	0.97	0.22	45,50,53,62	0
2	HEM	B	501	43/43	0.98	0.06	25,29,35,41	0
2	HEM	A	501	43/43	0.98	0.05	24,29,34,39	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.

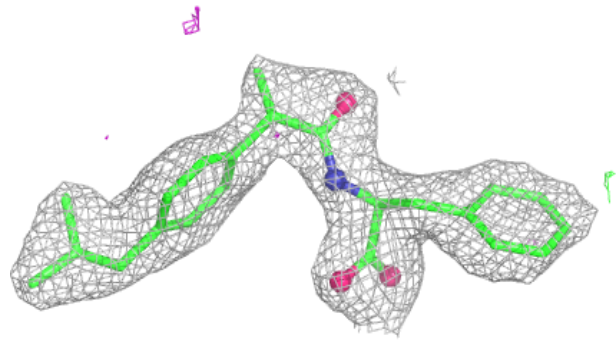
Electron density around BWX B 502:

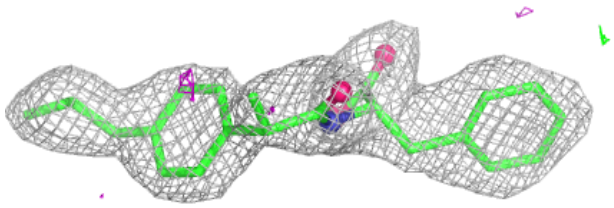
2mF$_o$-DF$_c$ (at 0.7 rmsd) in gray
mF$_o$-DF$_c$ (at 3 rmsd) in purple (negative)
and green (positive)

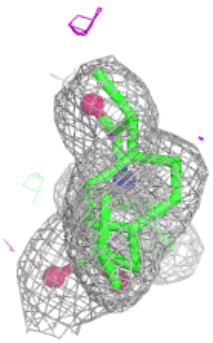


Electron density around BWX A 502:

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

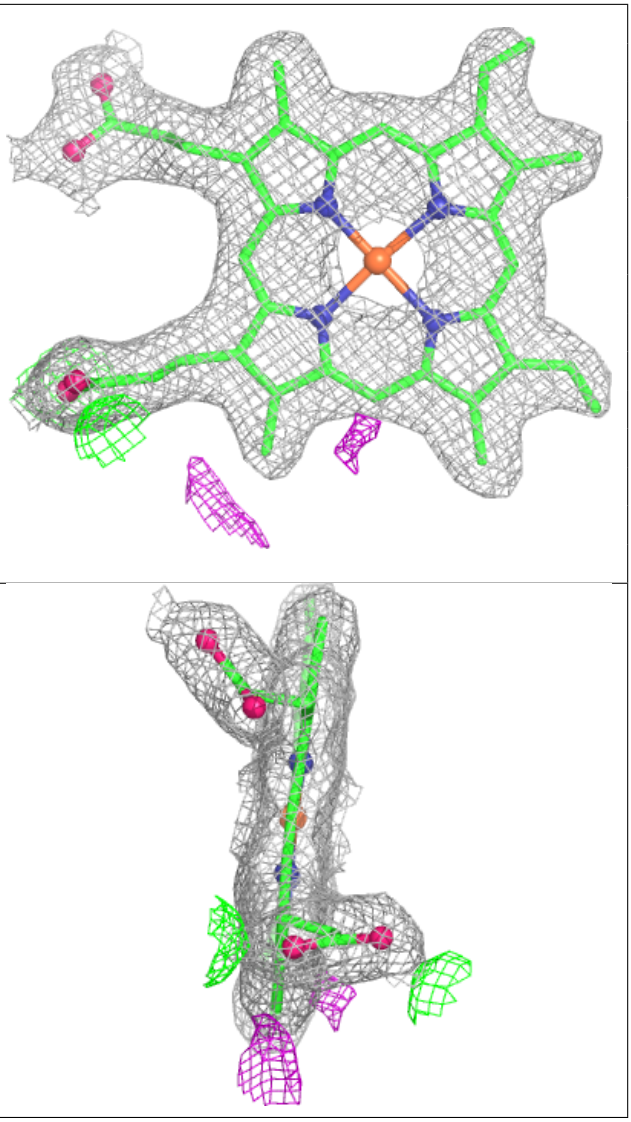






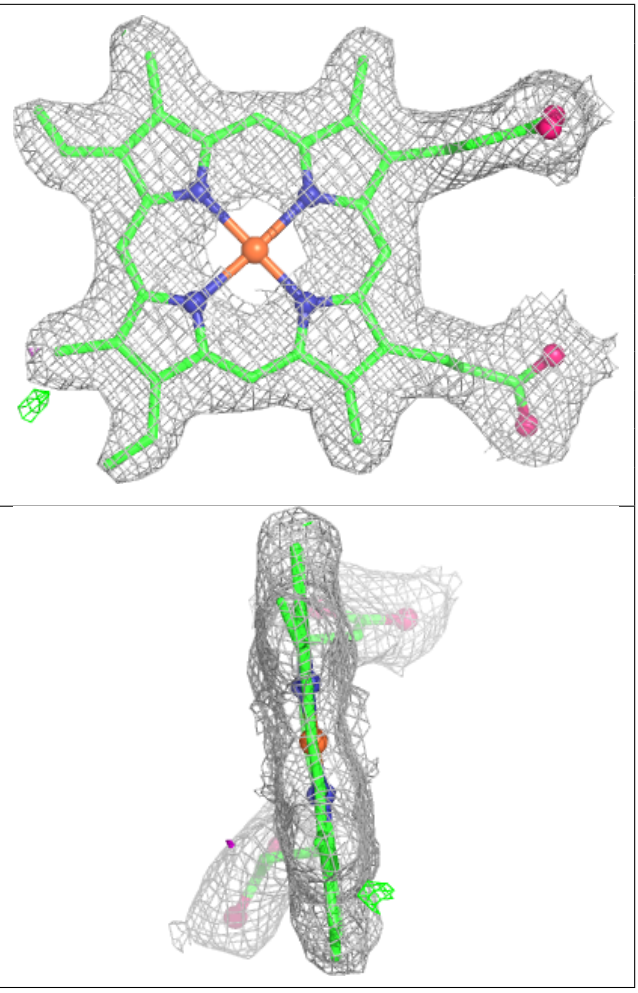
Electron density around HEM B 501:

$2mF_o$-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 501:

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.