Full wwPDB X-ray Structure Validation Report (i)

Apr 28, 2024 – 12:48 pm BST

PDB ID	:	5MJH
Title	:	X-ray generated oxyferrous/water mixed complex of DtpA from Streptomyces lividans
Authors	:	Moreno Chicano, T.; Chaplin, A.K.; Worrall, J.A.R.; Strange, R.W.; Hough, M.A.
Deposited on	:	2016-12-01
Resolution	:	1.45 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:

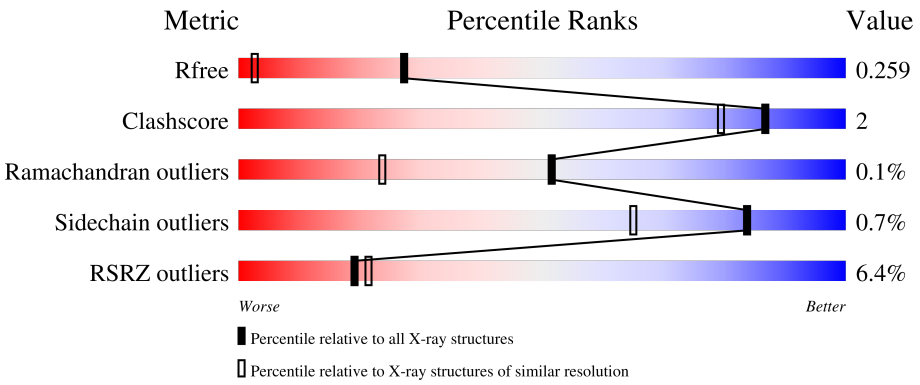
MolProbity	:	4.02b-467
Mogul	:	1.8.4, CSD as541be (2020)
Xtriage (Phenix)	:	1.13
EDS	:	2.36.2
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36.2

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 1.45 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	1156 (1.46-1.46)
Clashscore	141614	1202 (1.46-1.46)
Ramachandran outliers	138981	1178 (1.46-1.46)
Sidechain outliers	138945	1178 (1.46-1.46)
RSRZ outliers	127900	1139 (1.46-1.46)

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	377	7% 91% 7% •
1	B	377	6% 93% • •

2 Entry composition (i)

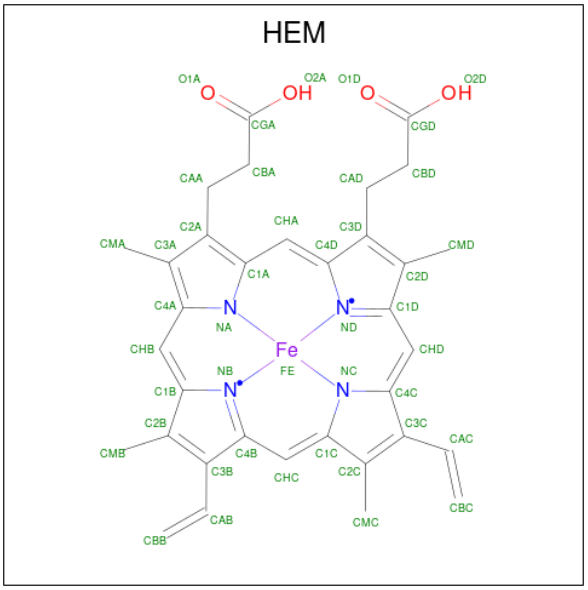
There are 4 unique types of molecules in this entry. The entry contains 6358 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 Dye type peroxidase A.

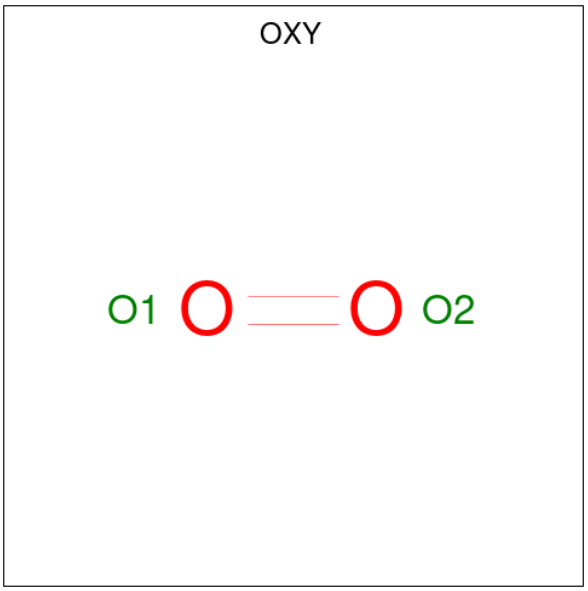
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	368	2812	1742	530	528	12	0	2	0
1	B	368	2833	1755	534	532	12	10	5	0

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



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
			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 OXYGEN MOLECULE (three-letter code: OXY) (formula: O_2).



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	1	Total O 2 2	0	1
3	B	1	Total O 2 2	0	1

- Molecule 4 is water.

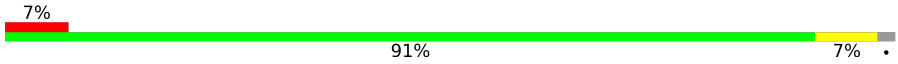
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	298	Total O 299 299	0	2
4	B	323	Total O 324 324	0	2

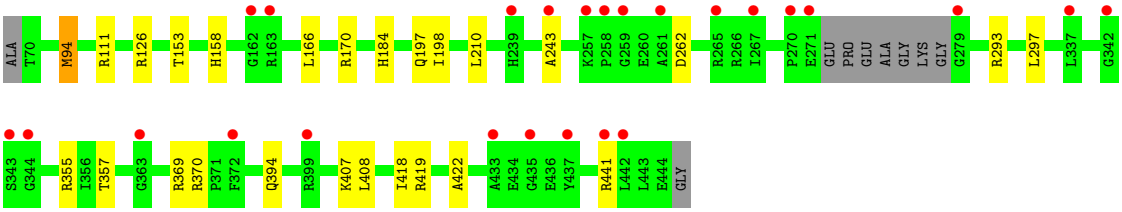
3 Residue-property plots (i)

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: Dye type peroxidase A

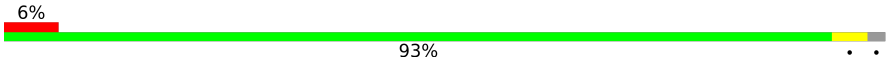
Chain A:

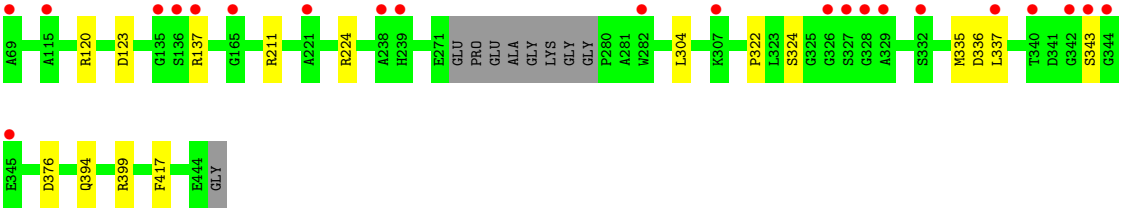




• Molecule 1: Dye type peroxidase A

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	59.78Å 70.63Å 77.64Å 90.00° 93.01° 90.00°	Depositor
Resolution (Å)	77.53 – 1.45 40.01 – 1.45	Depositor EDS
% Data completeness (in resolution range)	98.1 (77.53-1.45) 98.2 (40.01-1.45)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) >$ [1]	1.69 (at 1.45Å)	Xtriage
Refinement program	REFMAC 5.8.0155	Depositor
R, R_{free}	0.221 , 0.254 0.228 , 0.259	Depositor DCC
R_{free} test set	5527 reflections (4.93%)	wwPDB-VP
Wilson B-factor (Å^2)	10.5	Xtriage
Anisotropy	0.170	Xtriage
Bulk solvent k_{sol}(e/Å^3), B_{sol}(Å^2)	0.37 , 46.9	EDS
L-test for twinning[2]	$< \|L\| > = 0.48$, $< L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o,F_c correlation	0.92	EDS
Total number of atoms	6358	wwPDB-VP
Average B, all atoms (Å^2)	13.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.64% of the height of the origin peak. No significant pseudotranslation is detected.*

[1]Intensities estimated from amplitudes.

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

5 Model quality (i)

5.1 Standard geometry (i)

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

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	0.67	0/2874	0.90	5/3888 (0.1%)
1	B	0.71	0/2901	0.90	5/3922 (0.1%)
All	All	0.69	0/5775	0.90	10/7810 (0.1%)

There are no bond length outliers.

All (10) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
1	A	94[A]	MET	CG-SD-CE	6.85	111.17	100.20
1	A	94[B]	MET	CG-SD-CE	6.85	111.17	100.20
1	B	137	ARG	NE-CZ-NH1	5.73	123.17	120.30
1	B	211	ARG	NE-CZ-NH2	-5.70	117.45	120.30
1	B	399	ARG	NE-CZ-NH1	5.53	123.06	120.30
1	A	355	ARG	NE-CZ-NH1	5.52	123.06	120.30
1	A	369	ARG	NE-CZ-NH1	5.52	123.06	120.30
1	A	126	ARG	NE-CZ-NH1	5.44	123.02	120.30
1	B	123	ASP	CB-CG-OD1	5.39	123.15	118.30
1	B	224	ARG	NE-CZ-NH1	5.23	122.91	120.30

There are no chirality outliers.

There are no planarity outliers.

5.2 Too-close contacts (i)

In the following table, the Non-H and H(model) columns list the number of non-hydrogen atoms and hydrogen atoms in the chain respectively. The H(added) column lists the number of hydrogen atoms added and optimized by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, whereas Symm-Clashes lists symmetry-related clashes.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2812	0	2723	17	0
1	B	2833	0	2748	6	0
2	A	43	0	30	1	0
2	B	43	0	30	0	0
3	A	2	0	0	0	0
3	B	2	0	0	0	0
4	A	299	0	0	4	0
4	B	324	0	0	2	0
All	All	6358	0	5531	24	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 2.

All (24) 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:94[A]:MET:HE1	4:A:706:HOH:O	1.64	0.97
1:A:94[A]:MET:CE	4:A:706:HOH:O	2.21	0.86
1:A:197:GLN:OE1	4:A:501:HOH:O	2.07	0.71
1:B:322:PRO:HB3	1:B:335[B]:MET:SD	2.45	0.57
1:A:94[A]:MET:HE3	1:A:370:ARG:HG2	1.88	0.56
1:A:153:THR:HG21	4:A:501:HOH:O	2.07	0.53
1:A:94[B]:MET:SD	1:A:243:ALA:HB2	2.50	0.51
1:B:376:ASP:OD1	4:B:501:HOH:O	2.19	0.50
1:A:184:HIS:CG	1:A:419:ARG:HD2	2.49	0.47
1:A:94[A]:MET:HE3	1:A:370:ARG:CG	2.45	0.47
1:B:304:LEU:HD11	1:B:417:PHE:CZ	2.49	0.46
1:A:111:ARG:CZ	1:A:166:LEU:HD21	2.46	0.46
1:A:357:THR:HG22	1:A:408:LEU:HD21	1.98	0.45
1:A:94[A]:MET:CE	1:A:370:ARG:HG2	2.47	0.44
1:A:158:HIS:CD2	1:A:170:ARG:HD2	2.53	0.44
1:A:262:ASP:OD1	1:A:262:ASP:N	2.48	0.43
1:A:94[A]:MET:HE3	1:A:370:ARG:C	2.38	0.43
2:A:1401:HEM:HMB2	2:A:1401:HEM:HBB2	2.01	0.42
1:B:335[A]:MET:SD	1:B:337:LEU:HD21	2.60	0.42
1:A:297:LEU:HD13	1:A:418:ILE:CA	2.49	0.41
1:B:324:SER:HB2	1:B:336:ASP:HB3	2.02	0.41
1:A:293:ARG:HB3	1:A:422:ALA:HB3	2.02	0.41
1:B:120:ARG:NH1	4:B:514:HOH:O	2.48	0.40
1:A:198:ILE:CD1	1:A:210:LEU:HA	2.52	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	366/377 (97%)	358 (98%)	8 (2%)	0	100	100
1	B	369/377 (98%)	361 (98%)	7 (2%)	1 (0%)	41	18
All	All	735/754 (98%)	719 (98%)	15 (2%)	1 (0%)	51	24

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	343	SER

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	283/285 (99%)	280 (99%)	3 (1%)	73	48
1	B	286/285 (100%)	285 (100%)	1 (0%)	92	82
All	All	569/570 (100%)	565 (99%)	4 (1%)	84	65

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	394	GLN
1	A	407	LYS
1	A	441	ARG
1	B	394	GLN

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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)

4 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	B	1401	1,3	41,50,50	1.40	5 (12%)	45,82,82	2.30	17 (37%)
2	HEM	A	1401	1,3	41,50,50	1.39	8 (19%)	45,82,82	1.83	10 (22%)
3	OXY	B	1402[A]	2	1,1,1	0.03	0	-		
3	OXY	A	1402[A]	2	1,1,1	0.01	0	-		

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	HEM	B	1401	1,3	-	4/12/54/54	-
2	HEM	A	1401	1,3	-	4/12/54/54	-

All (13) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	1401	HEM	C1B-NB	-4.23	1.33	1.40
2	B	1401	HEM	CHB-C1B	3.35	1.43	1.35
2	B	1401	HEM	C1B-NB	-3.33	1.34	1.40
2	A	1401	HEM	C4D-ND	-2.90	1.35	1.40
2	B	1401	HEM	C4B-NB	-2.65	1.33	1.38
2	B	1401	HEM	FE-NB	2.55	2.09	1.96
2	B	1401	HEM	C4D-ND	-2.54	1.35	1.40
2	A	1401	HEM	FE-NB	2.50	2.09	1.96
2	A	1401	HEM	C3B-C4B	2.50	1.49	1.44
2	A	1401	HEM	O1D-CGD	2.43	1.30	1.22
2	A	1401	HEM	C4D-C3D	2.30	1.49	1.45
2	A	1401	HEM	CHB-C1B	2.17	1.40	1.35
2	A	1401	HEM	O2A-CGA	-2.03	1.23	1.30

All (27) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1401	HEM	C1B-NB-C4B	6.29	111.57	105.07
2	B	1401	HEM	CBA-CAA-C2A	5.54	122.08	112.62
2	A	1401	HEM	CHC-C4B-NB	5.01	129.88	124.43
2	A	1401	HEM	C1B-NB-C4B	4.45	109.67	105.07
2	B	1401	HEM	CHC-C4B-NB	3.96	128.74	124.43
2	B	1401	HEM	C3B-C2B-C1B	3.92	109.40	106.49
2	B	1401	HEM	C4B-C3B-C2B	-3.70	104.18	107.11
2	B	1401	HEM	CHB-C1B-NB	3.41	128.60	124.38
2	A	1401	HEM	CHA-C4D-ND	3.37	128.54	124.38
2	A	1401	HEM	CMD-C2D-C1D	3.25	129.98	125.04
2	B	1401	HEM	C2C-C3C-C4C	-3.23	104.64	106.90
2	A	1401	HEM	CBA-CAA-C2A	3.16	118.00	112.62
2	A	1401	HEM	CHA-C4D-C3D	-3.14	119.44	125.33
2	B	1401	HEM	CHA-C4D-C3D	-2.87	119.94	125.33
2	A	1401	HEM	C4A-C3A-C2A	2.81	108.95	107.00
2	B	1401	HEM	CMA-C3A-C4A	-2.71	124.31	128.46
2	B	1401	HEM	C3D-C4D-ND	2.51	112.95	110.17
2	A	1401	HEM	CMC-C2C-C3C	2.50	129.36	124.68
2	A	1401	HEM	CHD-C1D-C2D	-2.50	121.07	124.98
2	B	1401	HEM	C2B-C1B-NB	-2.49	106.89	109.84
2	B	1401	HEM	CAD-C3D-C4D	2.39	128.83	124.66
2	A	1401	HEM	CMA-C3A-C4A	-2.37	124.82	128.46
2	B	1401	HEM	CHA-C4D-ND	2.20	127.10	124.38
2	B	1401	HEM	CHD-C1D-C2D	-2.17	121.59	124.98
2	B	1401	HEM	CAA-CBA-CGA	-2.15	107.72	113.76

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Mol	Chain	Res	Type	Atoms	Z	Observed(^{o})	Ideal(^{o})
2	B	1401	HEM	C4C-CHD-C1D	2.14	125.39	122.56
2	B	1401	HEM	O2D-CGD-O1D	-2.05	118.20	123.30

There are no chirality outliers.

All (8) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	1401	HEM	CAA-CBA-CGA-O2A
2	A	1401	HEM	CAA-CBA-CGA-O1A
2	B	1401	HEM	CAD-CBD-CGD-O1D
2	B	1401	HEM	CAA-CBA-CGA-O2A
2	B	1401	HEM	CAD-CBD-CGD-O2D
2	B	1401	HEM	CAA-CBA-CGA-O1A
2	A	1401	HEM	CAD-CBD-CGD-O2D
2	A	1401	HEM	CAD-CBD-CGD-O1D

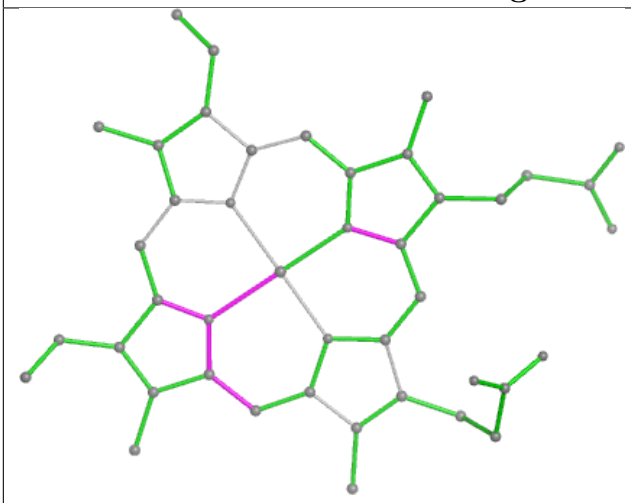
There are no ring outliers.

1 monomer is involved in 1 short contact:

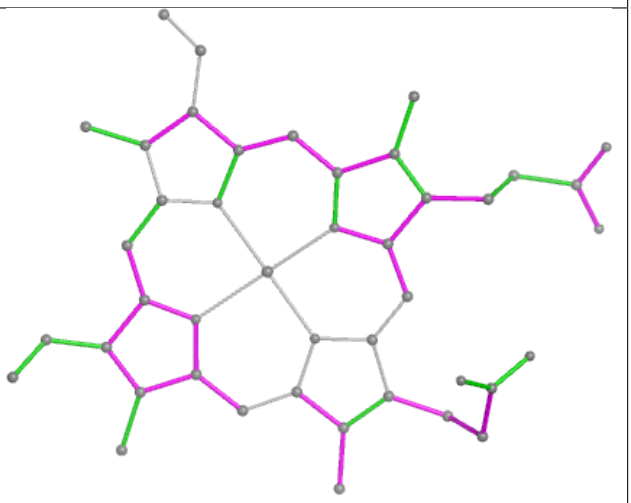
Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	1401	HEM	1	0

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.

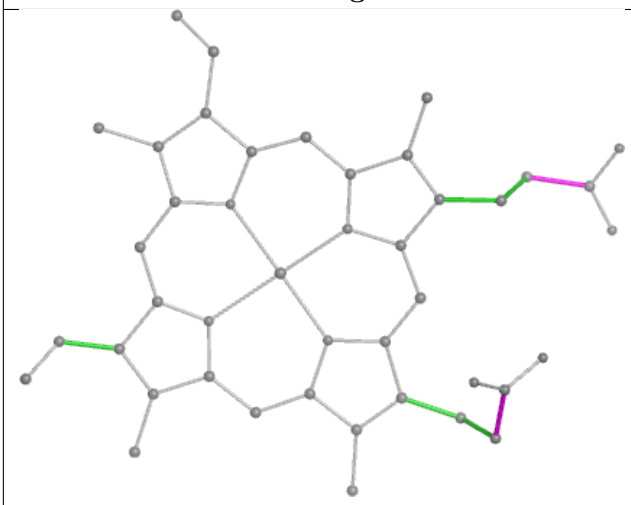
Ligand HEM B 1401



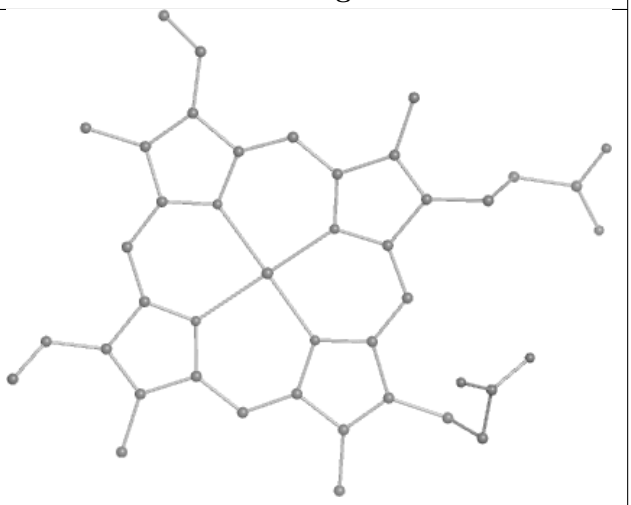
Bond lengths



Bond angles

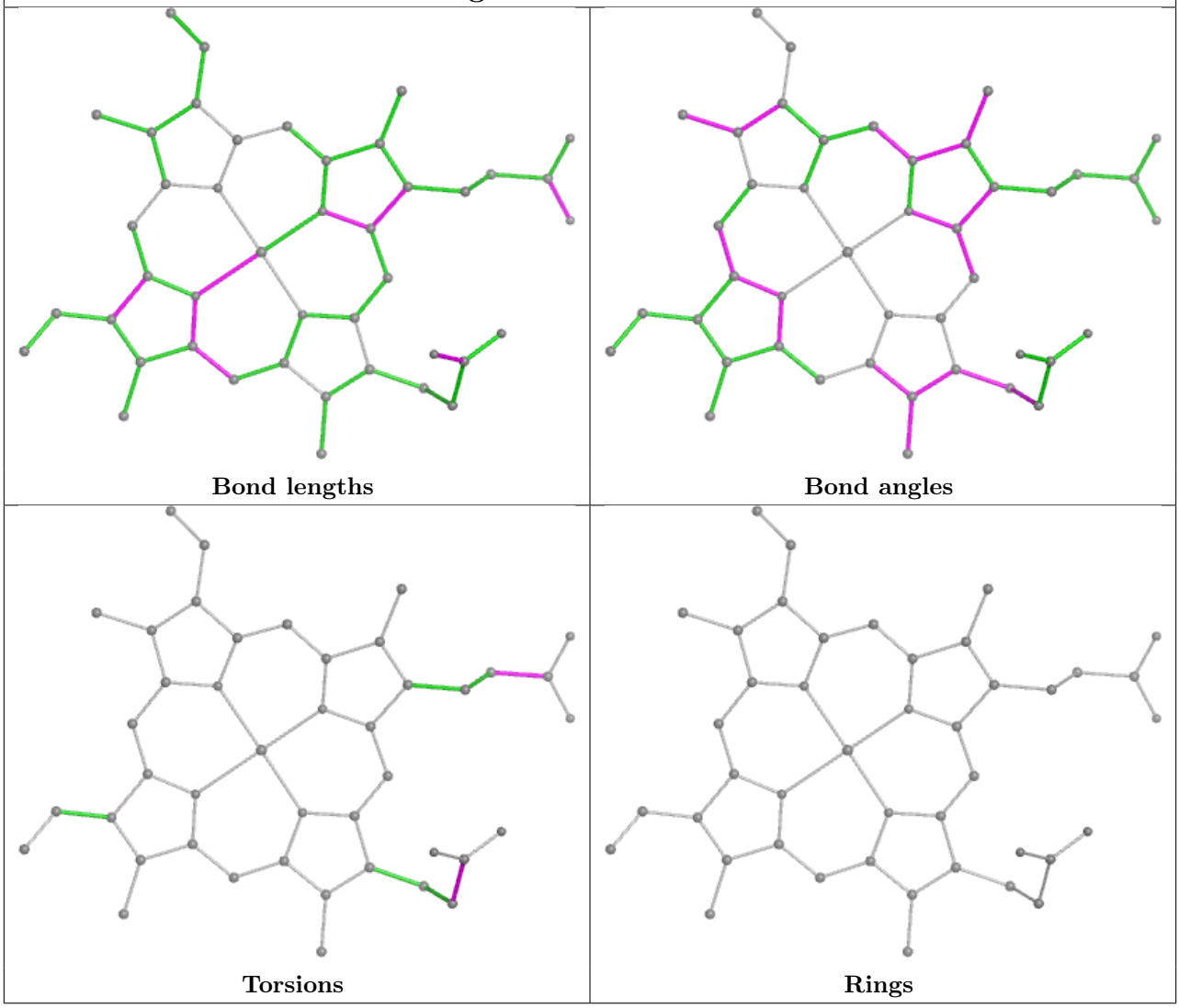


Torsions



Rings

Ligand HEM A 1401



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	368/377 (97%)	0.79	25 (6%)	17	19	6, 12, 24, 34	24 (6%)
1	B	368/377 (97%)	0.71	22 (5%)	21	24	5, 11, 23, 34	16 (4%)
All	All	736/754 (97%)	0.75	47 (6%)	19	21	5, 12, 24, 34	40 (5%)

All (47) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	136	SER	7.8
1	A	261	ALA	7.4
1	A	279	GLY	6.7
1	A	342	GLY	5.0
1	A	343	SER	4.7
1	B	332	SER	4.5
1	B	239	HIS	4.3
1	B	337	LEU	4.3
1	B	69	ALA	4.2
1	B	135	GLY	4.1
1	B	342	GLY	3.9
1	A	441	ARG	3.4
1	B	343	SER	3.2
1	B	340	THR	2.9
1	A	442	LEU	2.9
1	B	344	GLY	2.8
1	A	257	LYS	2.8
1	B	326	GLY	2.7
1	A	344	GLY	2.7
1	B	345	GLU	2.6
1	A	162	GLY	2.5
1	A	337	LEU	2.5
1	A	437	TYR	2.5
1	A	243	ALA	2.4

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Mol	Chain	Res	Type	RSRZ
1	B	329	ALA	2.4
1	A	435	GLY	2.4
1	A	163	ARG	2.4
1	A	259	GLY	2.3
1	B	115	ALA	2.3
1	B	137	ARG	2.2
1	B	165	GLY	2.2
1	A	270	PRO	2.2
1	A	433	ALA	2.2
1	B	282	TRP	2.2
1	A	372	PHE	2.2
1	A	399	ARG	2.2
1	B	327	SER	2.1
1	A	258	PRO	2.1
1	B	307	LYS	2.1
1	A	267	ILE	2.1
1	A	239	HIS	2.1
1	A	271	GLU	2.0
1	A	363	GLY	2.0
1	B	328	GLY	2.0
1	A	265	ARG	2.0
1	B	221	ALA	2.0
1	B	238	ALA	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 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.

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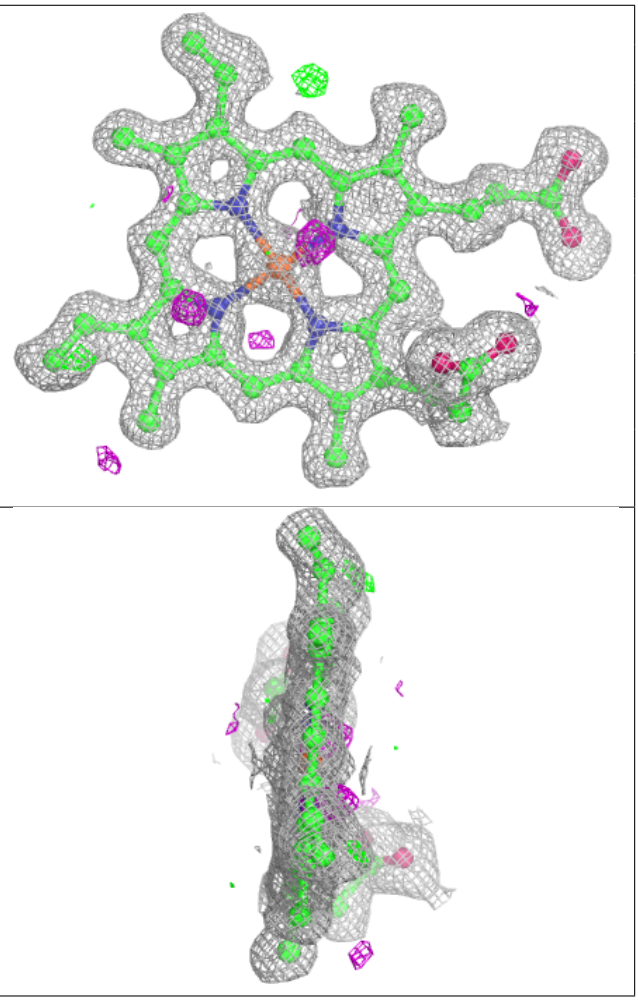
Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9

Mol	Type	Chain	Res	Atoms	RSCC	RSR	B-factors(Å^2)	Q<0.9
3	OXY	A	1402[A]	2/2	0.93	0.15	4,4,4,4	2
3	OXY	B	1402[A]	2/2	0.94	0.14	6,6,6,9	2
2	HEM	A	1401	43/43	0.97	0.10	7,7,8,9	0
2	HEM	B	1401	43/43	0.97	0.10	6,7,9,10	0

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

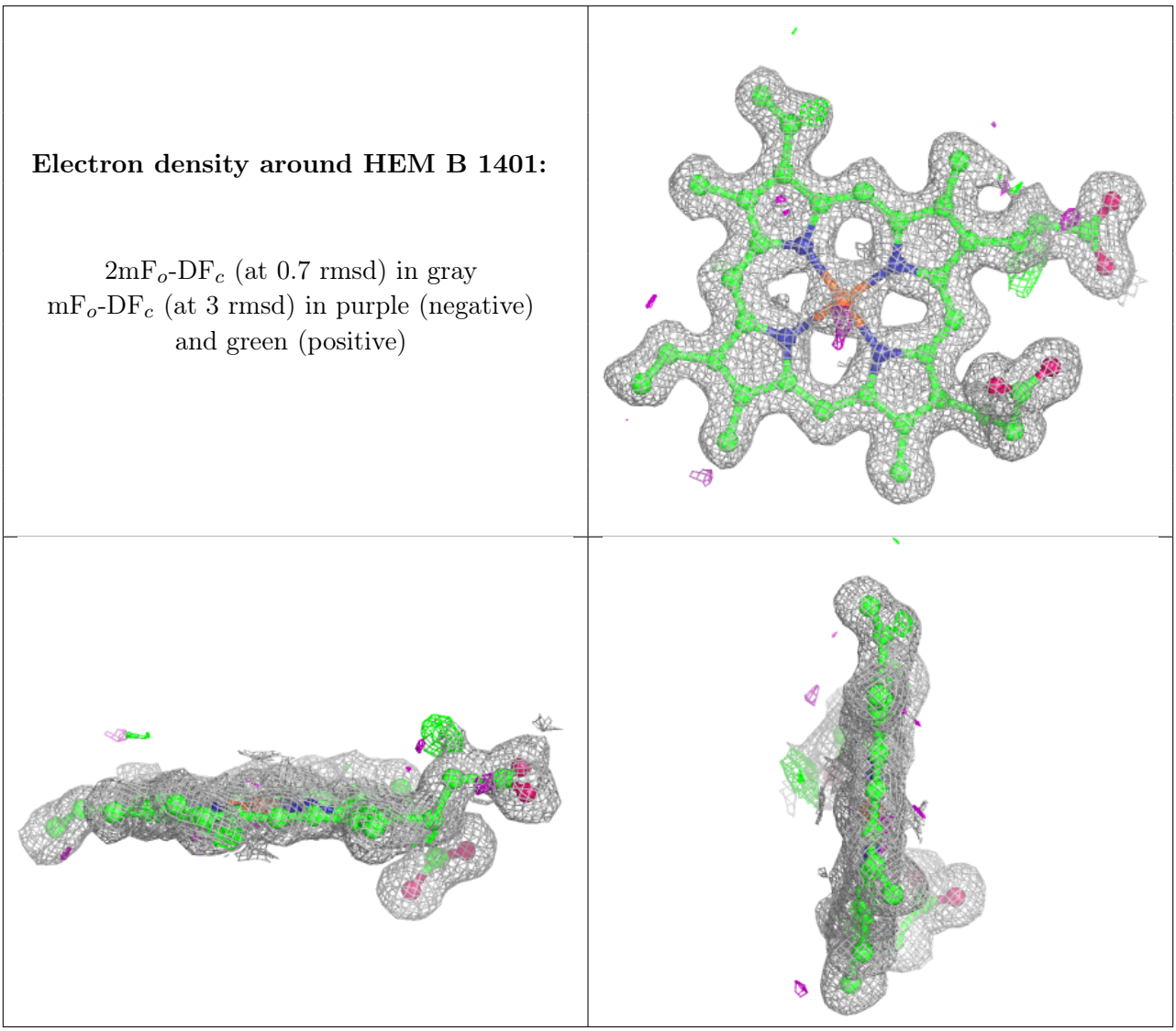
Electron density around HEM A 1401:

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 1401:

$2mF_o$-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.