



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 4, 2023 – 09:58 PM EDT

PDB ID : 6VDY
Title : Crystal Structure of Dehaloperoxidase B wild type in Complex with Substrate Trichlorophenol
Authors : Ghiladi, R.A.; de Serrano, V.S.; McGuire, A.; Malewschik, T.
Deposited on : 2019-12-27
Resolution : 1.70 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 : **FAILED**
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : 1.13
EDS : **FAILED**
buster-report : 1.1.7 (2018)
Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.35.1

1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 1.70 Å.

There are no overall percentile quality scores available for this entry.

MolProbity and EDS failed to run properly - the sequence quality summary graphics cannot be shown.

2 Entry composition i

There are 6 unique types of molecules in this entry. The entry contains 2726 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 Dehaloperoxidase B.

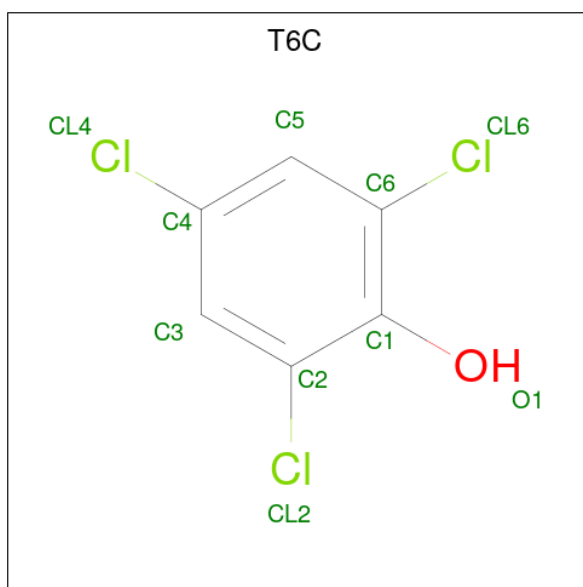
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	137	Total	C	N	O	S	0	25	0
			1239	775	208	248	8			
1	B	137	Total	C	N	O	S	0	21	0
			1196	749	200	239	8			

- Molecule 2 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$) (labeled as "Ligand of Interest" by depositor).



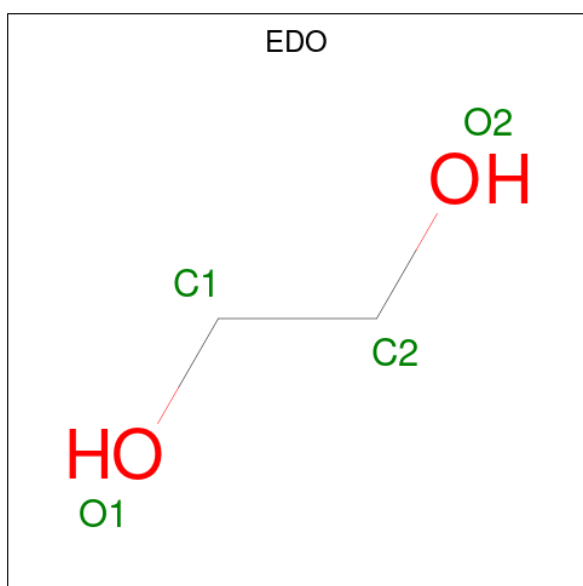
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	Fe	N			O
2	A	1	Total	C	Fe	N	O	0	0
			39	32	1	4	2		
2	B	1	Total	C	Fe	N	O	0	0
			39	32	1	4	2		

- Molecule 3 is 2,4,6-trichlorophenol (three-letter code: T6C) (formula: $C_6H_3Cl_3O$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	
			Total	C	Cl			O
3	A	1	10	6	3	1	0	0
3	B	1	10	6	3	1	0	0

- Molecule 4 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: $C_2H_6O_2$).



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

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

- Molecule 5 is SULFATE ION (three-letter code: SO4) (formula: O₄S).



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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	79	Total	O	0	1
			80	80		
6	B	87	Total	O	0	0
			87	87		

MolProbity and EDS failed to run properly - this section is therefore empty.

3 Data and refinement statistics

EDS failed to run properly - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	59.23Å 67.46Å 68.13Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	37.26 – 1.70	Depositor
% Data completeness (in resolution range)	98.8 (37.26-1.70)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.40 (at 1.70Å)	Xtrriage
Refinement program	PHENIX dev_3714, PHENIX dev_3714	Depositor
R, R_{free}	0.179 , 0.214	Depositor
Wilson B-factor (Å ²)	23.1	Xtrriage
Anisotropy	0.193	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	0.078 for -h,l,k	Xtrriage
Total number of atoms	2726	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 9.44% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

²Theoretical values of $\langle |L| \rangle$, $\langle L^2 \rangle$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.

4 Model quality [i](#)

4.1 Standard geometry [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.2 Too-close contacts [i](#)

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4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3.2 Protein sidechains [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.3.3 RNA [i](#)

MolProbity failed to run properly - this section is therefore empty.

4.4 Non-standard residues in protein, DNA, RNA chains [i](#)

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

4.5 Carbohydrates [i](#)

There are no monosaccharides in this entry.

4.6 Ligand geometry [i](#)

10 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
5	SO4	A	204	-	4,4,4	0.57	0	6,6,6	1.14	1 (16%)
2	HEM	B	402	1	37,46,50	1.81	12 (32%)	40,77,82	2.11	14 (35%)
4	EDO	A	203	-	3,3,3	0.79	0	2,2,2	0.67	0
2	HEM	A	201	1	37,46,50	1.76	9 (24%)	40,77,82	1.93	11 (27%)
4	EDO	B	401	-	3,3,3	0.56	0	2,2,2	0.25	0
3	T6C	A	202	-	10,10,10	1.18	1 (10%)	14,14,14	1.46	4 (28%)
4	EDO	B	405	-	3,3,3	0.36	0	2,2,2	0.34	0
3	T6C	B	403	-	10,10,10	1.30	1 (10%)	14,14,14	1.55	5 (35%)
5	SO4	A	205	-	4,4,4	0.40	0	6,6,6	0.81	0
4	EDO	B	404	-	3,3,3	0.82	0	2,2,2	0.87	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	402	1	-	3/7/49/54	-
4	EDO	A	203	-	-	0/1/1/1	-
2	HEM	A	201	1	-	2/7/49/54	-
4	EDO	B	401	-	-	0/1/1/1	-
3	T6C	A	202	-	-	-	0/1/1/1
4	EDO	B	405	-	-	0/1/1/1	-
3	T6C	B	403	-	-	-	0/1/1/1
4	EDO	B	404	-	-	0/1/1/1	-

All (23) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	402	HEM	C3C-CAC	4.14	1.56	1.47
2	B	402	HEM	C3C-C2C	-3.98	1.34	1.40
2	A	201	HEM	C3C-CAC	3.89	1.55	1.47
2	A	201	HEM	FE-ND	3.69	2.15	1.96
2	A	201	HEM	C3C-C2C	-3.62	1.35	1.40
2	A	201	HEM	FE-NB	3.39	2.13	1.96
2	B	402	HEM	FE-NB	3.38	2.13	1.96
2	B	402	HEM	CAB-C3B	3.07	1.55	1.47

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	402	HEM	FE-ND	3.01	2.11	1.96
2	A	201	HEM	CAB-C3B	2.99	1.55	1.47
2	B	402	HEM	CHB-C1B	2.58	1.41	1.35
3	B	403	T6C	C4-CL4	2.50	1.80	1.74
2	B	402	HEM	CHA-C4D	2.39	1.41	1.35
2	A	201	HEM	CAD-C3D	2.34	1.55	1.50
2	B	402	HEM	CAD-C3D	2.30	1.55	1.50
2	A	201	HEM	C4A-NA	2.28	1.40	1.36
2	B	402	HEM	C2C-C1C	2.14	1.47	1.42
2	B	402	HEM	CMD-C2D	2.12	1.55	1.50
2	A	201	HEM	CHB-C1B	2.08	1.40	1.35
2	A	201	HEM	CMD-C2D	2.05	1.55	1.50
2	B	402	HEM	C4D-C3D	2.03	1.48	1.44
3	A	202	T6C	C4-CL4	2.03	1.78	1.74
2	B	402	HEM	CMB-C2B	2.00	1.55	1.50

All (35) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	402	HEM	C1B-NB-C4B	5.41	110.67	105.07
2	A	201	HEM	C4D-ND-C1D	4.64	109.86	105.07
2	A	201	HEM	C1B-NB-C4B	4.58	109.80	105.07
2	B	402	HEM	C4B-CHC-C1C	4.32	128.26	122.56
2	B	402	HEM	C4D-ND-C1D	4.31	109.53	105.07
2	A	201	HEM	C4C-CHD-C1D	3.81	127.58	122.56
2	A	201	HEM	C4B-CHC-C1C	3.71	127.45	122.56
2	B	402	HEM	C3D-C4D-ND	-3.41	105.80	109.84
2	B	402	HEM	C2B-C1B-NB	-3.39	105.82	109.84
2	B	402	HEM	CHD-C1D-ND	3.23	127.94	124.43
2	A	201	HEM	C3D-C4D-ND	-3.00	106.28	109.84
2	B	402	HEM	C4C-CHD-C1D	2.99	126.50	122.56
2	B	402	HEM	CAA-CBA-CGA	-2.85	105.76	113.76
2	B	402	HEM	C3B-C2B-C1B	2.82	108.58	106.49
2	A	201	HEM	C2B-C1B-NB	-2.73	106.60	109.84
3	A	202	T6C	C5-C6-CL6	2.65	122.77	118.49
2	B	402	HEM	CAD-C3D-C4D	2.61	129.01	125.04
3	B	403	T6C	C3-C4-CL4	2.49	122.26	119.15
2	B	402	HEM	CBA-CAA-C2A	-2.45	108.45	112.62
3	B	403	T6C	C1-C2-CL2	-2.44	115.72	118.78
2	A	201	HEM	C3B-C2B-C1B	2.37	108.24	106.49
2	A	201	HEM	C2D-C1D-ND	-2.29	107.14	109.88
3	B	403	T6C	C5-C6-CL6	2.28	122.17	118.49

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	402	HEM	CAD-C3D-C2D	-2.22	122.13	126.75
5	A	204	SO4	O4-S-O3	2.22	118.52	109.06
2	A	201	HEM	CAA-CBA-CGA	-2.21	107.56	113.76
3	B	403	T6C	C3-C2-CL2	2.19	122.03	118.49
3	B	403	T6C	C6-C5-C4	2.17	121.14	118.71
3	A	202	T6C	C1-C2-CL2	-2.15	116.08	118.78
2	B	402	HEM	CHC-C4B-C3B	2.14	127.85	124.57
2	A	201	HEM	CAD-C3D-C4D	2.13	128.28	125.04
3	A	202	T6C	C3-C2-CL2	2.13	121.92	118.49
2	A	201	HEM	C3C-C4C-NC	-2.12	106.95	110.94
3	A	202	T6C	C5-C6-C1	-2.07	118.98	121.78
2	B	402	HEM	C4A-C3A-C2A	2.01	108.39	107.00

There are no chirality outliers.

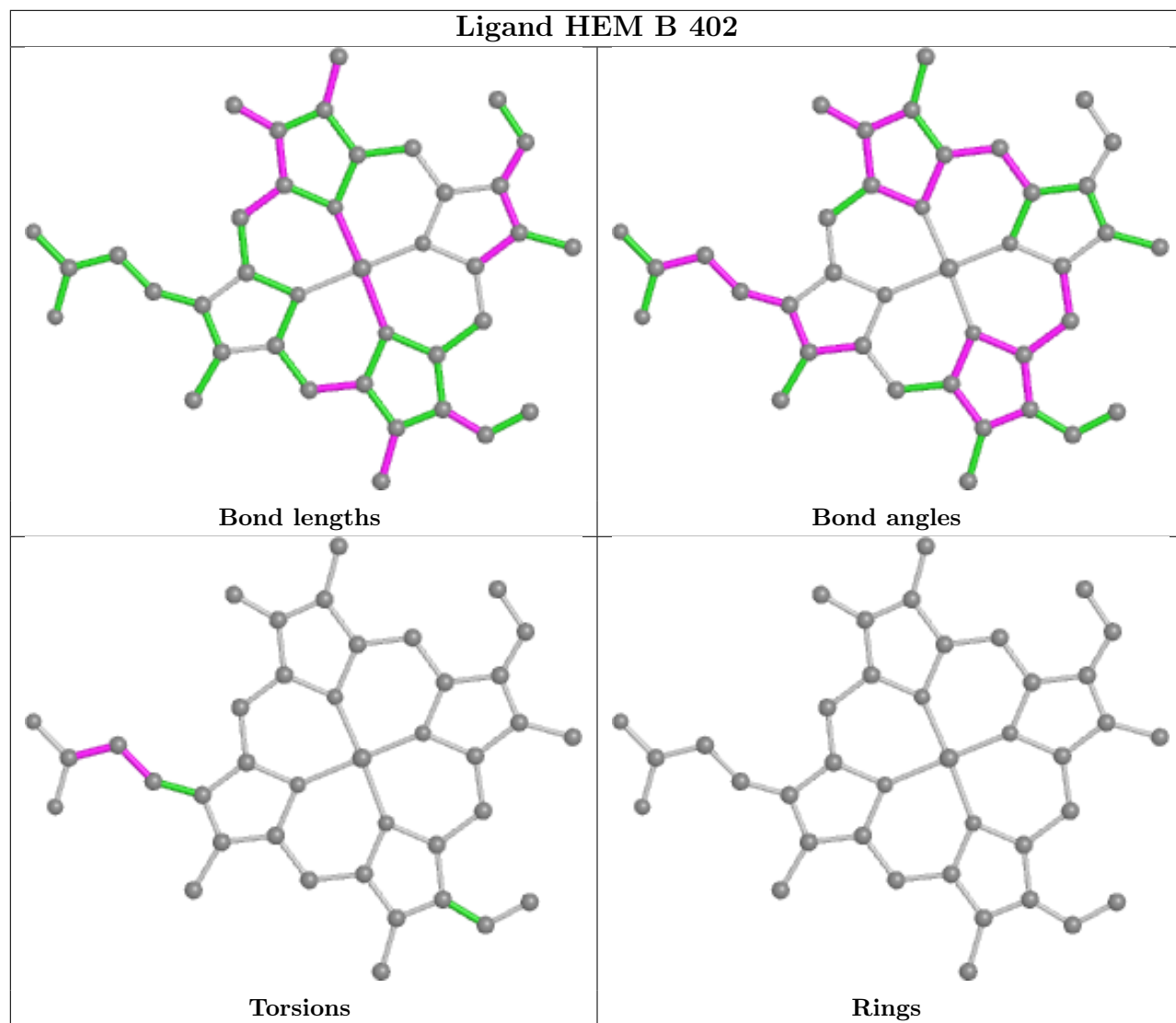
All (5) torsion outliers are listed below:

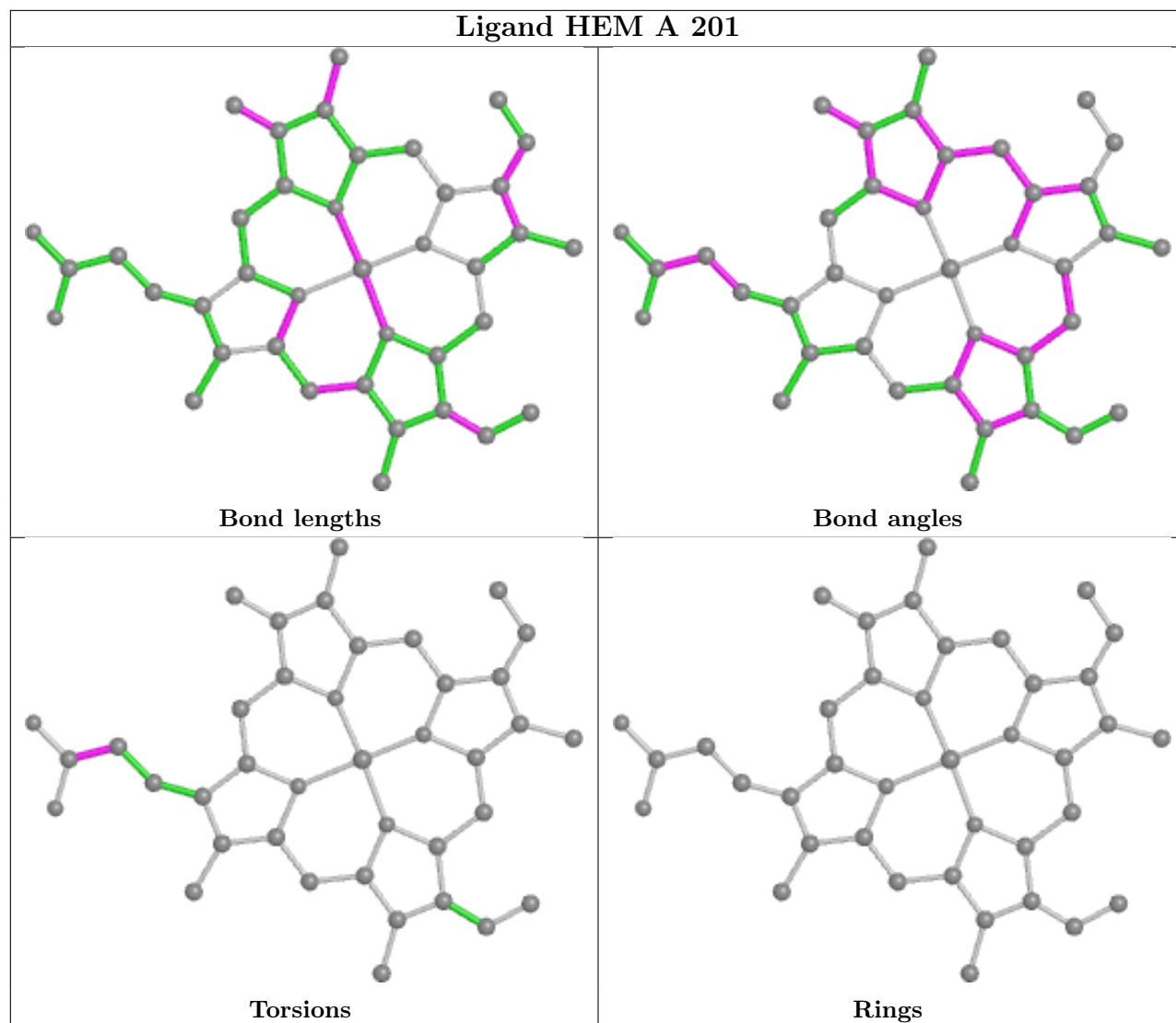
Mol	Chain	Res	Type	Atoms
2	B	402	HEM	C2A-CAA-CBA-CGA
2	B	402	HEM	CAA-CBA-CGA-O1A
2	B	402	HEM	CAA-CBA-CGA-O2A
2	A	201	HEM	CAA-CBA-CGA-O1A
2	A	201	HEM	CAA-CBA-CGA-O2A

There are no ring outliers.

No monomer is involved in short contacts.

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 than 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.





4.7 Other polymers [\(i\)](#)

There are no such residues in this entry.

4.8 Polymer linkage issues [\(i\)](#)

There are no chain breaks in this entry.

5 Fit of model and data

5.1 Protein, DNA and RNA chains

EDS failed to run properly - this section is therefore empty.

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

EDS failed to run properly - this section is therefore empty.

5.3 Carbohydrates

EDS failed to run properly - this section is therefore empty.

5.4 Ligands

EDS failed to run properly - this section is therefore empty.

5.5 Other polymers

EDS failed to run properly - this section is therefore empty.