



Full wwPDB X-ray Structure Validation Report ⓘ

Oct 2, 2023 – 11:59 PM EDT

PDB ID : 6PMZ
Title : Structure of rat neuronal nitric oxide synthase heme domain in complex with
7-(5-(Aminomethyl)pyridin-3-yl)-4-methylquinolin-2-amine
Authors : Li, H.; Poulos, T.L.
Deposited on : 2019-07-02
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 ⓘ 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 2.10 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 7 unique types of molecules in this entry. The entry contains 7012 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 Nitric oxide synthase, brain.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	409	Total 3338	C 2138	N 571	O 607	S 22	0	3	0
1	B	411	Total 3351	C 2144	N 574	O 612	S 21	0	2	0

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



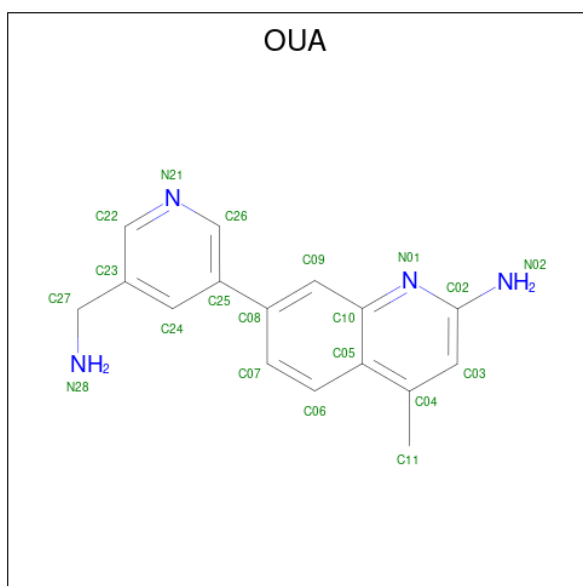
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
			Total	C	Fe	N			O
2	A	1	Total 43	C 34	Fe 1	N 4	O 4	0	0
2	B	1	Total 43	C 34	Fe 1	N 4	O 4	0	0

- Molecule 3 is 5,6,7,8-TETRAHYDROBIOPTERIN (three-letter code: H4B) (formula: $C_9H_{15}N_5O_3$).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
3	A	1	17	9	5	3	0	0
3	B	1	17	9	5	3	0	0

- Molecule 4 is 7-[5-(aminomethyl)pyridin-3-yl]-4-methylquinolin-2-amine (three-letter code: OUA) (formula: $C_{16}H_{16}N_4$).



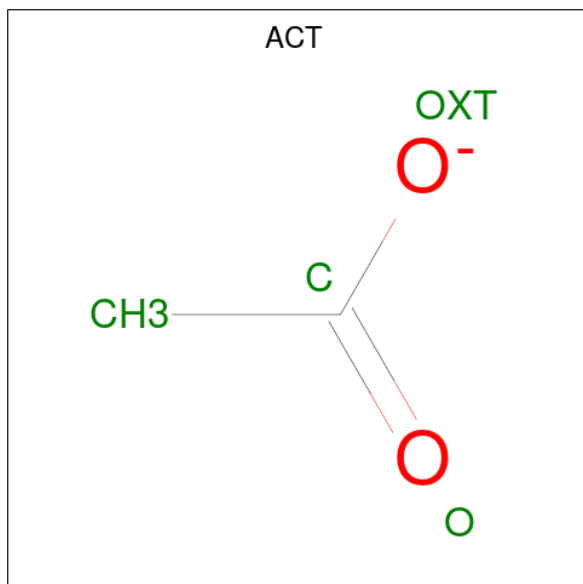
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
			Total	C	N	O		
4	A	1	20	16	4	0	0	

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Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
			Total	C	N		
4	B	1	20	16	4	0	0

- Molecule 5 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).



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

- Molecule 6 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	Zn		
6	A	1	1	1	0	0

- Molecule 7 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
7	A	69	69	69	0	0
7	B	85	85	85	0	0

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	51.44Å 110.87Å 164.38Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.06 – 2.10	Depositor
% Data completeness (in resolution range)	95.5 (36.06-2.10)	Depositor
R_{merge}	0.13	Depositor
R_{sym}	0.13	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.55 (at 2.10Å)	Xtrriage
Refinement program	PHENIX 1.9_1692	Depositor
R, R_{free}	0.185 , 0.246	Depositor
Wilson B-factor (Å ²)	37.2	Xtrriage
Anisotropy	1.368	Xtrriage
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	7012	wwPDB-VP
Average B, all atoms (Å ²)	68.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.86% of the height of the origin peak. No significant pseudotranslation is detected.*

¹ Intensities estimated from amplitudes.

² Theoretical 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](#)

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

4.3 Torsion angles [i](#)

4.3.1 Protein backbone [i](#)

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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](#)

Of 9 ligands modelled in this entry, 1 is monoatomic - leaving 8 for Mogul analysis.

In the following table, the Counts columns list the number of bonds (or angles) for which Mogul statistics could be retrieved, the number of bonds (or angles) that are observed in the model and the number of bonds (or angles) that are defined in the Chemical Component Dictionary. The Link column lists molecule types, if any, to which the group is linked. The Z score for a bond

length (or angle) is the number of standard deviations the observed value is removed from the expected value. A bond length (or angle) with $|Z| > 2$ is considered an outlier worth inspection. RMSZ is the root-mean-square of all Z scores of the bond lengths (or angles).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
4	OUA	A	803	-	22,22,22	0.87	1 (4%)	30,31,31	1.82	8 (26%)
2	HEM	B	801	1	41,50,50	1.91	7 (17%)	45,82,82	1.98	16 (35%)
4	OUA	B	803	-	22,22,22	0.92	1 (4%)	30,31,31	1.58	7 (23%)
5	ACT	B	804	-	3,3,3	0.77	0	3,3,3	0.71	0
5	ACT	A	804	-	3,3,3	0.77	0	3,3,3	0.72	0
2	HEM	A	801	1	41,50,50	1.90	5 (12%)	45,82,82	1.69	9 (20%)
3	H4B	B	802	-	16,18,18	0.86	0	11,26,26	2.41	5 (45%)
3	H4B	A	802	-	16,18,18	0.78	0	11,26,26	2.69	6 (54%)

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

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
4	OUA	A	803	-	-	4/6/6/6	0/3/3/3
2	HEM	B	801	1	-	2/12/54/54	-
4	OUA	B	803	-	-	4/6/6/6	0/3/3/3
2	HEM	A	801	1	-	8/12/54/54	-
3	H4B	B	802	-	-	0/8/17/17	0/2/2/2
3	H4B	A	802	-	-	0/8/17/17	0/2/2/2

All (14) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	801	HEM	C3D-C2D	7.87	1.53	1.36
2	B	801	HEM	C3D-C2D	7.54	1.52	1.36
2	B	801	HEM	C3C-C2C	-4.00	1.34	1.40
2	A	801	HEM	C3C-C2C	-3.84	1.35	1.40
2	B	801	HEM	C3C-CAC	3.57	1.55	1.47
2	A	801	HEM	C3C-CAC	3.39	1.54	1.47
2	B	801	HEM	CAB-C3B	2.96	1.55	1.47
2	A	801	HEM	CAB-C3B	2.87	1.55	1.47
2	B	801	HEM	FE-NB	2.65	2.10	1.96
2	A	801	HEM	FE-NB	2.56	2.09	1.96
2	B	801	HEM	CMB-C2B	2.28	1.55	1.50

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Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
4	A	803	OUA	C02-N01	2.17	1.36	1.33
2	B	801	HEM	CMD-C2D	2.17	1.55	1.50
4	B	803	OUA	C05-C10	-2.01	1.39	1.42

All (51) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	801	HEM	C4D-ND-C1D	6.30	111.58	105.07
3	B	802	H4B	C8A-C4A-C4	5.08	119.08	114.57
3	A	802	H4B	C8A-C4A-C4	4.87	118.90	114.57
2	A	801	HEM	C4D-ND-C1D	4.66	109.89	105.07
4	A	803	OUA	C04-C05-C10	4.18	120.28	118.01
2	B	801	HEM	C4B-CHC-C1C	3.96	127.78	122.56
3	A	802	H4B	C4-C4A-N5	3.65	122.18	119.12
4	A	803	OUA	C26-N21-C22	3.59	122.37	117.48
2	B	801	HEM	CBA-CAA-C2A	-3.53	106.59	112.62
2	A	801	HEM	C4C-CHD-C1D	3.46	127.12	122.56
3	A	802	H4B	N1-C2-N3	-3.44	120.02	125.42
4	A	803	OUA	C05-C10-N01	-3.42	119.18	122.81
2	A	801	HEM	CAD-CBD-CGD	-3.28	106.55	113.60
4	B	803	OUA	C26-N21-C22	3.23	121.89	117.48
4	B	803	OUA	C24-C25-C08	-3.16	115.64	120.86
3	A	802	H4B	C2-N1-C8A	3.15	121.59	114.54
3	A	802	H4B	C2-N3-C4	3.03	120.74	115.93
3	B	802	H4B	N1-C2-N3	-2.92	120.83	125.42
2	A	801	HEM	C3B-C2B-C1B	2.85	108.60	106.49
2	B	801	HEM	CBD-CAD-C3D	-2.82	104.80	112.63
3	B	802	H4B	C2-N3-C4	2.80	120.38	115.93
2	B	801	HEM	C3D-C4D-ND	-2.68	107.18	110.17
3	B	802	H4B	C2-N1-C8A	2.67	120.53	114.54
4	B	803	OUA	C04-C05-C10	2.61	119.42	118.01
2	A	801	HEM	CHA-C4D-ND	2.60	127.60	124.38
4	B	803	OUA	C05-C10-N01	-2.59	120.07	122.81
4	B	803	OUA	C26-C25-C08	2.59	126.20	121.69
2	B	801	HEM	CMD-C2D-C1D	2.58	128.96	125.04
3	B	802	H4B	C4-C4A-N5	2.53	121.25	119.12
2	B	801	HEM	CAD-CBD-CGD	-2.50	108.23	113.60
4	A	803	OUA	N02-C02-N01	2.49	120.32	118.26
2	B	801	HEM	CHA-C4D-ND	2.48	127.44	124.38
4	B	803	OUA	C24-C23-C22	2.47	119.10	116.71
2	B	801	HEM	C1B-NB-C4B	2.47	107.63	105.07
2	A	801	HEM	C4B-CHC-C1C	2.47	125.82	122.56

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	A	802	H4B	N2-C2-N3	2.45	121.07	117.25
2	A	801	HEM	C1B-NB-C4B	2.43	107.58	105.07
4	A	803	OUA	C24-C25-C08	-2.42	116.86	120.86
4	A	803	OUA	C07-C08-C25	-2.38	117.23	121.36
2	B	801	HEM	C4C-CHD-C1D	2.33	125.63	122.56
2	B	801	HEM	CAD-C3D-C4D	2.24	128.57	124.66
4	A	803	OUA	C06-C05-C04	-2.24	119.39	123.66
2	B	801	HEM	C1D-C2D-C3D	-2.21	104.63	106.96
2	B	801	HEM	C3B-C2B-C1B	2.17	108.10	106.49
2	A	801	HEM	C1D-C2D-C3D	-2.16	104.69	106.96
4	A	803	OUA	C24-C23-C22	2.13	118.77	116.71
2	B	801	HEM	CHD-C1D-ND	2.06	126.67	124.43
4	B	803	OUA	C23-C22-N21	-2.06	119.70	123.72
2	B	801	HEM	CMC-C2C-C3C	2.05	128.51	124.68
2	A	801	HEM	CMC-C2C-C3C	2.02	128.46	124.68
2	B	801	HEM	CAD-C3D-C2D	-2.00	124.14	127.88

There are no chirality outliers.

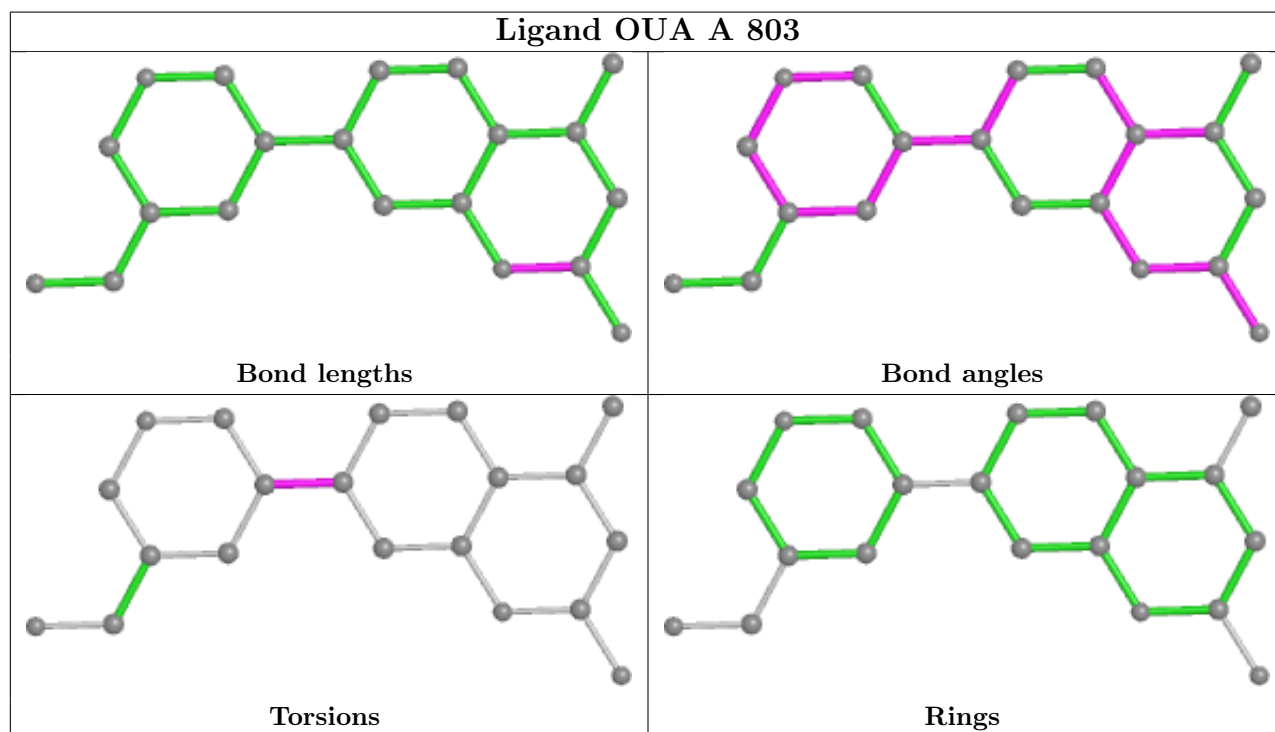
All (18) torsion outliers are listed below:

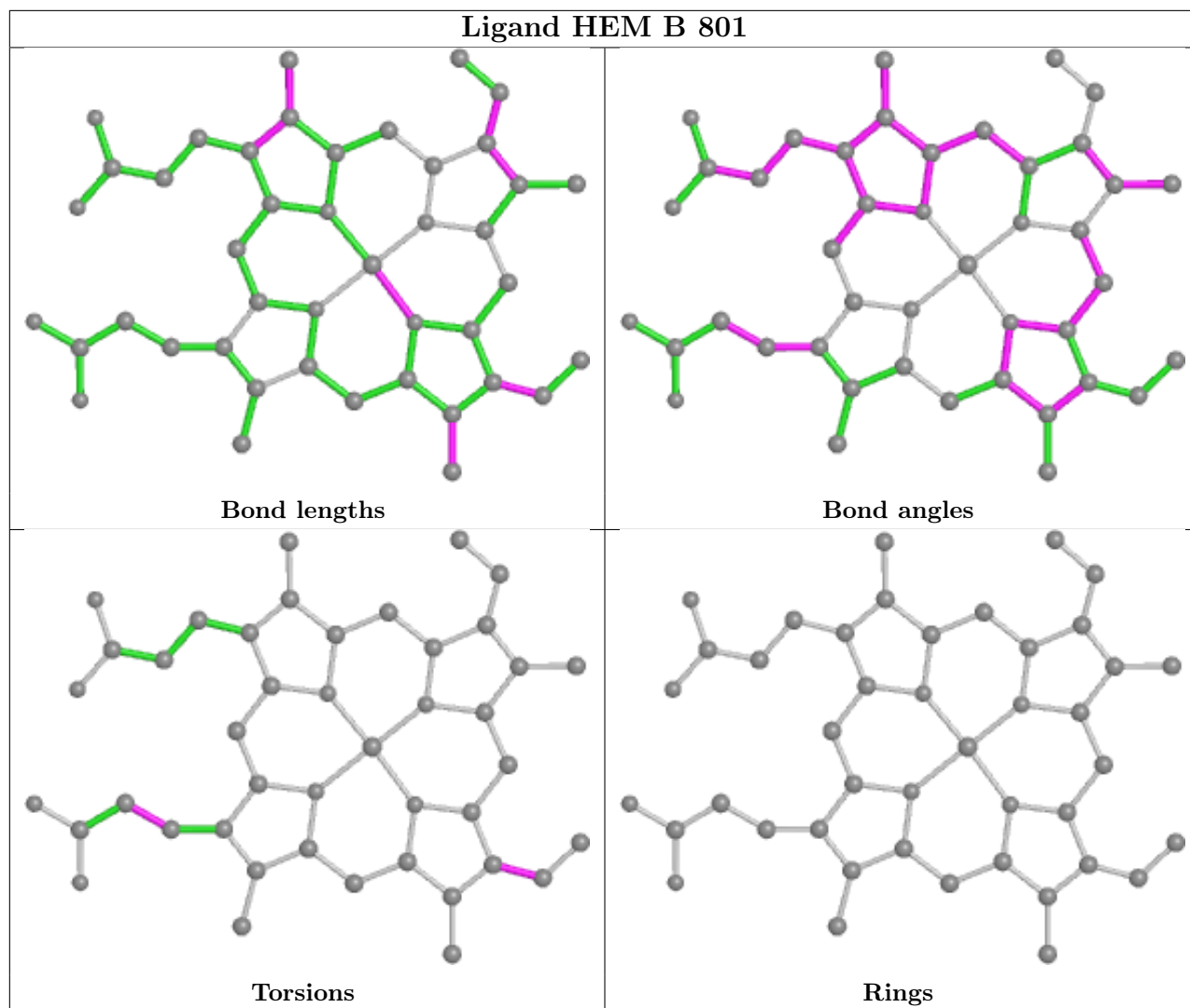
Mol	Chain	Res	Type	Atoms
2	A	801	HEM	C2A-CAA-CBA-CGA
2	A	801	HEM	C2D-C3D-CAD-CBD
4	B	803	OUA	C07-C08-C25-C24
2	A	801	HEM	C4D-C3D-CAD-CBD
4	B	803	OUA	C09-C08-C25-C24
4	B	803	OUA	C09-C08-C25-C26
4	B	803	OUA	C07-C08-C25-C26
4	A	803	OUA	C07-C08-C25-C24
2	B	801	HEM	C2A-CAA-CBA-CGA
4	A	803	OUA	C09-C08-C25-C24
2	A	801	HEM	C4B-C3B-CAB-CBB
2	B	801	HEM	C4B-C3B-CAB-CBB
4	A	803	OUA	C09-C08-C25-C26
4	A	803	OUA	C07-C08-C25-C26
2	A	801	HEM	C1A-C2A-CAA-CBA
2	A	801	HEM	C3A-C2A-CAA-CBA
2	A	801	HEM	CAD-CBD-CGD-O1D
2	A	801	HEM	CAD-CBD-CGD-O2D

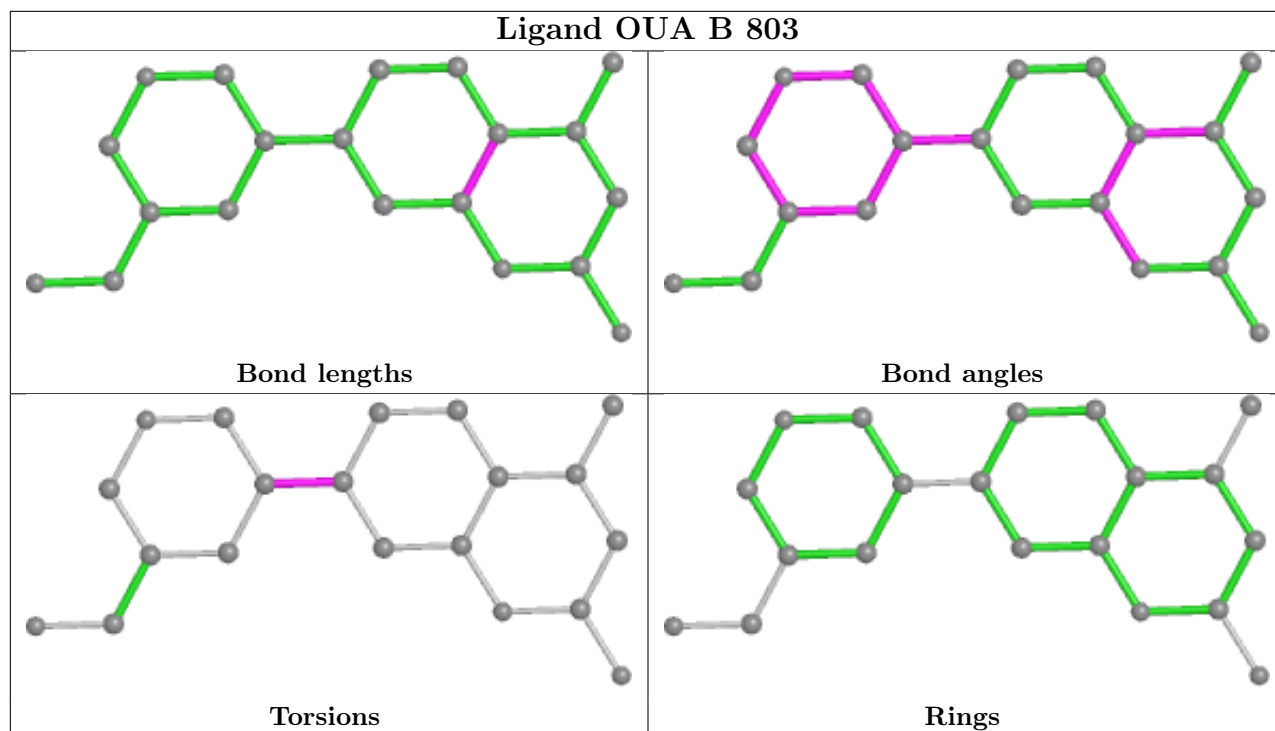
There are no ring outliers.

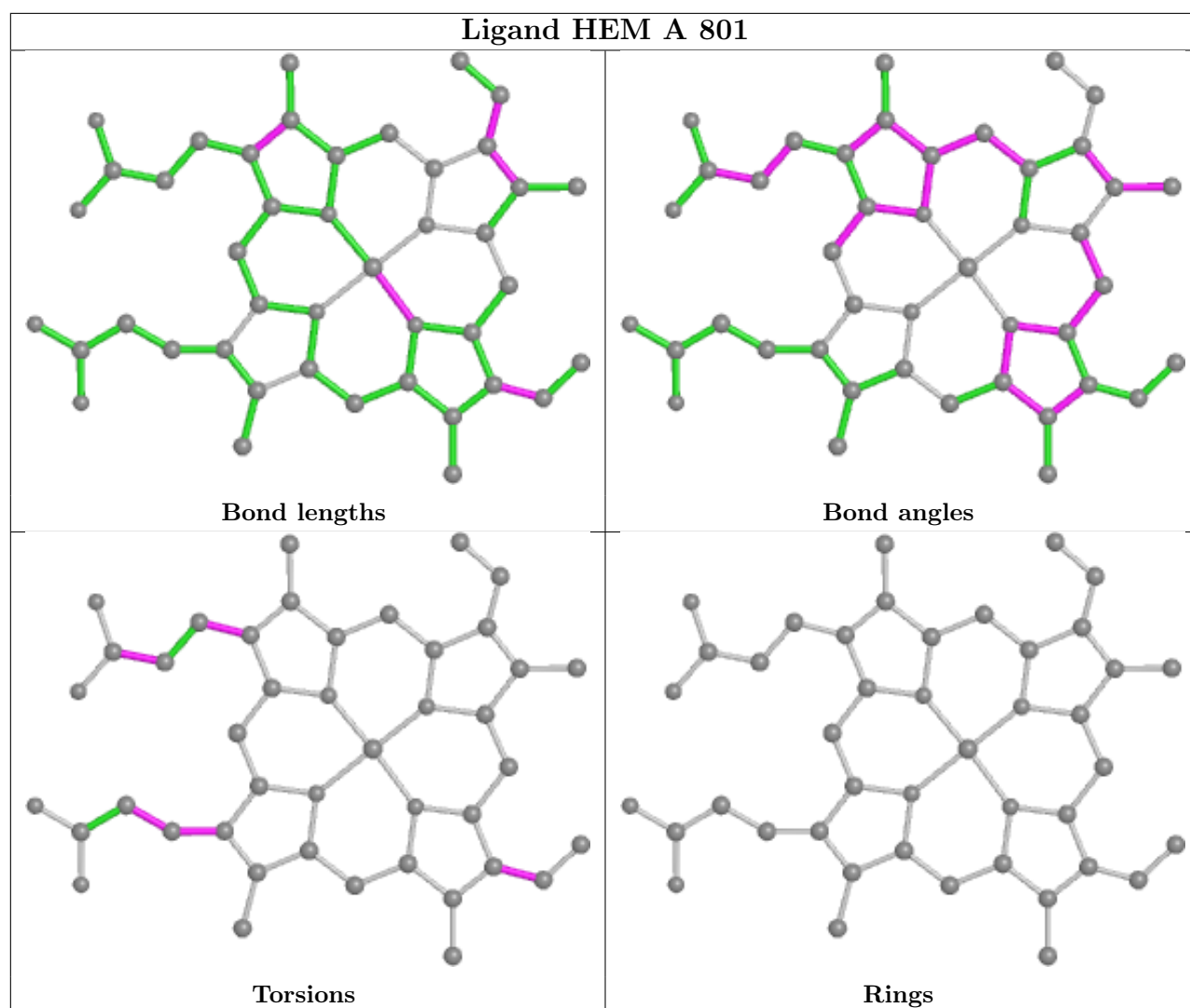
No monomer is involved in short contacts.

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 [i](#)

5.1 Protein, DNA and RNA chains [i](#)

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

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

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

5.3 Carbohydrates [i](#)

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

5.4 Ligands [i](#)

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

5.5 Other polymers [i](#)

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