



wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 1CYC
Title : THE CRYSTAL STRUCTURE OF BONITO (KATSUO) FERROCYTOCHROME C AT 2.3 ANGSTROMS RESOLUTION. II. STRUCTURE AND FUNCTION
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Deposited on : 1976-08-01
Resolution : 2.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary 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 following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtrriage (Phenix) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
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) : 2.17.1

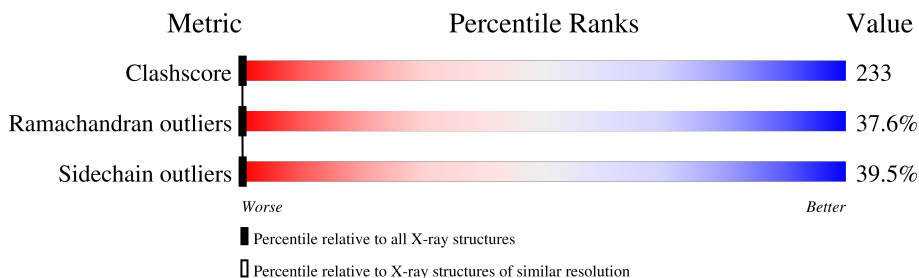
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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))
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)

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 $\leq 5\%$

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	103	32% 65%
1	B	103	32% 65%

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	HEC	A	104	-	-	X	-
2	HEC	B	104	-	-	X	-

2 Entry composition [i](#)

There are 3 unique types of molecules in this entry. The entry contains 1677 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 FERROCYTOCHROME C.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	103	Total 795	C 505	N 139	O 147	S 4	0	0	0
1	B	103	Total 795	C 505	N 139	O 147	S 4	0	0	0

- Molecule 2 is HEME C (three-letter code: HEC) (formula: $C_{34}H_{34}FeN_4O_4$).



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 water.

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

3 Residue-property plots

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. 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. 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.

Note EDS was not executed.

- Molecule 1: FERROCYTOCHROME C

Chain A:  32% 65%



- Molecule 1: FERROCYTOCHROME C

Chain B:  32% 65%



4 Data and refinement statistics

Xtrriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	57.68Å 84.58Å 37.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	(Not available) – 2.30	Depositor
% Data completeness (in resolution range)	(Not available) ((Not available)-2.30)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	unknown	Depositor
R, R_{free}	(Not available) , (Not available)	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	1677	wwPDB-VP
Average B, all atoms (Å ²)	0.0	wwPDB-VP

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: HEC

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	7.94	366/811 (45.1%)	6.18	322/1084 (29.7%)
1	B	7.94	366/811 (45.1%)	6.18	322/1084 (29.7%)
All	All	7.94	732/1622 (45.1%)	6.18	644/2168 (29.7%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	2	35
1	B	2	35
All	All	4	70

The worst 5 of 732 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	3	VAL	CA-CB	60.22	2.81	1.54
1	B	3	VAL	CA-CB	60.22	2.81	1.54
1	A	44	GLU	CG-CD	37.82	2.08	1.51
1	B	44	GLU	CG-CD	37.82	2.08	1.51
1	A	46	TYR	CG-CD1	-37.48	0.90	1.39

The worst 5 of 644 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	3	VAL	CA-CB-CG2	-38.56	53.07	110.90
1	B	3	VAL	CA-CB-CG2	-38.56	53.07	110.90
1	A	3	VAL	CA-CB-CG1	-33.90	60.05	110.90
1	B	3	VAL	CA-CB-CG1	-33.90	60.05	110.90

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	18	HIS	CG-ND1-CE1	33.74	155.43	108.20

All (4) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	40	THR	CB
1	A	78	THR	CB
1	B	40	THR	CB
1	B	78	THR	CB

5 of 70 planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	10	PHE	Mainchain
1	A	12	GLN	Sidechain
1	A	16	GLN	Sidechain
1	A	2	ASP	Mainchain
1	A	5	LYS	Mainchain

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	795	0	773	388	47
1	B	795	0	773	388	47
2	A	43	0	27	23	0
2	B	43	0	27	23	0
3	A	1	0	0	0	0
All	All	1677	0	1600	761	54

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 233.

The worst 5 of 761 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:66:GLU:HB3	1:A:74:TYR:CE2	1.24	1.67
1:B:66:GLU:HB3	1:B:74:TYR:CE2	1.24	1.67
1:A:26:HIS:CB	1:A:26:HIS:CG	1.78	1.64
1:B:26:HIS:CG	1:B:26:HIS:CB	1.78	1.64
1:A:49:THR:CA	1:A:49:THR:CB	1.74	1.63

The worst 5 of 54 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)
1:A:88:LYS:CE	1:B:88:LYS:CD[4_455]	0.41	1.79
1:A:88:LYS:CD	1:B:88:LYS:CE[4_455]	0.41	1.79
1:A:16:GLN:OE1	1:A:47:SER:OG[2_454]	0.69	1.51
1:B:16:GLN:OE1	1:B:47:SER:OG[2_554]	0.69	1.51
1:A:103:SER:CB	1:B:86:LYS:CE[1_455]	0.77	1.43

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	101/103 (98%)	42 (42%)	21 (21%)	38 (38%)	0 0
1	B	101/103 (98%)	42 (42%)	21 (21%)	38 (38%)	0 0
All	All	202/206 (98%)	84 (42%)	42 (21%)	76 (38%)	0 0

5 of 76 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	15	ALA
1	A	18	HIS
1	A	20	VAL
1	A	24	GLY
1	A	25	LYS

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	81/83 (98%)	49 (60%)	32 (40%)	0	0
1	B	81/83 (98%)	49 (60%)	32 (40%)	0	0
All	All	162/166 (98%)	98 (60%)	64 (40%)	0	0

5 of 64 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	B	81	ILE
1	B	86	LYS
1	A	80	MET
1	A	78	THR
1	B	88	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 12 such sidechains are listed below:

Mol	Chain	Res	Type
1	B	16	GLN
1	B	31	ASN
1	B	70	ASN
1	B	42	GLN
1	A	42	GLN

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](#)

2 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	HEC	B	104	1	26,50,50	7.07	19 (73%)	18,82,82	5.79	10 (55%)
2	HEC	A	104	1	26,50,50	7.07	19 (73%)	18,82,82	5.79	10 (55%)

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	HEC	B	104	1	-	4/6/54/54	-
2	HEC	A	104	1	-	4/6/54/54	-

The worst 5 of 38 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	104	HEC	C3C-C2C	-18.47	1.21	1.40
2	B	104	HEC	C3C-C2C	-18.47	1.21	1.40
2	A	104	HEC	C3B-C4B	-14.84	1.16	1.43
2	B	104	HEC	C3B-C4B	-14.84	1.16	1.43
2	A	104	HEC	CAA-C2A	-14.70	1.26	1.52

The worst 5 of 20 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	104	HEC	C4C-C3C-C2C	16.56	124.23	106.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	104	HEC	C4C-C3C-C2C	16.56	124.23	106.35
2	A	104	HEC	CAD-CBD-CGD	8.82	127.47	112.67
2	B	104	HEC	CAD-CBD-CGD	8.82	127.47	112.67
2	A	104	HEC	C3C-C4C-NC	-8.03	95.79	110.94

There are no chirality outliers.

5 of 8 torsion outliers are listed below:

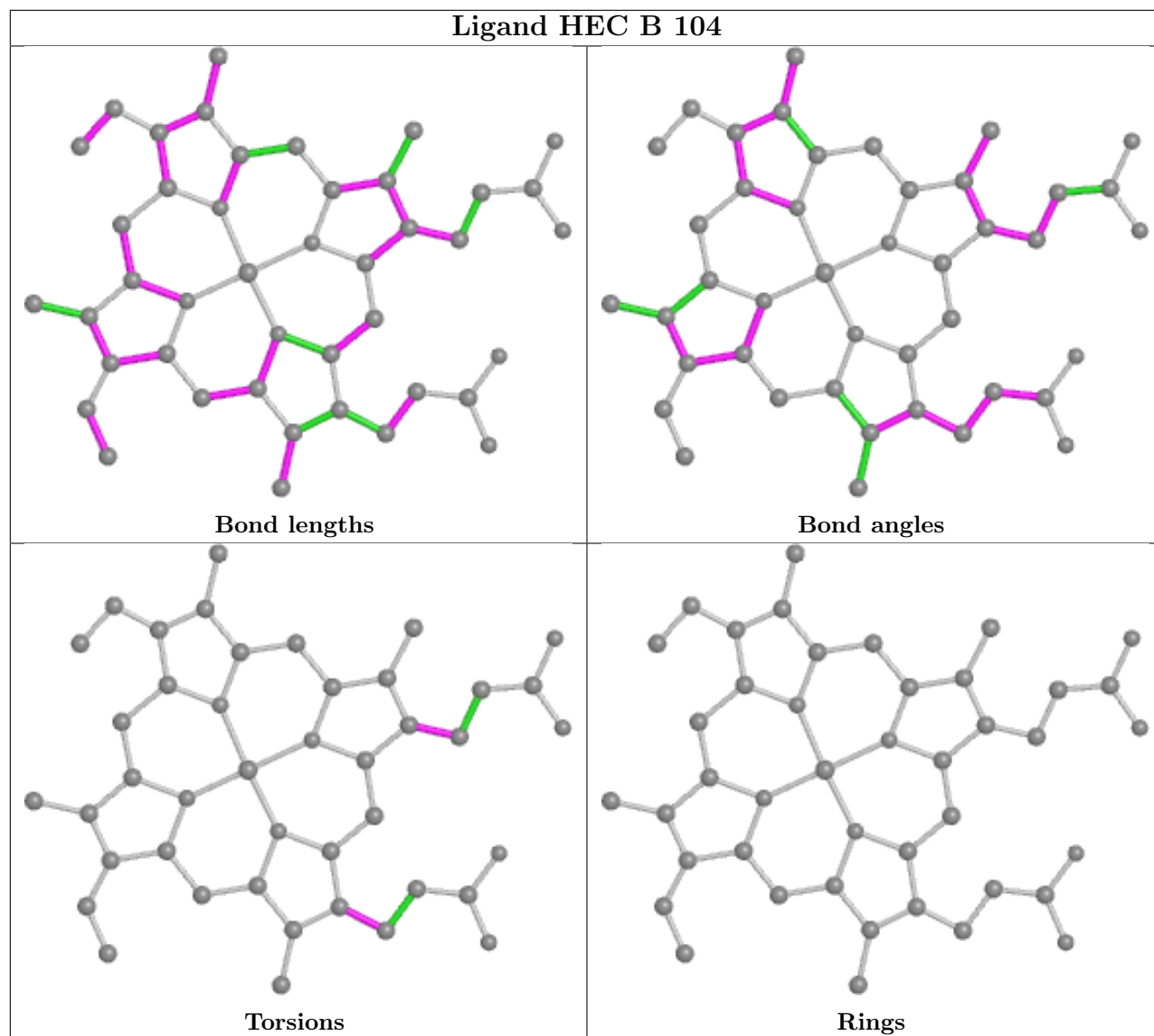
Mol	Chain	Res	Type	Atoms
2	A	104	HEC	C1A-C2A-CAA-CBA
2	A	104	HEC	C3A-C2A-CAA-CBA
2	A	104	HEC	C2D-C3D-CAD-CBD
2	A	104	HEC	C4D-C3D-CAD-CBD
2	B	104	HEC	C1A-C2A-CAA-CBA

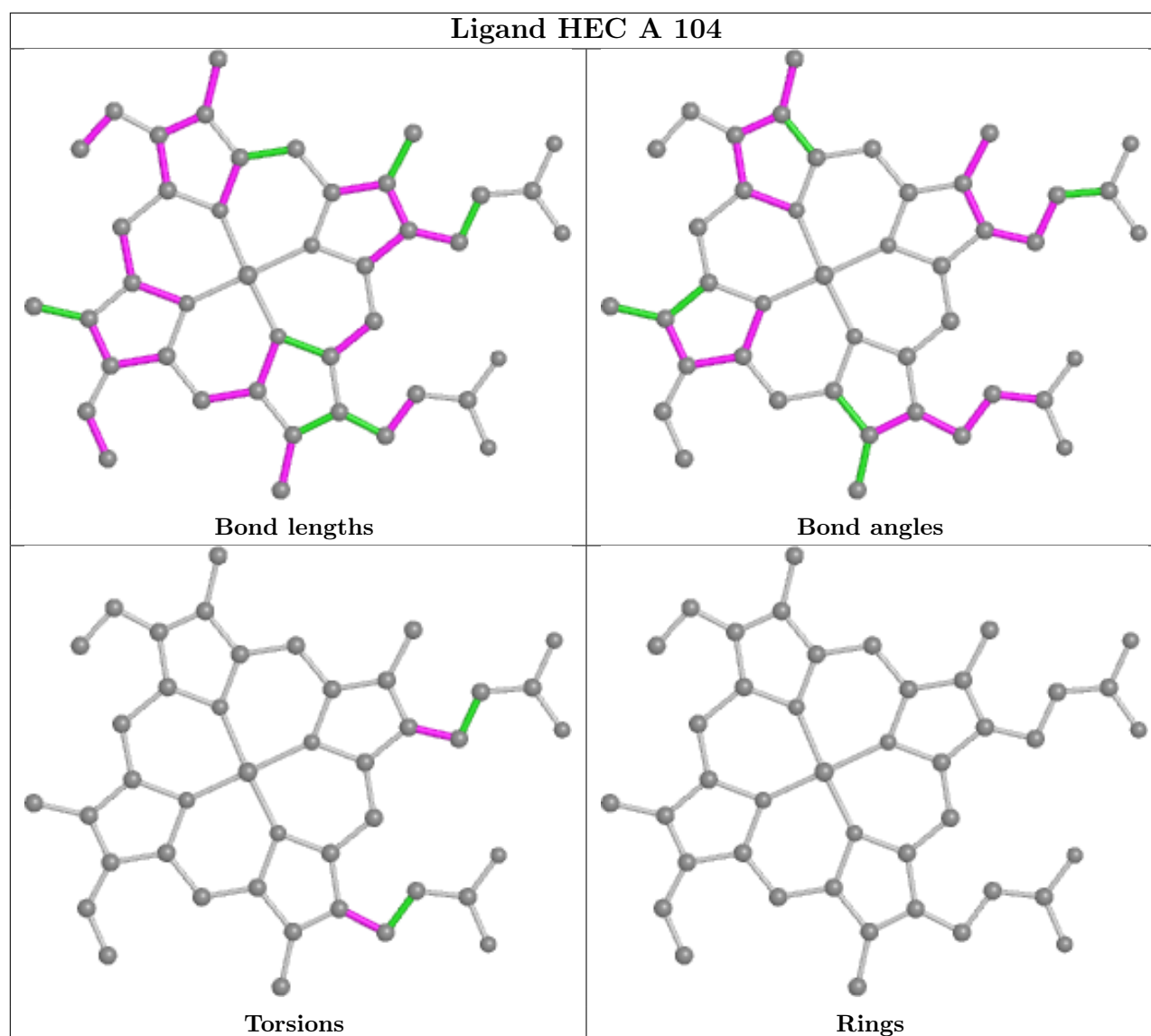
There are no ring outliers.

2 monomers are involved in 46 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	B	104	HEC	23	0
2	A	104	HEC	23	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 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.





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

There are no such residues in this entry.

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

The following chains have linkage breaks:

Mol	Chain	Number of breaks
1	A	29
1	B	29

The worst 5 of 58 chain breaks are listed below:

Model	Chain	Residue-1	Atom-1	Residue-2	Atom-2	Distance (Å)
1	A	26:HIS	C	27:LYS	N	1.20
1	A	31:ASN	C	32:LEU	N	1.20
1	A	50:ASP	C	51:ALA	N	1.20
1	A	80:MET	C	81:ILE	N	1.20
1	B	26:HIS	C	27:LYS	N	1.20

6 Fit of model and data

6.1 Protein, DNA and RNA chains

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.