



# Full wwPDB X-ray Structure Validation Report ⓘ

Oct 5, 2023 – 03:11 AM EDT

PDB ID : 6UZI  
Title : Crystal structure of Dihydrolipoyl dehydrogenase from Elizabethkingia anophelis NUHP1  
Authors : Seattle Structural Genomics Center for Infectious Disease (SSGCID)  
Deposited on : 2019-11-15  
Resolution : 2.80 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](mailto: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>.

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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 2.80 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 14122 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 Dihydrolipoyl dehydrogenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	467	Total 3478	C 2186	N 594	O 681	S 17	0	0	0
1	B	467	Total 3393	C 2132	N 578	O 666	S 17	0	0	0
1	C	470	Total 3432	C 2157	N 583	O 675	S 17	0	0	0
1	D	466	Total 3355	C 2111	N 571	O 657	S 16	0	0	0

There are 32 discrepancies between the modelled and reference sequences:

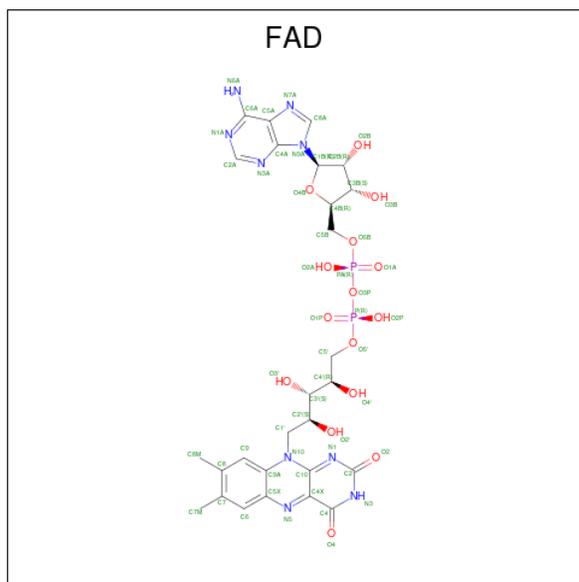
Chain	Residue	Modelled	Actual	Comment	Reference
A	-7	MET	-	expression tag	UNP A0A077ELH4
A	-6	ALA	-	expression tag	UNP A0A077ELH4
A	-5	HIS	-	expression tag	UNP A0A077ELH4
A	-4	HIS	-	expression tag	UNP A0A077ELH4
A	-3	HIS	-	expression tag	UNP A0A077ELH4
A	-2	HIS	-	expression tag	UNP A0A077ELH4
A	-1	HIS	-	expression tag	UNP A0A077ELH4
A	0	HIS	-	expression tag	UNP A0A077ELH4
B	-7	MET	-	expression tag	UNP A0A077ELH4
B	-6	ALA	-	expression tag	UNP A0A077ELH4
B	-5	HIS	-	expression tag	UNP A0A077ELH4
B	-4	HIS	-	expression tag	UNP A0A077ELH4
B	-3	HIS	-	expression tag	UNP A0A077ELH4
B	-2	HIS	-	expression tag	UNP A0A077ELH4
B	-1	HIS	-	expression tag	UNP A0A077ELH4
B	0	HIS	-	expression tag	UNP A0A077ELH4
C	-7	MET	-	expression tag	UNP A0A077ELH4
C	-6	ALA	-	expression tag	UNP A0A077ELH4
C	-5	HIS	-	expression tag	UNP A0A077ELH4
C	-4	HIS	-	expression tag	UNP A0A077ELH4
C	-3	HIS	-	expression tag	UNP A0A077ELH4

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Chain	Residue	Modelled	Actual	Comment	Reference
C	-2	HIS	-	expression tag	UNP A0A077ELH4
C	-1	HIS	-	expression tag	UNP A0A077ELH4
C	0	HIS	-	expression tag	UNP A0A077ELH4
D	-7	MET	-	expression tag	UNP A0A077ELH4
D	-6	ALA	-	expression tag	UNP A0A077ELH4
D	-5	HIS	-	expression tag	UNP A0A077ELH4
D	-4	HIS	-	expression tag	UNP A0A077ELH4
D	-3	HIS	-	expression tag	UNP A0A077ELH4
D	-2	HIS	-	expression tag	UNP A0A077ELH4
D	-1	HIS	-	expression tag	UNP A0A077ELH4
D	0	HIS	-	expression tag	UNP A0A077ELH4

- Molecule 2 is FLAVIN-ADENINE DINUCLEOTIDE (three-letter code: FAD) (formula:  $C_{27}H_{33}N_9O_{15}P_2$ ) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf		
			Total	C	N	O			P	
2	A	1	Total	53	27	9	15	2	0	0
2	B	1	Total	53	27	9	15	2	0	0
2	C	1	Total	53	27	9	15	2	0	0
2	D	1	Total	53	27	9	15	2	0	0

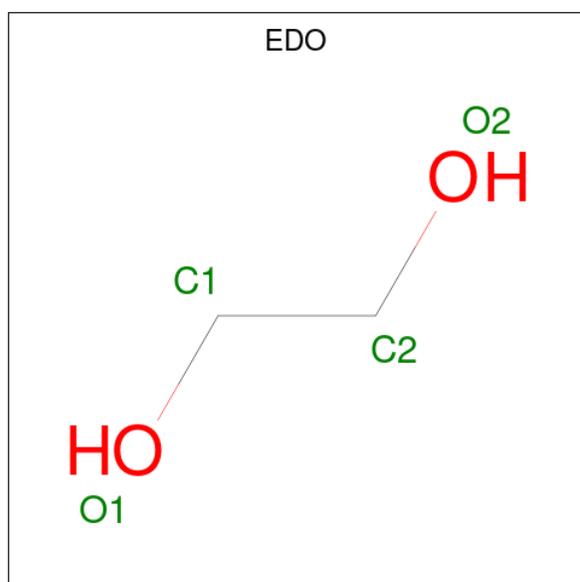
- Molecule 3 is ZINC ION (three-letter code: ZN) (formula: Zn).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	2	Total	Zn	0	0
			2	2		
3	C	2	Total	Zn	0	0
			2	2		

- Molecule 4 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

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

- Molecule 5 is 1,2-ETHANEDIOL (three-letter code: EDO) (formula: C<sub>2</sub>H<sub>6</sub>O<sub>2</sub>).



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

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

- Molecule 6 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
6	A	104	Total	O	0	0
			104	104		
6	B	61	Total	O	0	0
			61	61		
6	C	39	Total	O	0	0
			39	39		
6	D	17	Total	O	0	0
			17	17		

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, $\alpha$ , $\beta$ , $\gamma$	74.06Å 106.80Å 255.70Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	45.25 – 2.80	Depositor
% Data completeness (in resolution range)	98.9 (45.25-2.80)	Depositor
$R_{merge}$	0.09	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.89 (at 2.81Å)	Xtriage
Refinement program	PHENIX 1.17.1 -3660	Depositor
R, $R_{free}$	0.169 , 0.230	Depositor
Wilson B-factor (Å <sup>2</sup> )	54.9	Xtriage
Anisotropy	0.402	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.50$ , $\langle L^2 \rangle = 0.34$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	14122	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	58.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.44% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

<sup>2</sup>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](#)

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](#)

Of 17 ligands modelled in this entry, 7 are monoatomic - leaving 10 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
5	EDO	A	504	-	3,3,3	0.43	0	2,2,2	0.37	0
2	FAD	D	501	-	53,58,58	0.46	0	68,89,89	0.59	1 (1%)
5	EDO	A	506	-	3,3,3	0.61	0	2,2,2	0.03	0
2	FAD	B	501	-	53,58,58	0.45	0	68,89,89	0.61	2 (2%)
5	EDO	B	503	-	3,3,3	0.56	0	2,2,2	0.07	0
5	EDO	A	505	-	3,3,3	0.52	0	2,2,2	0.25	0
5	EDO	A	507	-	3,3,3	0.43	0	2,2,2	0.37	0
5	EDO	C	504	-	3,3,3	0.49	0	2,2,2	0.31	0
2	FAD	C	501	-	53,58,58	0.44	0	68,89,89	0.65	1 (1%)
2	FAD	A	501	-	53,58,58	0.45	0	68,89,89	0.61	1 (1%)

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	EDO	A	504	-	-	1/1/1/1	-
2	FAD	D	501	-	-	8/30/50/50	0/6/6/6
5	EDO	A	506	-	-	1/1/1/1	-
2	FAD	B	501	-	-	11/30/50/50	0/6/6/6
5	EDO	B	503	-	-	1/1/1/1	-
5	EDO	A	505	-	-	0/1/1/1	-
5	EDO	A	507	-	-	0/1/1/1	-
5	EDO	C	504	-	-	1/1/1/1	-
2	FAD	C	501	-	-	9/30/50/50	0/6/6/6
2	FAD	A	501	-	-	10/30/50/50	0/6/6/6

There are no bond length outliers.

All (5) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	D	501	FAD	C5A-C6A-N6A	2.32	123.88	120.35
2	A	501	FAD	C5A-C6A-N6A	2.32	123.88	120.35
2	C	501	FAD	C5A-C6A-N6A	2.32	123.87	120.35
2	B	501	FAD	C5A-C6A-N6A	2.30	123.85	120.35

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	B	501	FAD	O5'-P-O1P	2.29	118.00	109.07

There are no chirality outliers.

All (42) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	A	501	FAD	C5B-O5B-PA-O2A
2	A	501	FAD	C5B-O5B-PA-O3P
2	A	501	FAD	O4B-C4B-C5B-O5B
2	A	501	FAD	C3'-C4'-C5'-O5'
2	A	501	FAD	O4'-C4'-C5'-O5'
2	A	501	FAD	C5'-O5'-P-O1P
2	A	501	FAD	C5'-O5'-P-O2P
2	B	501	FAD	C5B-O5B-PA-O3P
2	B	501	FAD	C3'-C4'-C5'-O5'
2	B	501	FAD	O4'-C4'-C5'-O5'
2	B	501	FAD	C5'-O5'-P-O1P
2	B	501	FAD	C5'-O5'-P-O2P
2	C	501	FAD	C5B-O5B-PA-O3P
2	C	501	FAD	C3'-C4'-C5'-O5'
2	C	501	FAD	O4'-C4'-C5'-O5'
2	C	501	FAD	C5'-O5'-P-O1P
2	D	501	FAD	C5'-O5'-P-O2P
2	B	501	FAD	O4B-C4B-C5B-O5B
2	C	501	FAD	O4B-C4B-C5B-O5B
2	B	501	FAD	C3B-C4B-C5B-O5B
2	C	501	FAD	C3B-C4B-C5B-O5B
5	A	504	EDO	O1-C1-C2-O2
5	A	506	EDO	O1-C1-C2-O2
2	D	501	FAD	O4B-C4B-C5B-O5B
2	D	501	FAD	C3B-C4B-C5B-O5B
2	B	501	FAD	PA-O3P-P-O5'
2	C	501	FAD	PA-O3P-P-O5'
2	D	501	FAD	PA-O3P-P-O5'
2	D	501	FAD	C5'-O5'-P-O3P
2	A	501	FAD	C5B-O5B-PA-O1A
2	B	501	FAD	C5B-O5B-PA-O1A
2	B	501	FAD	C5B-O5B-PA-O2A
2	C	501	FAD	C5B-O5B-PA-O1A
2	C	501	FAD	C5B-O5B-PA-O2A
5	C	504	EDO	O1-C1-C2-O2
5	B	503	EDO	O1-C1-C2-O2

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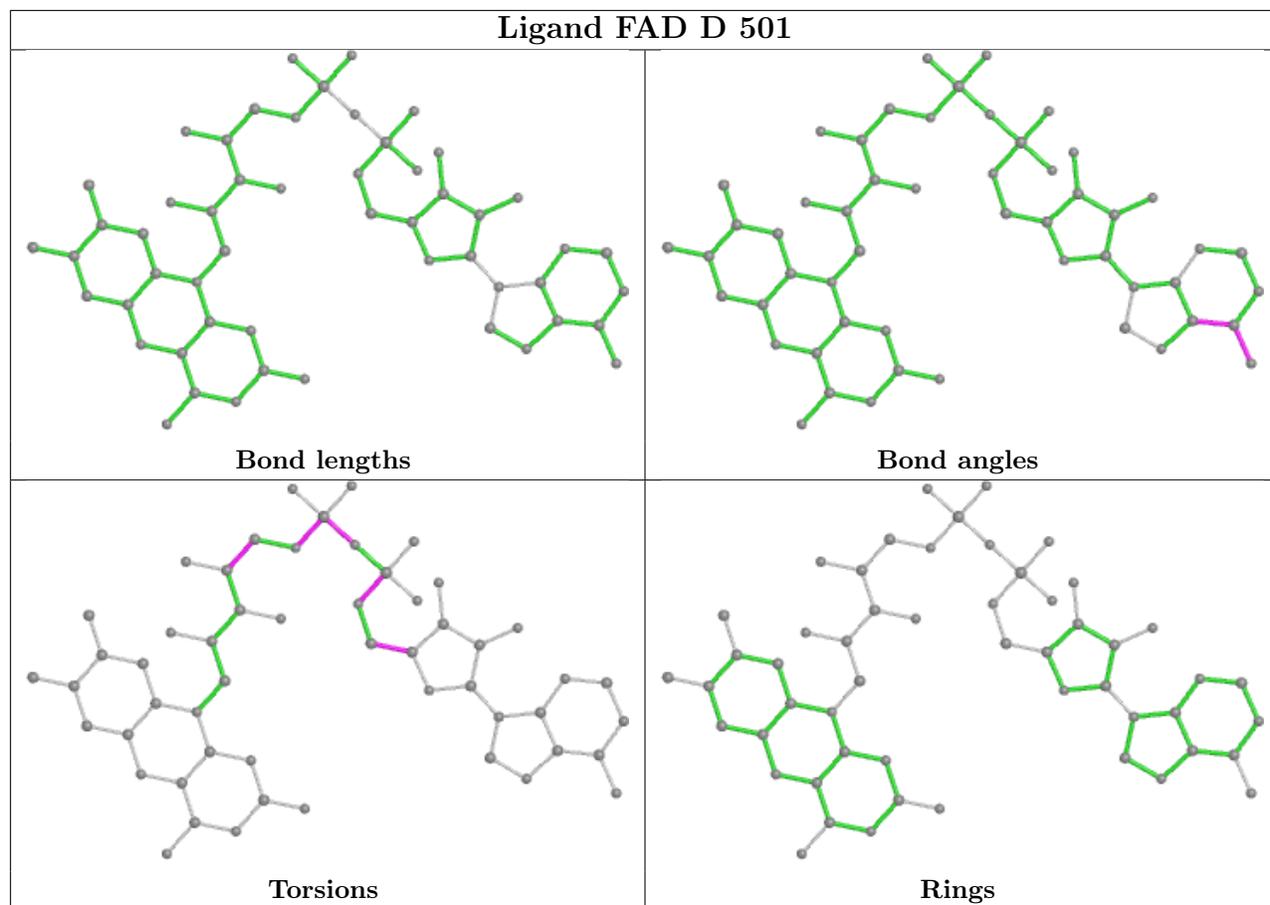
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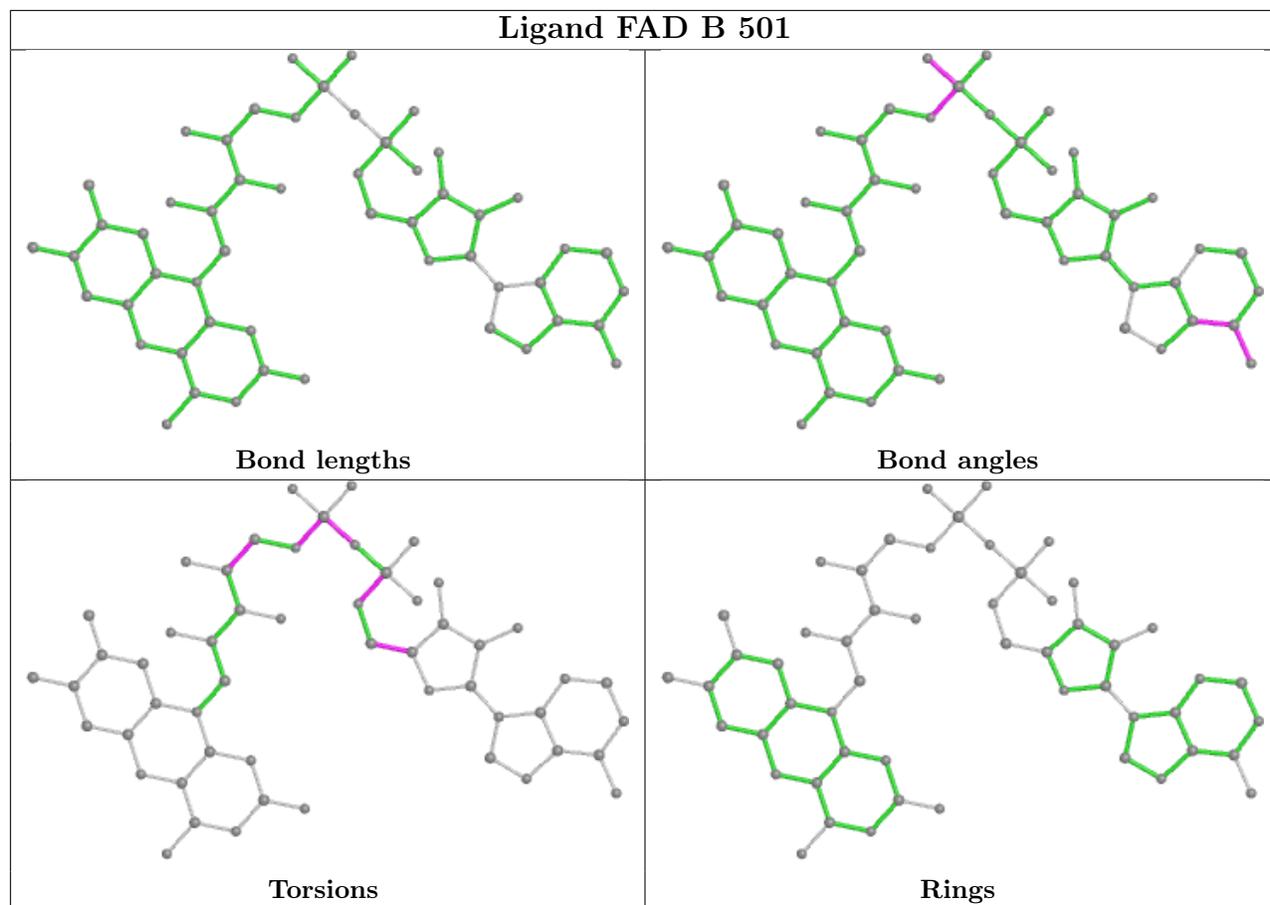
Mol	Chain	Res	Type	Atoms
2	A	501	FAD	PA-O3P-P-O5'
2	D	501	FAD	O4'-C4'-C5'-O5'
2	A	501	FAD	C5'-O5'-P-O3P
2	B	501	FAD	C5'-O5'-P-O3P
2	D	501	FAD	C5B-O5B-PA-O1A
2	D	501	FAD	C5'-O5'-P-O1P

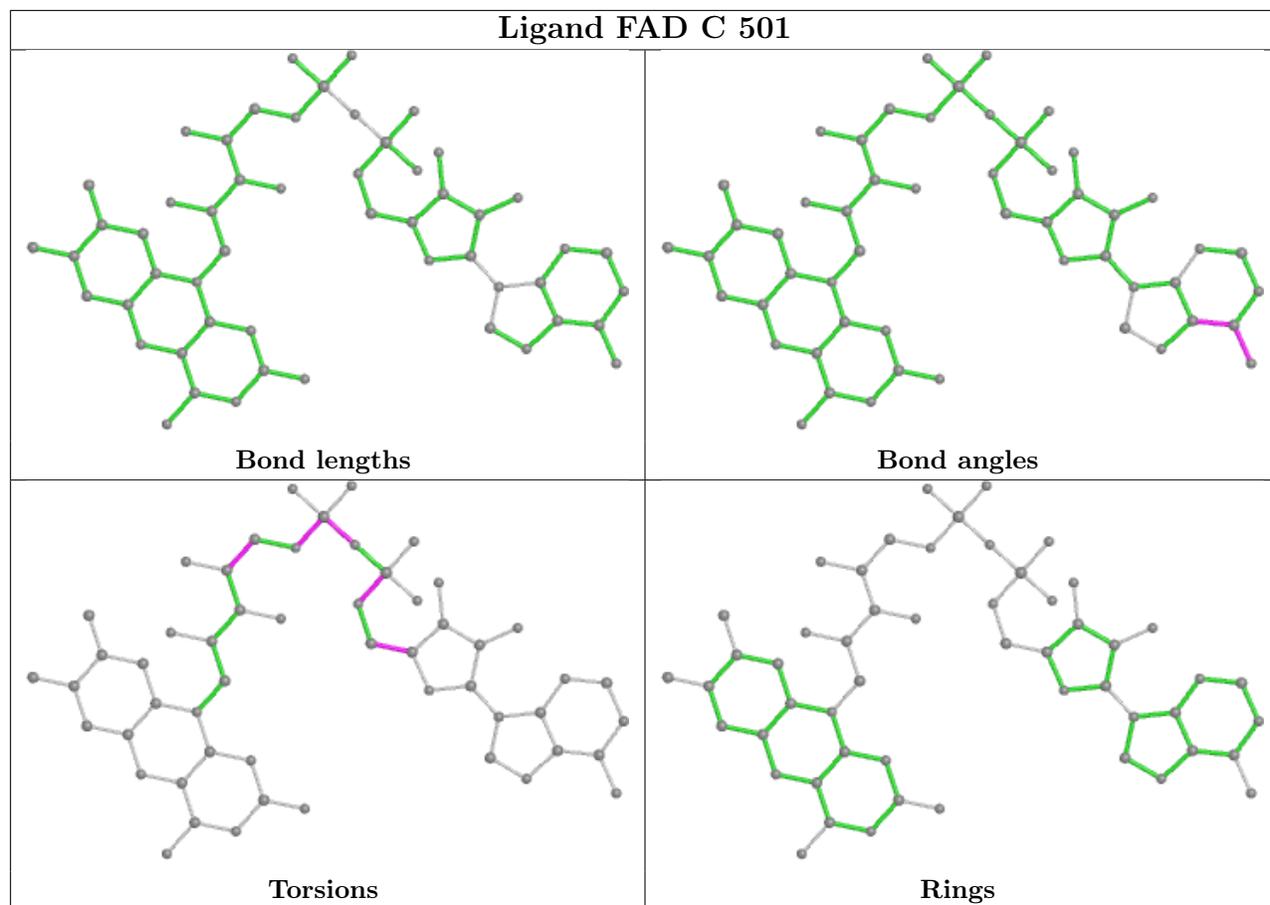
There are no ring outliers.

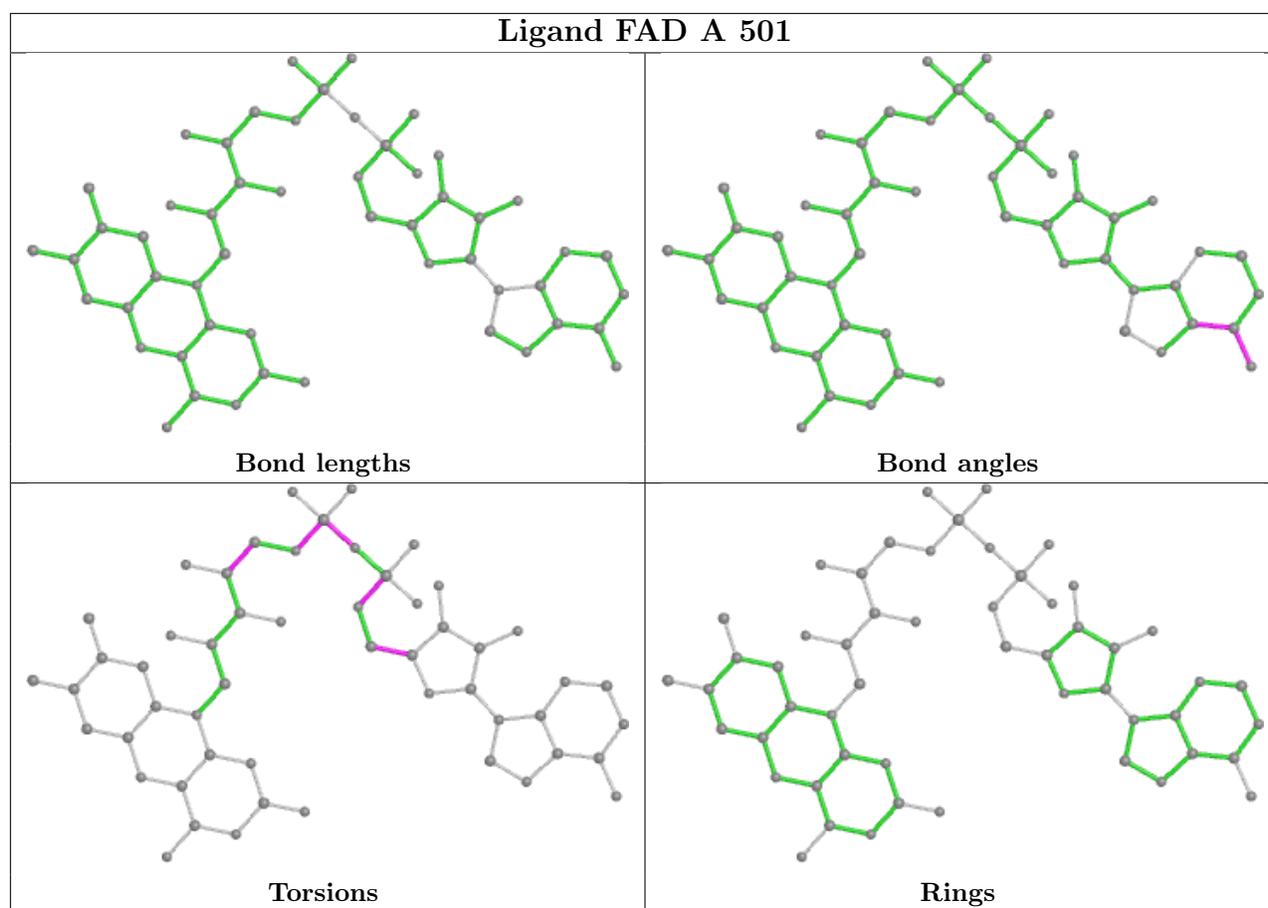
No monomer is involved in short contacts.

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.