

# wwPDB X-ray Structure Validation Summary Report (i)

#### Oct 3, 2023 – 12:09 AM EDT

PDB ID	:	6OC0
Title	:	Crystal structure of human DHODH with OSU-03012
Authors	:	Durst, M.A.; Lavie, A.
Deposited on		
Resolution	:	1.40  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 (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 (1)) were used in the production of this report:

MolProbity	:	FAILED
Mogul	:	1.8.5 (274361), CSD as541be (2020)
Xtriage (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 (i)

The following experimental techniques were used to determine the structure:  $X\hbox{-}RAY\,DIFFRACTION$ 

The reported resolution of this entry is 1.40 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.



#### 6OC0

# 2 Entry composition (i)

There are 6 unique types of molecules in this entry. The entry contains 3168 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.

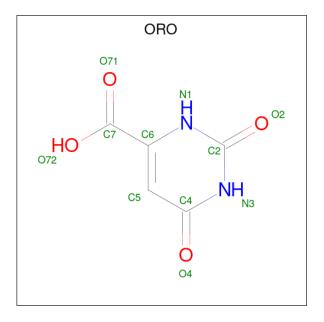
• Molecule 1 is a protein called Dihydroorotate dehydrogenase (quinone), mitochondrial.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	А	365	Total	C 1784	N 594	0	S 4	0	9	0
			2845	1784	524	533	4			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	27	HIS	-	expression tag	UNP Q02127
А	28	MET	-	expression tag	UNP Q02127

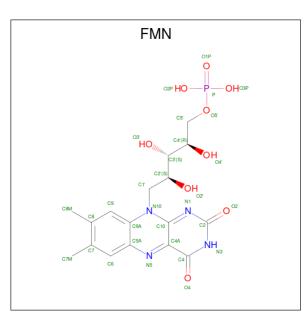
• Molecule 2 is OROTIC ACID (three-letter code: ORO) (formula:  $C_5H_4N_2O_4$ ).



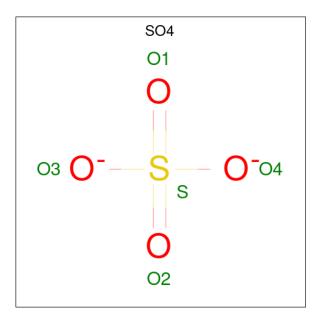
Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	А	1	Total 11	С 5	N 2	0 4	0	0

• Molecule 3 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula:  $C_{17}H_{21}N_4O_9P$ ).





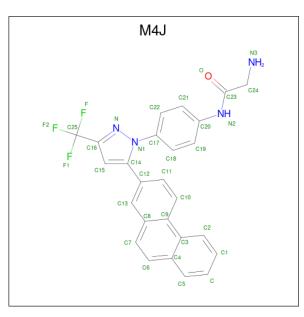
Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	А	1	Total	С	Ν	0	Р	0	0
5			31	17	4	9	1	0	



Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
4	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0
4	А	1	$\begin{array}{ccc} \text{Total} & \text{O} & \text{S} \\ 5 & 4 & 1 \end{array}$	0	0



• Molecule 5 is N-{4-[5-(phenanthren-2-yl)-3-(trifluoromethyl)-1H-pyrazol-1-yl]phenyl}glycin amide (three-letter code: M4J) (formula:  $C_{26}H_{19}F_3N_4O$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
Б	Λ	1	Total	С	F	Ν	Ο	0	0
5	А	1	34	26	3	4	1	0	

• Molecule 6 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
6	А	232	Total O   232 232	0	0

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



# 3 Data and refinement statistics (i)

Property	Value	Source		
Space group	P 32 2 1	Depositor		
Cell constants	91.01Å 91.01Å 123.07Å	Depositor		
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $120.00^{\circ}$	Depositor		
Resolution (Å)	48.55 - 1.40	Depositor		
% Data completeness	97.7 (48.55-1.40)	Depositor		
(in resolution range)		Depositor		
$R_{merge}$	(Not available)	Depositor		
R <sub>sym</sub>	(Not available)	Depositor		
$< I/\sigma(I) > 1$	$1.18 (at 1.41 \text{\AA})$	Xtriage		
Refinement program	REFMAC 5.8.0238	Depositor		
$R, R_{free}$	0.151 , $0.184$	Depositor		
Wilson B-factor $(Å^2)$	13.9	Xtriage		
Anisotropy	0.922	Xtriage		
L-test for twinning <sup>2</sup>	$<  L  > = 0.48, < L^2 > = 0.31$	Xtriage		
Estimated twinning fraction	0.028 for -h,-k,l	Xtriage		
Total number of atoms	3168	wwPDB-VP		
Average B, all atoms $(Å^2)$	24.0	wwPDB-VP		

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

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.18% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;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.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

# 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)

6 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



Mol	Trune	ype Chain Res Link		Tinle	Bo	ond leng	ths	Bond angles		
	Type	Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z >2
2	ORO	А	401	-	9,11,11	1.36	1 (11%)	8,15,15	<mark>3.32</mark>	3 (37%)
4	SO4	А	405	-	4,4,4	0.39	0	6,6,6	0.09	0
4	SO4	А	403	-	4,4,4	0.32	0	6,6,6	0.29	0
3	FMN	А	402	-	33,33,33	1.28	2 (6%)	48,50,50	1.33	7 (14%)
5	M4J	А	406	-	36,38,38	1.18	3 (8%)	48,56,56	1.39	5 (10%)
4	SO4	А	404	-	4,4,4	0.28	0	6,6,6	0.09	0

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).

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	ORO	А	401	-	-	4/4/4/4	0/1/1/1
5	M4J	А	406	-	-	0/20/20/20	0/5/5/5
3	FMN	A	402	-	-	5/18/18/18	0/3/3/3

The worst 5 of 6 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
3	А	402	FMN	C4A-N5	3.79	1.38	1.30
5	А	406	M4J	C17-N1	-3.72	1.32	1.44
3	А	402	FMN	C10-N1	2.88	1.39	1.33
5	А	406	M4J	C9-C8	2.44	1.47	1.42
2	А	401	ORO	C4-N3	2.43	1.37	1.33

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$\mathbf{Ideal}(^{o})$
2	А	401	ORO	C5-C4-N3	-7.33	115.52	124.08
2	А	401	ORO	C6-C5-C4	5.20	120.09	116.73
5	А	406	M4J	C15-C14-C12	-4.93	120.26	128.10
3	А	402	FMN	O2-C2-N1	-3.75	115.61	121.83
5	А	406	M4J	C16-N-N1	-3.42	101.69	105.66

There are no chirality outliers.

5 of 9 torsion outliers are listed below:



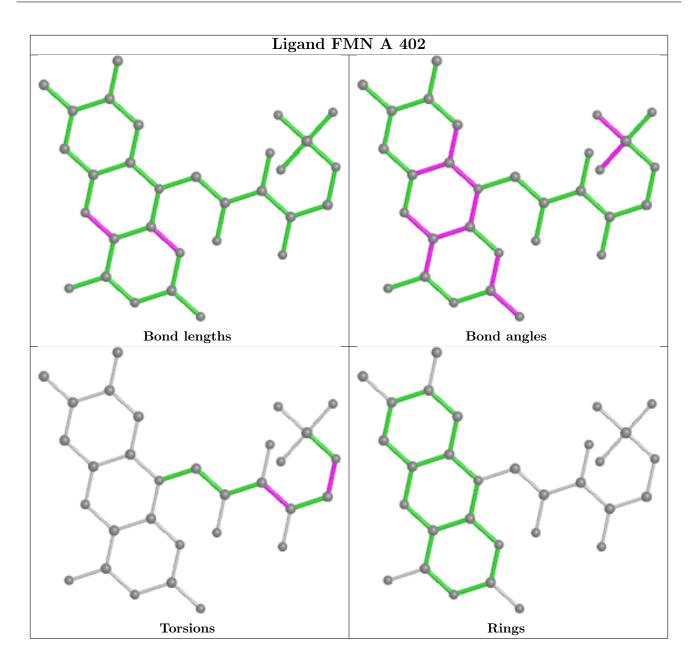
Mol	Chain	Res	Type	Atoms
2	А	401	ORO	C5-C6-C7-O71
2	А	401	ORO	C5-C6-C7-O72
2	А	401	ORO	N1-C6-C7-O72
3	А	402	FMN	C2'-C3'-C4'-O4'
2	А	401	ORO	N1-C6-C7-O71

There are no ring outliers.

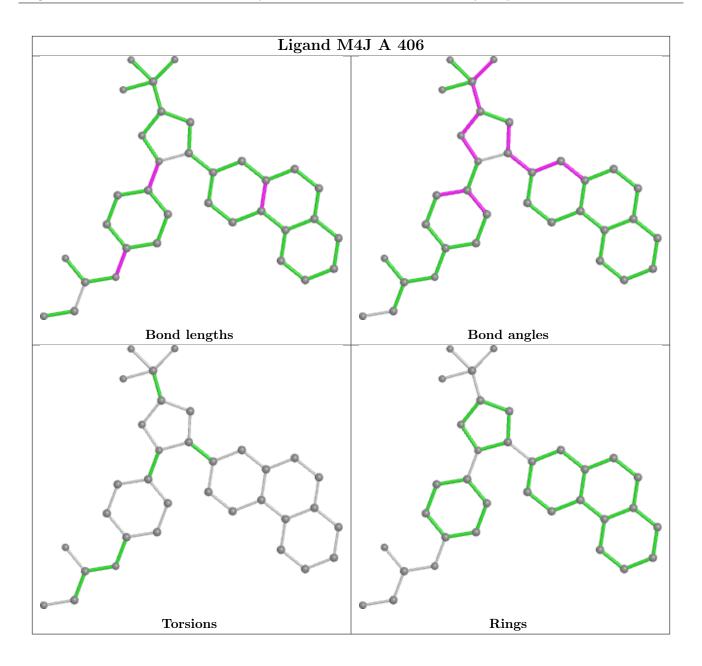
No monomer is involved in short contacts.

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 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 sufficient must be 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.

