



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:18 PM GMT

PDB ID : 1EY2  
Title : HUMAN HOMOGENITISATE DIOXYGENASE WITH FE(II)  
Authors : Timm, D.E.; Titus, G.P.; Penalva, M.A.; Mueller, H.A.; de Cordoba, S.M.  
Deposited on : 2000-05-05  
Resolution : 2.30 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  
<http://wwpdb.org/validation/2016/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 : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : 1.9-1692  
EDS : rb-20026688  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Refmac : 5.8.0135  
CCP4 : 6.5.0  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

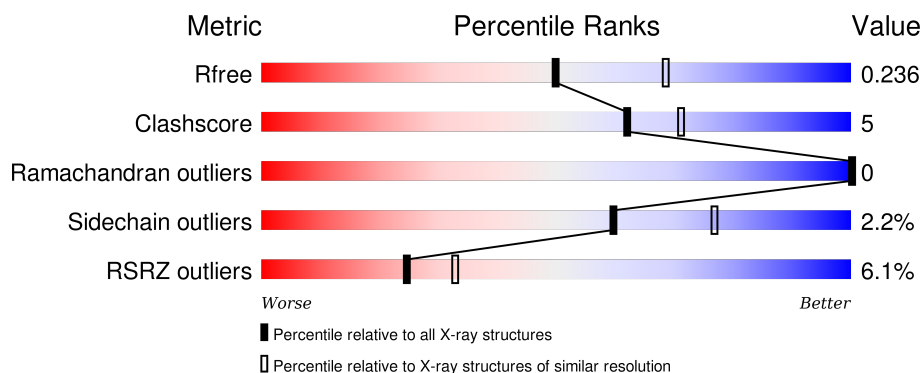
# 1 Overall quality at a glance

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))
$R_{free}$	91344	3852 (2.30-2.30)
Clashscore	102246	4452 (2.30-2.30)
Ramachandran outliers	100387	4410 (2.30-2.30)
Sidechain outliers	100360	4409 (2.30-2.30)
RSRZ outliers	91569	3857 (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 on the lower bar indicate the fraction of residues that contain outliers for  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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  $\leq 5\%$ . The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density. The numeric value is given above the bar.

Mol	Chain	Length	Quality of chain
1	A	471	<div> <div>5%</div> <div>79%</div> <div>8%</div> <div>11%</div> </div>

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3544 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 HOMOGENITISATE 1,2-DIOXYGENASE.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	419	Total	C	N	O	S	Se	0	0	0
			3311	2123	566	603	10	9			

There are 36 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-25	MSE	-	EXPRESSION TAG	UNP Q93099
A	-24	GLY	-	EXPRESSION TAG	UNP Q93099
A	-23	HIS	-	EXPRESSION TAG	UNP Q93099
A	-22	HIS	-	EXPRESSION TAG	UNP Q93099
A	-21	HIS	-	EXPRESSION TAG	UNP Q93099
A	-20	HIS	-	EXPRESSION TAG	UNP Q93099
A	-19	HIS	-	EXPRESSION TAG	UNP Q93099
A	-18	HIS	-	EXPRESSION TAG	UNP Q93099
A	-17	HIS	-	EXPRESSION TAG	UNP Q93099
A	-16	HIS	-	EXPRESSION TAG	UNP Q93099
A	-15	HIS	-	EXPRESSION TAG	UNP Q93099
A	-14	HIS	-	EXPRESSION TAG	UNP Q93099
A	-13	SER	-	EXPRESSION TAG	UNP Q93099
A	-12	SER	-	EXPRESSION TAG	UNP Q93099
A	-11	GLY	-	EXPRESSION TAG	UNP Q93099
A	-10	HIS	-	EXPRESSION TAG	UNP Q93099
A	-9	ILE	-	EXPRESSION TAG	UNP Q93099
A	-8	ASP	-	EXPRESSION TAG	UNP Q93099
A	-7	ASP	-	EXPRESSION TAG	UNP Q93099
A	-6	ASP	-	EXPRESSION TAG	UNP Q93099
A	-5	ASP	-	EXPRESSION TAG	UNP Q93099
A	-4	LYS	-	EXPRESSION TAG	UNP Q93099
A	-3	HIS	-	EXPRESSION TAG	UNP Q93099
A	-2	MSE	-	EXPRESSION TAG	UNP Q93099
A	-1	GLY	-	EXPRESSION TAG	UNP Q93099
A	0	SER	-	EXPRESSION TAG	UNP Q93099
A	1	MSE	MET	ENGINEERED	UNP Q93099

*Continued on next page...*

*Continued from previous page...*

Chain	Residue	Modelled	Actual	Comment	Reference
A	142	MSE	MET	ENGINEERED	UNP Q93099
A	172	MSE	MET	ENGINEERED	UNP Q93099
A	186	MSE	MET	ENGINEERED	UNP Q93099
A	283	MSE	MET	ENGINEERED	UNP Q93099
A	339	MSE	MET	ENGINEERED	UNP Q93099
A	343	MSE	MET	ENGINEERED	UNP Q93099
A	368	MSE	MET	ENGINEERED	UNP Q93099
A	396	MSE	MET	ENGINEERED	UNP Q93099
A	399	MSE	MET	ENGINEERED	UNP Q93099

- Molecule 2 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	1	Total Fe 1 1	0	0

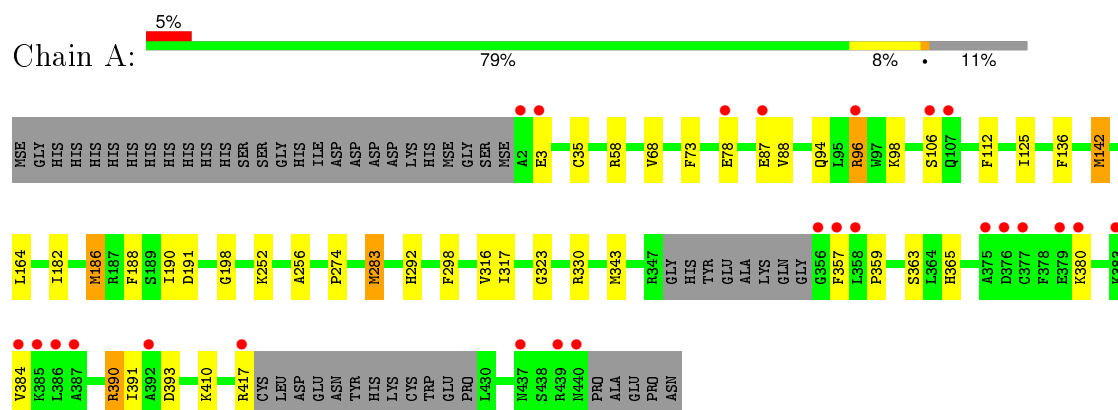
- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	232	Total O 232 232	0	0

### 3 Residue-property plots [i](#)

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors displayed in the second graphic. The second graphic shows the sequence view annotated by issues in geometry and electron density. 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. A red dot above a residue indicates a poor fit to the electron density ( $RSRZ > 2$ ). 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.

#### • Molecule 1: HOMOGENITISATE 1,2-DIOXYGENASE



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 63 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	157.81Å 157.81Å 95.15Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	30.00 – 2.30 29.82 – 2.29	Depositor EDS
% Data completeness (in resolution range)	98.7 (30.00-2.30) 98.9 (29.82-2.29)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.16 (at 2.29Å)	Xtriage
Refinement program	X-PLOR 3.851	Depositor
R, $R_{free}$	0.193 , 0.242 0.189 , 0.236	Depositor DCC
$R_{free}$ test set	1576 reflections (5.02%)	DCC
Wilson B-factor (Å <sup>2</sup> )	32.5	Xtriage
Anisotropy	0.460	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.36 , 49.6	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.51$ , $\langle L^2 \rangle = 0.34$	Xtriage
Outliers	0 of 31486 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3544	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.76% 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, 0.375 respectively for untwinned datasets, and 0.333, 0.2 for perfectly twinned datasets.

## 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: FE2

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	0.51	2/3400 (0.1%)	0.75	1/4600 (0.0%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	142	MSE	SE-CE	-6.07	1.59	1.95
1	A	186	MSE	CG-SE	-5.14	1.77	1.95

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	96	ARG	NE-CZ-NH1	5.88	123.24	120.30

There are no chirality outliers.

There are no planarity outliers.

### 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	3311	0	3213	35	0
2	A	1	0	0	0	0
3	A	232	0	0	4	0
All	All	3544	0	3213	35	0

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 5.

All (35) 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:112:PHE:HB3	1:A:142:MSE:HE2	1.45	0.98
1:A:94:GLN:HE22	1:A:96:ARG:HH21	1.30	0.78
1:A:142:MSE:HE3	1:A:190:ILE:H	1.51	0.75
1:A:283:MSE:HG3	3:A:2049:HOH:O	1.86	0.75
1:A:142:MSE:HE3	1:A:190:ILE:N	2.09	0.67
1:A:94:GLN:HE22	1:A:96:ARG:NH2	1.92	0.67
1:A:96:ARG:HH11	1:A:357:PHE:HD2	1.44	0.64
1:A:96:ARG:HD3	1:A:363:SER:HB3	1.81	0.61
1:A:98:LYS:HG2	1:A:417:ARG:HH21	1.72	0.54
1:A:292:HIS:HD2	1:A:330:ARG:O	1.91	0.53
1:A:380:LYS:O	1:A:384:VAL:HG13	2.09	0.52
1:A:390:ARG:HD3	1:A:393:ASP:OD2	2.11	0.51
1:A:125:ILE:HG22	3:A:2208:HOH:O	2.10	0.51
1:A:142:MSE:HE3	1:A:190:ILE:HB	1.94	0.49
1:A:298:PHE:HB3	1:A:317:ILE:O	2.12	0.48
1:A:87:GLU:HB2	1:A:410:LYS:HE3	1.95	0.48
1:A:78:GLU:HG3	3:A:2134:HOH:O	2.14	0.48
1:A:98:LYS:HG3	1:A:417:ARG:HE	1.80	0.46
1:A:87:GLU:HG3	1:A:410:LYS:HG3	1.98	0.46
1:A:164:LEU:HB3	1:A:191:ASP:HB2	1.98	0.46
1:A:94:GLN:NE2	1:A:365:HIS:NE2	2.64	0.46
1:A:136:PHE:CZ	1:A:198:GLY:HA3	2.50	0.46
1:A:87:GLU:HG3	1:A:88:VAL:HG23	1.97	0.45
1:A:323:GLY:HA3	1:A:391:ILE:HD11	1.98	0.45
1:A:292:HIS:CD2	1:A:330:ARG:O	2.70	0.45
1:A:73:PHE:CD2	1:A:256:ALA:HB2	2.52	0.44
1:A:186:MSE:SE	1:A:274:PRO:HG3	2.68	0.44
1:A:142:MSE:CE	1:A:190:ILE:HB	2.48	0.42
1:A:35:CYS:HB3	3:A:2102:HOH:O	2.19	0.42
1:A:98:LYS:HG2	1:A:417:ARG:NH2	2.35	0.42
1:A:323:GLY:HA3	1:A:391:ILE:CD1	2.50	0.42
1:A:182:ILE:HG21	1:A:188:PHE:CD2	2.56	0.41
1:A:78:GLU:OE2	1:A:252:LYS:HA	2.21	0.40
1:A:96:ARG:NH1	1:A:357:PHE:CD2	2.86	0.40
1:A:298:PHE:HB2	1:A:316:VAL:HG13	2.04	0.40

There are no symmetry-related clashes.



## 5.3 Torsion angles

### 5.3.1 Protein backbone

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	413/471 (88%)	400 (97%)	13 (3%)	0	100	100

There are no Ramachandran outliers to report.

### 5.3.2 Protein sidechains

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	361/394 (92%)	353 (98%)	8 (2%)	60	77

All (8) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	3	GLU
1	A	58	ARG
1	A	68	VAL
1	A	106	SER
1	A	283	MSE
1	A	343	MSE
1	A	359	PRO
1	A	390	ARG

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (4) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	12	ASN
1	A	31	ASN
1	A	94	GLN
1	A	292	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules in this entry.

## 5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

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

## 5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 5.6 Ligand geometry ⓘ

Of 1 ligands modelled in this entry, 1 is monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 5.7 Other polymers ⓘ

There are no such residues in this entry.

## 5.8 Polymer linkage issues ⓘ

There are no chain breaks in this entry.

## 6 Fit of model and data ⓘ

### 6.1 Protein, DNA and RNA chains ⓘ

In the following table, the column labelled ‘#RSRZ> 2’ contains the number (and percentage) of RSRZ outliers, followed by percent RSRZ outliers for the chain as percentile scores relative to all X-ray entries and entries of similar resolution. The OWAB column contains the minimum, median, 95<sup>th</sup> percentile and maximum values of the occupancy-weighted average B-factor per residue. The column labelled ‘Q< 0.9’ lists the number of (and percentage) of residues with an average occupancy less than 0.9.

Mol	Chain	Analysed	<RSRZ>	#RSRZ>2		OWAB(Å <sup>2</sup> )	Q<0.9
1	A	410/471 (87%)	-0.13	25 (6%)	25 33	14, 29, 60, 75	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	356	GLY	7.6
1	A	379	GLU	5.2
1	A	376	ASP	4.5
1	A	2	ALA	4.4
1	A	380	LYS	4.4
1	A	78	GLU	3.7
1	A	385	LYS	3.6
1	A	384	VAL	3.5
1	A	383	LYS	3.5
1	A	377	CYS	3.2
1	A	106	SER	3.1
1	A	107	GLN	3.0
1	A	386	LEU	2.8
1	A	96	ARG	2.8
1	A	437	ASN	2.7
1	A	439	ARG	2.5
1	A	375	ALA	2.5
1	A	3	GLU	2.5
1	A	358	LEU	2.4
1	A	417	ARG	2.3
1	A	392	ALA	2.2
1	A	87	GLU	2.2
1	A	440	ASN	2.1
1	A	357	PHE	2.1
1	A	387	ALA	2.0

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

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

## 6.3 Carbohydrates [i](#)

There are no carbohydrates in this entry.

## 6.4 Ligands [i](#)

In the following table, the Atoms column lists the number of modelled atoms in the group and the number defined in the chemical component dictionary. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> percentile and maximum values of B factors of atoms in the group. The column labelled 'Q< 0.9' lists the number of atoms with occupancy less than 0.9.

Mol	Type	Chain	Res	Atoms	RSCC	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	FE2	A	1001	1/1	0.88	0.07	-	39,39,39,39	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.