

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

#### Nov 7, 2023 – 06:20 PM JST

PDB ID	:	8H85
Title	:	Trans-3/4-proline-hydroxylase H11 with 3-hydroxyl-proline
Authors	:	Gong, W.M.; Hu, X.Y.
Deposited on		
Resolution	:	2.38  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 (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:

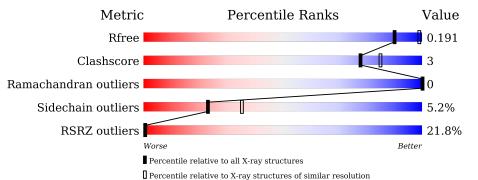
Xtriage (Phenix) EDS buster-report Percentile statistics Refmac CCP4 Ideal geometry (proteins) Ideal geometry (DNA, RNA)	:::::::::::::::::::::::::::::::::::::::	20191225.v01 (using entries in the PDB archive December 25th 2019) 5.8.0158 7.0.044 (Gargrove) Engh & Huber (2001) Parkinson et al. (1996)
Ideal geometry (DNA, RNA) Validation Pipeline (wwPDB-VP)		Parkinson et al. (1996) 2.36

# 1 Overall quality at a glance (i)

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

The reported resolution of this entry is 2.38 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	$\begin{array}{c} \textbf{Whole archive} \\ (\# \textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R <sub>free</sub>	130704	5509 (2.40-2.36)
Clashscore	141614	6082 (2.40-2.36)
Ramachandran outliers	138981	5973 (2.40-2.36)
Sidechain outliers	138945	5975 (2.40-2.36)
RSRZ outliers	127900	5397 (2.40-2.36)

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 <=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		
			21%		
1	А	268	84%	12%	••



# 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 2146 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.

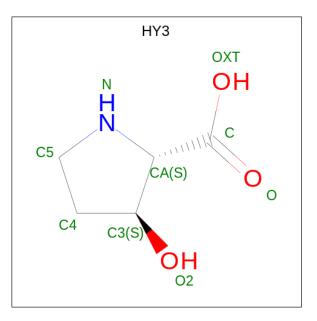
• Molecule 1 is a protein called Phytanoyl-CoA dioxygenase.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	А	262	Total 2076	C 1316	N 373	O 382	${ m S}{ m 5}$	0	2	0

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
А	-1	GLN	-	expression tag	UNP S5TUM1
А	0	ILE	-	expression tag	UNP S5TUM1

• Molecule 2 is 3-HYDROXYPROLINE (three-letter code: HY3) (formula: C<sub>5</sub>H<sub>9</sub>NO<sub>3</sub>) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Atoms				ZeroOcc	AltConf
2	А	1	Total 9	C 5	N 1	O 3	0	0

• Molecule 3 is water.

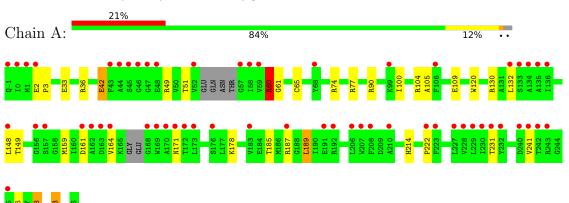


Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	61	Total O 61 61	0	0



# 3 Residue-property plots (i)

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 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: Phytanoyl-CoA dioxygenase



## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 2 2	Depositor
Cell constants	75.39Å 75.39Å 143.65Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$90.00^{\circ}$ $90.00^{\circ}$ $90.00^{\circ}$	Depositor
Resolution (Å)	32.82 - 2.38	Depositor
Resolution (A)	32.82 - 2.38	EDS
% Data completeness	99.8 (32.82-2.38)	Depositor
(in resolution range)	99.9 (32.82-2.38)	EDS
R <sub>merge</sub>	0.07	Depositor
R <sub>sym</sub>	(Not available)	Depositor
$< I/\sigma(I) > 1$	$4.27 (at 2.39 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0352	Depositor
D D.	0.164 , $0.188$	Depositor
$R, R_{free}$	0.172 , $0.191$	DCC
$R_{free}$ test set	854 reflections $(4.97%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	39.7	Xtriage
Anisotropy	0.168	Xtriage
Bulk solvent $k_{sol}(e/Å^3), B_{sol}(Å^2)$	0.38, $51.3$	EDS
L-test for twinning <sup>2</sup>	$ \langle L  \rangle = 0.48, \langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	2146	wwPDB-VP
Average B, all atoms $(Å^2)$	53.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.44% 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.

# 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:  $\rm HY3$ 

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).

Mal	Chain	Bo	nd lengths	Bond angles		
Mol	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	А	0.59	3/2130~(0.1%)	0.86	1/2893~(0.0%)	

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	А	0	9

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Ζ	Observed(Å)	Ideal(Å)
1	А	42	GLU	CD-OE1	6.08	1.32	1.25
1	А	248	GLU	CD-OE1	5.42	1.31	1.25
1	А	33	GLU	CD-OE2	5.07	1.31	1.25

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	$\mathbf{Z}$	$Observed(^{o})$	$Ideal(^{o})$
1	А	49	ARG	NE-CZ-NH2	-5.28	117.66	120.30

There are no chirality outliers.

All (9) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	А	104	ARG	Sidechain
1	А	130	ARG	Sidechain
1	А	187	ARG	Sidechain

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Mol	Chain	Res	Type	Group
1	А	246	ARG	Sidechain
1	А	253	ARG	Sidechain
1	А	60	ARG	Sidechain
1	А	74[A]	ARG	Sidechain
1	А	77[A]	ARG	Sidechain
1	А	90	ARG	Sidechain

 $Continued \ from \ previous \ page...$ 

## 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	А	2076	0	1985	12	0
2	А	9	0	2	0	0
3	А	61	0	0	0	0
All	All	2146	0	1987	12	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 3.

All (12) 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:42:GLU:OE1	1:A:65:CYS:SG	2.63	0.56
1:A:105:ALA:HB1	1:A:222:PRO:HA	1.90	0.54
1:A:60:ARG:NH2	1:A:109:GLU:OE1	2.42	0.53
1:A:120:TRP:CH2	1:A:231:THR:HG21	2.46	0.50
1:A:189:LEU:HD12	1:A:189:LEU:HA	1.77	0.47
1:A:149:THR:O	1:A:214:HIS:HA	2.15	0.46
1:A:61:GLY:HA2	1:A:100:ILE:O	2.17	0.45
1:A:253:ARG:HH11	1:A:253:ARG:HG3	1.82	0.44
1:A:164:VAL:HA	1:A:178:LYS:O	2.20	0.41
1:A:2:GLU:HB2	1:A:3:PRO:HD2	2.02	0.41
1:A:42:GLU:OE1	1:A:65:CYS:CB	2.69	0.41
1:A:51:THR:HG22	1:A:60:ARG:HD2	2.03	0.41

There are no symmetry-related clashes.



## 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	А	258/268~(96%)	250~(97%)	8 (3%)	0	100 100

There are no Ramachandran outliers to report.

#### 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	А	214/225~(95%)	203~(95%)	11 (5%)	24 36	

All (11) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	36	ARG
1	А	60	ARG
1	А	132	LEU
1	А	148	LEU
1	А	159	MET
1	А	161	ASP
1	А	171	ASN
1	А	185	THR
1	А	189	LEU
1	А	241	VAL
1	А	248	GLU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:



Mol	Chain	Res	Type
1	А	171	ASN

#### 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)

1 ligand is 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).

Mo	l Type	Chain	Res	Link	В	ond leng	gths	В	ond ang	gles
IVIC	I Iype	Ullalli	nes		Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z >2
2	HY3	А	301	-	8,9,9	1.39	1 (12%)	8,12,12	1.19	1 (12%)

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	HY3	А	301	-	-	0/4/14/14	0/1/1/1

All (1) bond length outliers are listed below:



Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	А	301	HY3	OXT-C	-3.49	1.19	1.30

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$Observed(^{o})$	$Ideal(^{o})$
2	А	301	HY3	C-CA-N	2.64	116.48	111.19

There are no chirality outliers.

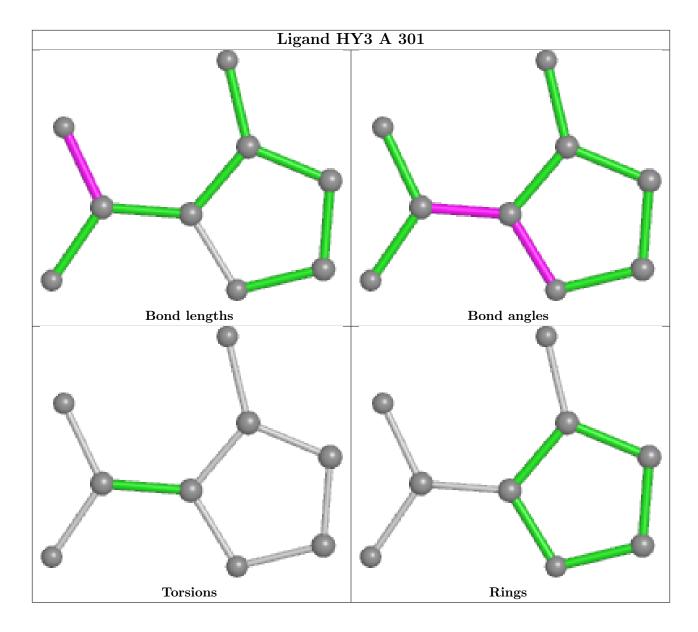
There are no torsion outliers.

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 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)

There are no chain breaks in this entry.



## 6 Fit of model and data (i)

## 6.1 Protein, DNA and RNA chains (i)

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,  $95^{th}$  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	$\mathbf{OWAB}(\mathbf{\AA}^2)$	Q<0.9
1	А	262/268~(97%)	0.98	57 (21%) 0 0	23, 45, 105, 138	0

All (57) RSRZ outliers are listed below:

Mol			Type	RSRZ	
1	А	170	ALA	9.1	
1	А	1	MET	8.4	
1	А	58	ILE	7.1	
1	А	161	ASP	6.3	
1	А	0	ILE	6.1	
1	А	168	GLY	5.7	
1	А	59	VAL	5.3	
1	А	134	ALA	5.3	
1	А	171	ASN	5.2	
1	А	169	TRP	5.1	
1	А	245	THR	5.0	
1	А	44	ALA	4.8	
1	А	229	LEU	4.8	
1	А	230	ILE	4.5	
1	А	163	ASP	4.5	
1	А	176	SER	4.4	
1	А	2	GLU	4.4	
1	А	57	GLY	4.3	
1	А	243	ARG	4.3	
1	А	45	SER	4.1	
1	А	183	VAL	4.0	
1	А	133	SER	3.9	
1	А	206	LEU	3.9	
1	А	135	ALA	3.8	
1	А	208	PHE	3.8	
1	А	172	THR	3.8	
1	A	148	LEU	3.8	

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Conti	Continued from previous page								
Mol	Chain	Res	Type	RSRZ					
1	А	46	GLY	3.7					
1	А	223	PHE	3.6					
1	А	157	SER	3.5					
1	А	242	THR	3.4					
1	А	228	VAL	3.4					
1	А	164	VAL	3.3					
1	А	231	THR	3.2					
1	А	173	LEU	3.2					
1	А	132	LEU	3.1					
1	А	240	ASP	2.8					
1	А	241	VAL	2.7					
1	А	222	PRO	2.7					
1	А	192	ARG	2.6					
1	А	43	PHE	2.6					
1	А	47	GLY	2.6					
1	А	156	GLY	2.5					
1	А	136	ILE	2.4					
1	А	99	LYS	2.3					
1	А	-1	GLN	2.3					
1	А	207	TRP	2.2					
1	А	48	GLU	2.2					
1	А	191	GLU	2.2					
1	А	232	TYR	2.2					
1	А	106	PHE	2.2					
1	А	187	ARG	2.2					
1	А	68	TYR	2.1					
1	А	227	LEU	2.1					

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

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

2.1

2.0

2.0

#### 6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

162

52

210

А

А

А

1

1

1

ALA

VAL

ALA



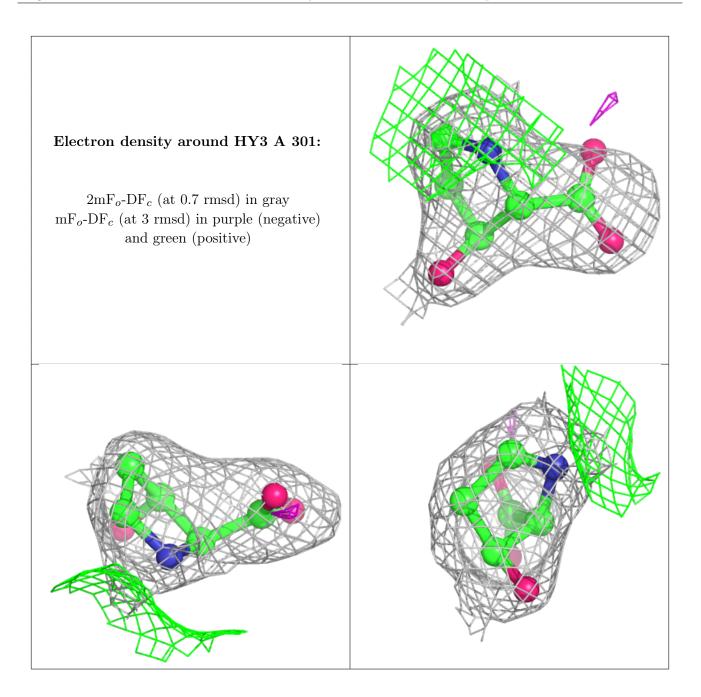
### 6.4 Ligands (i)

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. The B-factors column lists the minimum, median,  $95^{th}$  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	$\mathbf{B} ext{-factors}(\mathrm{\AA}^2)$	Q<0.9
2	HY3	А	301	9/9	0.89	0.36	55,70,78,78	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





## 6.5 Other polymers (i)

There are no such residues in this entry.

