



# Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 10:37 AM UTC

PDB ID : 1YIP / pdb\_00001yip  
Title : Oxidized Peptidylglycine Alpha-Hydroxylating Monooxygenase (PHM) in a New Crystal Form  
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Deposited on : 2005-01-12  
Resolution : 2.20 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 : 4-5-2 with Phenix2.0  
Xtrriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : 2.49

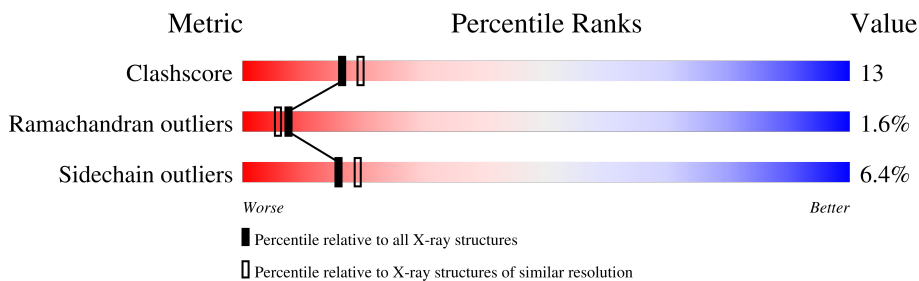
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.20 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	190562	6851 (2.20-2.20)
Ramachandran outliers	187476	6768 (2.20-2.20)
Sidechain outliers	187428	6769 (2.20-2.20)

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  $\geq 3$ , 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	311	 76% 20% . .

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 2585 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 Peptidyl-glycine alpha-amidating monooxygenase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	311	2425	1544	410	445	26	0	0	0

- Molecule 2 is COPPER (II) ION (CCD ID: CU) (formula: Cu).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	2	Total	Cu	0	0
			2	2		

- Molecule 3 is water.

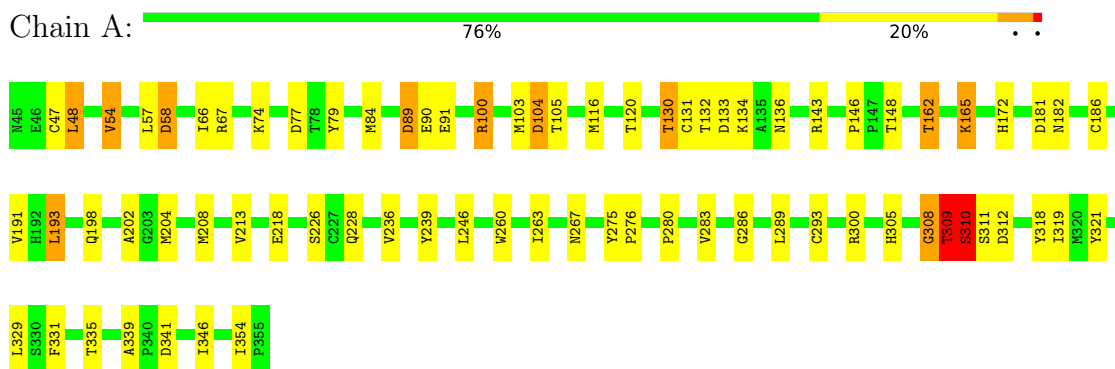
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	158	Total	O	0	0
			158	158		

### 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. 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: Peptidyl-glycine alpha-amidating monooxygenase



## 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, $\alpha$ , $\beta$ , $\gamma$	58.52Å 65.68Å 69.87Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.01 – 2.20	Depositor
% Data completeness (in resolution range)	98.9 (30.01-2.20)	Depositor
$R_{merge}$	0.06	Depositor
$R_{sym}$	0.06	Depositor
Refinement program	REFMAC refmac_5.2.0003	Depositor
R, $R_{free}$	0.190 , 0.239	Depositor
Estimated twinning fraction	No twinning to report.	Xtrriage
Total number of atoms	2585	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	42.0	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: CU

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	1.06	2/2494 (0.1%)	1.09	3/3393 (0.1%)

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	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	339	ALA	CA-CB	-7.69	1.44	1.53
1	A	283	VAL	CA-CB	-5.11	1.48	1.54

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	54	VAL	CB-CA-C	-8.23	98.31	110.81
1	A	146	PRO	N-CA-C	7.07	119.32	110.70
1	A	263	ILE	N-CA-C	-5.92	106.04	111.67

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	309	THR	Peptide

## 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	2425	0	2344	62	0
2	A	2	0	0	0	0
3	A	158	0	0	10	0
All	All	2585	0	2344	62	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 13.

All (62) 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:134:LYS:NZ	1:A:331:PHE:CG	2.21	1.08
1:A:267:ASN:HB2	1:A:354:ILE:HG23	1.47	0.93
1:A:134:LYS:NZ	1:A:331:PHE:CD1	2.36	0.92
1:A:89:ASP:OD2	1:A:165:LYS:NZ	2.05	0.88
1:A:136:ASN:HB3	3:A:427:HOH:O	1.81	0.79
1:A:116:MET:SD	1:A:133:ASP:HB3	2.26	0.75
1:A:57:LEU:O	1:A:58:ASP:HB3	1.90	0.71
1:A:198:GLN:O	3:A:409:HOH:O	2.09	0.70
1:A:133:ASP:OD2	3:A:377:HOH:O	2.11	0.68
1:A:134:LYS:NZ	1:A:331:PHE:CB	2.61	0.63
1:A:105:THR:HG21	3:A:499:HOH:O	1.98	0.62
1:A:134:LYS:HZ1	1:A:331:PHE:HB2	1.65	0.62
1:A:77:ASP:OD2	1:A:172:HIS:NE2	2.31	0.60
1:A:91:GLU:OE2	3:A:390:HOH:O	2.16	0.60
1:A:228:GLN:HE22	1:A:286:GLY:CA	2.15	0.59
1:A:228:GLN:NE2	1:A:286:GLY:C	2.62	0.57
1:A:239:TYR:HA	1:A:318:TYR:O	2.04	0.57
1:A:191:VAL:HG12	1:A:193:LEU:HD22	1.88	0.56
1:A:67:ARG:NH2	1:A:100:ARG:HD3	2.22	0.55
1:A:134:LYS:NZ	1:A:331:PHE:HB2	2.21	0.55
1:A:309:THR:HB	1:A:311:SER:HB2	1.88	0.55
1:A:309:THR:O	1:A:310:SER:C	2.50	0.54
1:A:74:LYS:NZ	3:A:469:HOH:O	2.41	0.54
1:A:57:LEU:O	1:A:58:ASP:CB	2.56	0.53

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:331:PHE:CD1	1:A:331:PHE:C	2.87	0.53
1:A:116:MET:O	1:A:131:CYS:HB3	2.10	0.52
1:A:181:ASP:HB3	3:A:478:HOH:O	2.10	0.52
1:A:182:ASN:OD1	1:A:182:ASN:C	2.53	0.51
1:A:208:MET:HG3	1:A:335:THR:HG22	1.91	0.51
1:A:79:TYR:CE1	1:A:172:HIS:CD2	3.00	0.50
1:A:228:GLN:NE2	1:A:286:GLY:CA	2.74	0.50
1:A:308:GLY:C	1:A:310:SER:N	2.69	0.49
1:A:103:MET:HE1	1:A:143:ARG:HG3	1.96	0.48
1:A:47:CYS:O	1:A:48:LEU:HD12	2.13	0.48
1:A:191:VAL:HG12	1:A:193:LEU:CD2	2.42	0.48
1:A:148:THR:OG1	1:A:276:PRO:HG3	2.15	0.47
1:A:74:LYS:HG2	1:A:181:ASP:HB2	1.96	0.47
1:A:202:ALA:HA	1:A:321:TYR:O	2.15	0.47
1:A:275:TYR:HA	1:A:276:PRO:HD3	1.78	0.47
1:A:67:ARG:HD2	1:A:186:CYS:O	2.15	0.46
1:A:213:VAL:HG22	1:A:305:HIS:CD2	2.50	0.46
1:A:162:THR:HB	3:A:515:HOH:O	2.14	0.46
1:A:329:LEU:H	1:A:329:LEU:HD23	1.79	0.46
1:A:79:TYR:CZ	1:A:172:HIS:CD2	3.04	0.45
1:A:236:VAL:O	1:A:280:PRO:HA	2.17	0.44
1:A:246:LEU:HD21	1:A:300:ARG:HD3	1.99	0.44
1:A:67:ARG:NH2	1:A:100:ARG:CD	2.80	0.44
1:A:260:TRP:CZ3	1:A:346:ILE:HG13	2.53	0.44
1:A:204:MET:HG2	3:A:472:HOH:O	2.18	0.43
1:A:228:GLN:HE22	1:A:286:GLY:N	2.17	0.43
1:A:354:ILE:HG23	1:A:354:ILE:O	2.19	0.43
1:A:321:TYR:C	1:A:321:TYR:CD1	2.97	0.43
1:A:130:THR:HG22	1:A:131:CYS:N	2.34	0.42
1:A:226:SER:HA	1:A:289:LEU:O	2.20	0.42
1:A:236:VAL:CG1	1:A:319:ILE:HG23	2.50	0.42
1:A:58:ASP:OD2	1:A:58:ASP:C	2.63	0.41
1:A:228:GLN:HE22	1:A:286:GLY:H	1.68	0.41
1:A:47:CYS:C	1:A:48:LEU:HD12	2.45	0.41
1:A:331:PHE:CD1	1:A:331:PHE:O	2.74	0.41
1:A:329:LEU:HD23	1:A:329:LEU:N	2.36	0.41
1:A:103:MET:O	1:A:104:ASP:C	2.63	0.41
1:A:218:GLU:HG3	3:A:487:HOH:O	2.21	0.40

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	A	309/311 (99%)	291 (94%)	13 (4%)	5 (2%)	<b>7</b> <b>6</b>

All (5) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	58	ASP
1	A	310	SER
1	A	309	THR
1	A	308	GLY
1	A	312	ASP

### 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	266/266 (100%)	249 (94%)	17 (6%)	<b>16</b> <b>19</b>

All (17) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	48	LEU
1	A	54	VAL
1	A	66	ILE
1	A	84	MET
1	A	89	ASP
1	A	90	GLU

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Mol	Chain	Res	Type
1	A	100	ARG
1	A	104	ASP
1	A	120	THR
1	A	130	THR
1	A	132	THR
1	A	162	THR
1	A	165	LYS
1	A	193	LEU
1	A	293	CYS
1	A	310	SER
1	A	341	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	305	HIS
1	A	316	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 oligosaccharides in this entry.

### 5.6 Ligand geometry [i](#)

Of 2 ligands modelled in this entry, 2 are 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 [i](#)

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

EDS was not executed - this section is therefore empty.

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

EDS was not executed - this section is therefore empty.

### 6.3 Carbohydrates [i](#)

EDS was not executed - this section is therefore empty.

### 6.4 Ligands [i](#)

EDS was not executed - this section is therefore empty.

### 6.5 Other polymers [i](#)

EDS was not executed - this section is therefore empty.