

wwPDB X-ray Structure Validation Summary Report (i)

Sep 22, 2021 – 10:02 am BST

PDB ID : 701X

Title: Unspecific peroxygenase from Hypoxylon sp. EC38 in complex with 1-

phenylimidazole

Authors : Rotilio, L.; Mattevi, A.

Deposited on : 2021-03-30

Resolution : 1.60 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 following versions of software and data (see references (1)) were used in the production of this report:

MolProbity: 4.02b-467

Mogul : 1.8.5 (274361), CSD as541be (2020)

Xtriage (Phenix) : NOT EXECUTED

EDS : NOT EXECUTED

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)

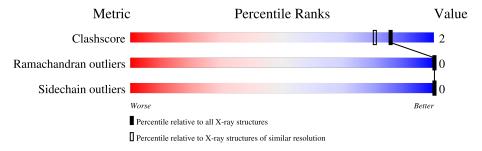
Validation Pipeline (wwPDB-VP) : 2.23.1

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X-RAY DIFFRACTION

The reported resolution of this entry is 1.60 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	Similar resolution
Metric	$(\# \mathrm{Entries})$	$(\# ext{Entries}, ext{resolution range}(ext{Å}))$
Clashscore	141614	3665 (1.60-1.60)
Ramachandran outliers	138981	3564 (1.60-1.60)
Sidechain outliers	138945	3563 (1.60-1.60)

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%

Note EDS was not executed.

Mol	Chain	Length	Quality of chain					
1	A	261	81%	5%	14%			



2 Entry composition (i)

There are 8 unique types of molecules in this entry. The entry contains 2108 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.

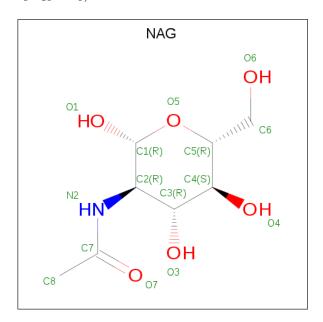
• Molecule 1 is a protein called Peroxygenase.

Mol	Chain	Residues		Ato	oms			ZeroOcc	AltConf	Trace
1	Δ	225	Total	С	N	О	S	0	6	0
1	Λ	220	1774	1127	297	346	4	0	О	

• Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

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

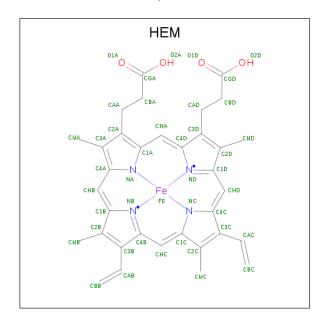
• Molecule 3 is 2-acetamido-2-deoxy-beta-D-glucopyranose (three-letter code: NAG) (formula: $C_8H_{15}NO_6$).



Mol	Chain	Residues	${f Atoms}$				ZeroOcc	AltConf
2	Λ	1	Total	С	N	О	0	0
)	A	1	14	8	1	5		
2	Λ	1	Total	С	N	О	0	0
3	А	1	14	8	1	5	0	

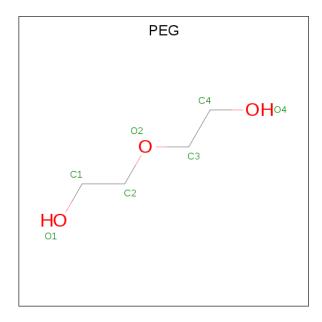


• Molecule 4 is PROTOPORPHYRIN IX CONTAINING FE (three-letter code: HEM) (formula: $C_{34}H_{32}FeN_4O_4$).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf		
4	Λ	1	Total	С	Fe	Ν	О	0	0
$\begin{array}{c c} 4 & A \end{array}$	1	43	34	1	4	4	U	U	

• Molecule 5 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: $C_4H_{10}O_3$).



Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
5	A	1	Total C (Э 3	0	0

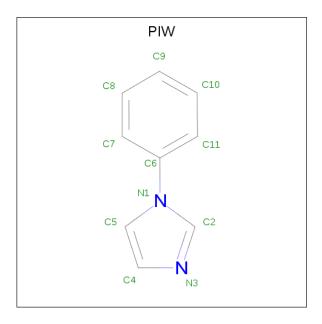
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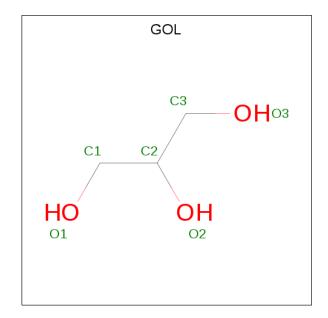
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
5	A	1	Total C O 7 4 3	0	0

• Molecule 6 is 1-phenyl-1H-imidazole (three-letter code: PIW) (formula: $C_9H_8N_2$) (labeled as "Ligand of Interest" by depositor).



Mol	Chain	Residues	Ato	oms		ZeroOcc	AltConf
6	A	1	Total 11	C 9	N 2	0	0

 \bullet Molecule 7 is GLYCEROL (three-letter code: GOL) (formula: $\mathrm{C_3H_8O_3}).$





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	1	Total C O 6 3 3	0	0

• Molecule 8 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
8	A	231	Total O 231 231	0	0

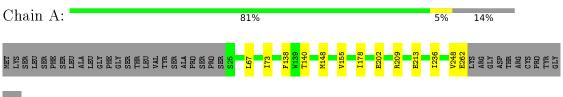


3 Residue-property plots (i)

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. 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: Peroxygenase







4 Data and refinement statistics (i)

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants	71.53Å 71.53Å 153.27Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.08 - 1.60	Depositor
% Data completeness	99.9 (48.08-1.60)	Depositor
(in resolution range)	33.3 (40.00 1.00)	Depositor
R_{merge}	0.16	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	REFMAC 5.8.0253	Depositor
R, R_{free}	0.134 , 0.159	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2108	wwPDB-VP
Average B, all atoms (Å ²)	22.0	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: HEM, GOL, MG, PEG, PIW, NAG

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	Boı	nd lengths	Bo	nd angles
MIOI	Chain	RMSZ	# Z > 5	RMSZ	# Z >5
1	Α	0.96	$5/1828 \; (0.3\%)$	0.95	1/2493 (0.0%)

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	${ m Observed(\AA)}$	$\operatorname{Ideal}(ext{\AA})$
1	A	202	GLU	CD-OE1	-8.83	1.16	1.25
1	A	262	GLU	CD-OE1	8.71	1.35	1.25
1	A	248	VAL	C-N	7.87	1.52	1.34
1	A	213	GLU	CD-OE1	6.31	1.32	1.25
1	A	209	ARG	NE-CZ	5.30	1.40	1.33

All (1) bond angle outliers are listed below:

Mol	Chain	${f Res}$	Type	${f Atoms}$	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\mathbf{Ideal}(^o)$
1	A	262	GLU	CA-C-O	-7.32	104.74	120.10

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	1774	0	1706	7	1	
2	A	1	0	0	0	0	

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-	110111	picolous	payc

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	28	0	26	0	0
4	A	43	0	30	1	0
5	A	14	0	20	0	0
6	A	11	0	8	0	0
7	A	6	0	8	0	0
8	A	231	0	0	0	0
All	All	2108	0	1798	8	1

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 2.

The worst 5 of 8 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} ({\rm \AA}) \end{array}$	Clash overlap (Å)
1:A:140:THR:N	1:A:148[B]:MET:HE3	2.17	0.60
1:A:138:PHE:CE1	1:A:155[B]:VAL:HG11	2.38	0.59
1:A:178:ILE:CD1	1:A:236[A]:ILE:HD11	2.40	0.52
1:A:67:LEU:HB3	1:A:73:ILE:HB	1.97	0.46
1:A:138:PHE:CE2	1:A:155[A]:VAL:HG21	2.54	0.42

All (1) symmetry-related close contacts are listed below. The label for Atom-2 includes the symmetry operator and encoded unit-cell translations to be applied.

Atom-1	Atom-2	$\begin{array}{c} \text{Interatomic} \\ \text{distance (Å)} \end{array}$	$egin{array}{c} ext{Clash} \ ext{overlap } (ext{Å}) \end{array}$
1:A:140:THR:OG1	1:A:140:THR:OG1[8_555]	2.13	0.07

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	228/261 (87%)	223 (98%)	5 (2%)	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	A	195/219 (89%)	195 (100%)	0	100 100

There are no protein residues with a non-rotameric sidechain to report.

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	28	GLN
1	A	93	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)

Of 8 ligands modelled in this entry, 1 is monoatomic - leaving 7 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	Tree	Chain	Res	Link	Bo	Bond lengths			Bond angles		
MIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2	
3	NAG	A	302	1	14,14,15	1.35	2 (14%)	17,19,21	1.34	4 (23%)	
3	NAG	A	304	1	14,14,15	1.14	1 (7%)	17,19,21	1.96	7 (41%)	
5	PEG	A	307	-	6,6,6	0.58	0	5,5,5	0.53	0	
7	GOL	A	308	-	5,5,5	0.35	0	5,5,5	0.55	0	
5	PEG	A	305	-	6,6,6	0.43	0	5,5,5	0.55	0	
6	PIW	A	306	4	12,12,12	2.89	4 (33%)	13,15,15	2.81	8 (61%)	
4	HEM	A	303	6,2,1	27,50,50	1.89	5 (18%)	17,82,82	1.78	6 (35%)	

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	\mathbf{Rings}
3	NAG	A	302	1	-	0/6/23/26	0/1/1/1
3	NAG	A	304	1	-	0/6/23/26	0/1/1/1
5	PEG	A	307	-	-	3/4/4/4	1
7	GOL	A	308	_	-	2/4/4/4	ı
5	PEG	A	305	-	-	2/4/4/4	-
6	PIW	A	306	4	-	0/4/4/4	0/2/2/2
4	HEM	A	303	6,2,1	_	0/6/54/54	_

The worst 5 of 12 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	$\operatorname{Ideal}(\text{\AA})$
6	A	306	PIW	C6-N1	-6.58	1.34	1.45
4	A	303	HEM	C3B-C2B	-6.00	1.32	1.40
6	A	306	PIW	C5-N1	-5.67	1.31	1.39
4	A	303	HEM	C3C-C2C	-4.17	1.34	1.40
4	A	303	HEM	C4D-C3D	4.00	1.51	1.42

The worst 5 of 25 bond angle outliers are listed below:

Mol	Chain	Res	Type	Type Atoms		$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^{o})$
6	A	306	PIW	C5-N1-C2	5.12	117.78	108.50
6	A	306	PIW	C5-N1-C6	-3.86	120.07	126.33
6	A	306	PIW	C4-C5-N1	-3.67	100.02	106.50

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Mol	Chain	${f Res}$	Type	${f Atoms}$	\mathbf{Z}	$\operatorname{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
3	A	304	NAG	O5-C5-C6	-3.39	101.88	107.20
3	A	304	NAG	O7-C7-C8	3.38	128.33	122.06

There are no chirality outliers.

5 of 7 torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
7	A	308	GOL	C1-C2-C3-O3
5	A	305	PEG	O1-C1-C2-O2
5	A	305	PEG	O2-C3-C4-O4
7	A	308	GOL	O2-C2-C3-O3
5	A	307	PEG	O1-C1-C2-O2

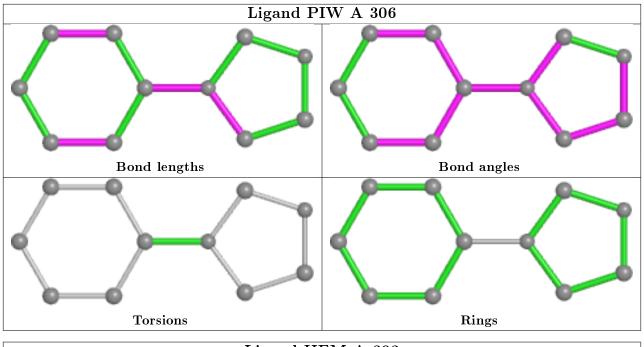
There are no ring outliers.

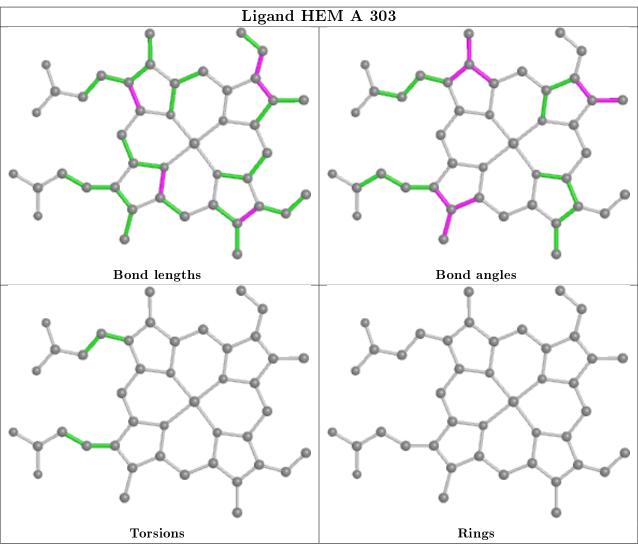
1 monomer is involved in 1 short contact:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
4	A	303	HEM	1	0

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)

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.

