



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 19, 2016 – 06:06 PM GMT

PDB ID : 1OYA  
Title : OLD YELLOW ENZYME AT 2 ANGSTROMS RESOLUTION: OVERALL  
STRUCTURE, LIGAND BINDING AND COMPARISON WITH RELATED  
FLAVOPROTEINS  
Authors : Fox, K.M.; Karplus, P.A.  
Deposited on : 1994-08-25  
Resolution : 2.00 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.

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The following versions of software and data (see [references ⓘ](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.7.1 (RC1), CSD as537be (2016)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : rb-20026982

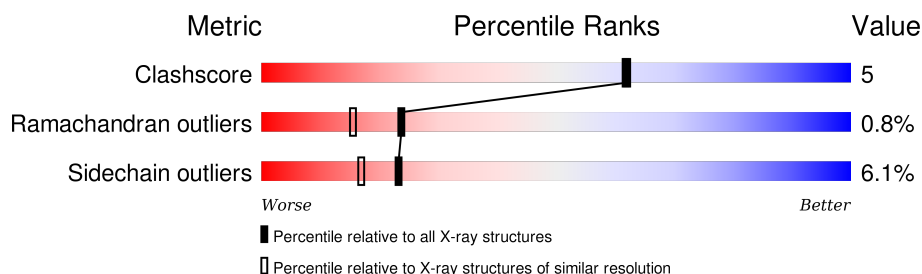
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.00 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	102246	7340 (2.00-2.00)
Ramachandran outliers	100387	7248 (2.00-2.00)
Sidechain outliers	100360	7247 (2.00-2.00)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	400	 80% 16% . .

## 2 Entry composition [i](#)

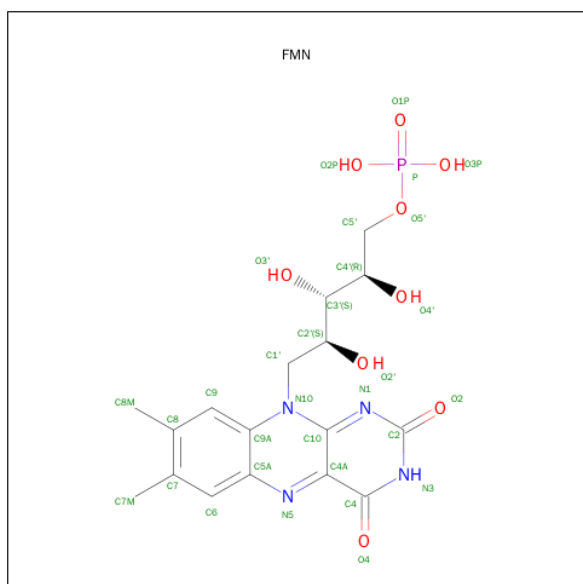
There are 3 unique types of molecules in this entry. The entry contains 3335 atoms, of which 0 are hydrogens and 0 are deuteriums.

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 OLD YELLOW ENZYME.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	399	Total	C	N	O	S	4	0	0
			3178	2027	551	594	6			

- Molecule 2 is FLAVIN MONONUCLEOTIDE (three-letter code: FMN) (formula:  $C_{17}H_{21}N_4O_9P$ ).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
2	A	1	Total	C	N	O	P	0	0
			31	17	4	9	1		

- Molecule 3 is water.

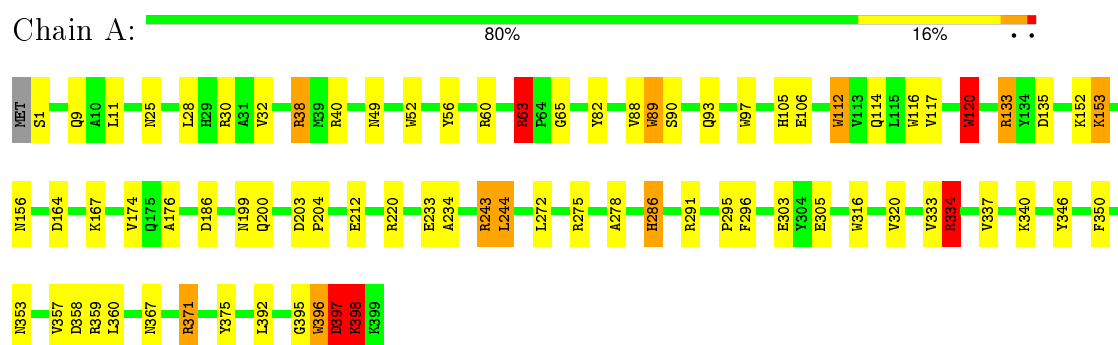
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	126	Total	O	0	0
			126	126		

### 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 ( $\text{RSRZ} > 2$ ). 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.

Note EDS was not executed.

#### • Molecule 1: OLD YELLOW ENZYME



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	142.88Å 142.88Å 43.01Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	10.00 – 2.00	Depositor
% Data completeness (in resolution range)	(Not available) (10.00-2.00)	Depositor
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.180 , 0.241	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	3335	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	19.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FMN

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	0.85	0/3260	1.55	52/4417 (1.2%)

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	2

There are no bond length outliers.

All (52) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	38	ARG	NE-CZ-NH2	-20.52	110.04	120.30
1	A	334	ARG	NE-CZ-NH1	15.48	128.04	120.30
1	A	38	ARG	NE-CZ-NH1	14.61	127.60	120.30
1	A	63	ARG	NE-CZ-NH1	12.28	126.44	120.30
1	A	397	ASP	CA-C-N	-11.72	91.42	117.20
1	A	334	ARG	NE-CZ-NH2	-10.81	114.90	120.30
1	A	220	ARG	NE-CZ-NH1	10.80	125.70	120.30
1	A	63	ARG	NE-CZ-NH2	-10.28	115.16	120.30
1	A	133	ARG	NE-CZ-NH1	10.05	125.33	120.30
1	A	38	ARG	CG-CD-NE	-9.58	91.69	111.80
1	A	220	ARG	NE-CZ-NH2	-8.61	116.00	120.30
1	A	52	TRP	CD1-CG-CD2	8.29	112.93	106.30
1	A	112	TRP	CD1-CG-CD2	7.84	112.57	106.30
1	A	133	ARG	NE-CZ-NH2	-7.70	116.45	120.30
1	A	316	TRP	CD1-CG-CD2	7.50	112.30	106.30
1	A	316	TRP	CE2-CD2-CG	-7.49	101.31	107.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	116	TRP	CD1-CG-CD2	7.32	112.16	106.30
1	A	116	TRP	CE2-CD2-CG	-7.20	101.54	107.30
1	A	112	TRP	CE2-CD2-CG	-7.19	101.55	107.30
1	A	52	TRP	CE2-CD2-CG	-7.13	101.59	107.30
1	A	359	ARG	NE-CZ-NH1	7.10	123.85	120.30
1	A	97	TRP	CE2-CD2-CG	-7.01	101.69	107.30
1	A	396	TRP	CE2-CD2-CG	-7.01	101.69	107.30
1	A	396	TRP	CD1-CG-CD2	6.82	111.75	106.30
1	A	89	TRP	CE2-CD2-CG	-6.65	101.98	107.30
1	A	97	TRP	CD1-CG-CD2	6.61	111.58	106.30
1	A	275	ARG	NE-CZ-NH1	6.44	123.52	120.30
1	A	375	TYR	CB-CG-CD2	-6.29	117.23	121.00
1	A	89	TRP	CD1-CG-CD2	6.24	111.29	106.30
1	A	233	GLU	CA-CB-CG	6.19	127.02	113.40
1	A	397	ASP	O-C-N	6.14	132.53	122.70
1	A	275	ARG	NE-CZ-NH2	-6.10	117.25	120.30
1	A	316	TRP	CG-CD2-CE3	5.99	139.29	133.90
1	A	212	GLU	N-CA-CB	-5.96	99.88	110.60
1	A	316	TRP	CB-CG-CD1	-5.93	119.29	127.00
1	A	60	ARG	NE-CZ-NH2	-5.86	117.37	120.30
1	A	243	ARG	NE-CZ-NH1	5.85	123.22	120.30
1	A	358	ASP	CB-CG-OD1	5.64	123.37	118.30
1	A	346	TYR	CB-CG-CD1	-5.60	117.64	121.00
1	A	244	LEU	CA-CB-CG	5.56	128.09	115.30
1	A	398	LYS	N-CA-C	-5.54	96.04	111.00
1	A	52	TRP	CG-CD1-NE1	-5.52	104.58	110.10
1	A	56	TYR	CB-CG-CD1	-5.48	117.71	121.00
1	A	52	TRP	CG-CD2-CE3	5.44	138.80	133.90
1	A	82	TYR	CB-CG-CD2	-5.37	117.78	121.00
1	A	30	ARG	NE-CZ-NH1	5.28	122.94	120.30
1	A	120	TRP	CE2-CD2-CG	-5.27	103.08	107.30
1	A	396	TRP	CG-CD2-CE3	5.24	138.62	133.90
1	A	112	TRP	CG-CD1-NE1	-5.23	104.87	110.10
1	A	117	VAL	CG1-CB-CG2	-5.22	102.55	110.90
1	A	359	ARG	NE-CZ-NH2	-5.10	117.75	120.30
1	A	296	PHE	CA-CB-CG	-5.04	101.81	113.90

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	396	TRP	Mainchain

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Mol	Chain	Res	Type	Group
1	A	397	ASP	Mainchain

## 5.2 Too-close contacts ⓘ

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	3178	0	3093	34	0
2	A	31	0	19	2	0
3	A	126	0	0	2	0
All	All	3335	0	3112	34	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 (34) 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:114:GLN:HE22	2:A:401:FMN:HN3	1.12	0.93
1:A:353:ASN:HD21	1:A:367:ASN:H	1.25	0.82
1:A:90:SER:H	1:A:93:GLN:NE2	1.97	0.63
1:A:114:GLN:NE2	2:A:401:FMN:HN3	1.93	0.62
1:A:9:GLN:HE21	1:A:334:ARG:NH1	1.99	0.60
1:A:63:ARG:HG2	1:A:357:VAL:HG11	1.85	0.59
1:A:392:LEU:HD22	1:A:398:LYS:HB3	1.85	0.59
1:A:371:ARG:CD	1:A:371:ARG:H	2.15	0.59
1:A:9:GLN:NE2	1:A:334:ARG:HH11	2.04	0.55
1:A:25:ASN:ND2	1:A:112:TRP:HE1	2.05	0.54
1:A:397:ASP:O	1:A:398:LYS:HB2	2.06	0.54
1:A:243:ARG:HH21	1:A:286:HIS:CE1	2.26	0.54
1:A:9:GLN:HE21	1:A:334:ARG:HH11	1.55	0.53
1:A:105:HIS:HE1	1:A:186:ASP:OD2	1.93	0.51
1:A:164:ASP:HA	1:A:167:LYS:HE3	1.92	0.50
1:A:333:VAL:O	1:A:337:VAL:HG22	2.12	0.50
1:A:120:TRP:H	1:A:200:GLN:HE22	1.61	0.49
1:A:152:LYS:HA	1:A:152:LYS:HD2	1.50	0.48
1:A:28:LEU:HB2	1:A:65:GLY:HA3	1.95	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:133:ARG:HD2	1:A:135:ASP:OD1	2.13	0.48
1:A:291:ARG:HA	1:A:303:GLU:HG3	1.96	0.47
1:A:89:TRP:H	1:A:93:GLN:HE22	1.63	0.46
1:A:89:TRP:H	1:A:93:GLN:NE2	2.14	0.45
1:A:371:ARG:H	1:A:371:ARG:HD2	1.82	0.43
1:A:174:VAL:HG13	1:A:234:ALA:HB2	2.00	0.43
1:A:295:PRO:HG3	3:A:616:HOH:O	2.18	0.43
1:A:88:VAL:HG13	1:A:176:ALA:HB1	2.01	0.43
1:A:395:GLY:O	1:A:397:ASP:N	2.49	0.43
1:A:153:LYS:HB2	1:A:153:LYS:NZ	2.35	0.42
1:A:90:SER:H	1:A:93:GLN:HE21	1.66	0.42
1:A:40:ARG:O	1:A:49:ASN:HB2	2.20	0.42
1:A:38:ARG:NH2	3:A:548:HOH:O	2.53	0.41
1:A:203:ASP:HA	1:A:204:PRO:HD2	1.85	0.41
1:A:353:ASN:ND2	1:A:367:ASN:H	2.03	0.41

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	397/400 (99%)	383 (96%)	11 (3%)	3 (1%)	24 15

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	397	ASP
1	A	398	LYS
1	A	278	ALA

### 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	328/329 (100%)	308 (94%)	20 (6%)	23	17

All (20) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	1	SER
1	A	11	LEU
1	A	32	VAL
1	A	63	ARG
1	A	106	GLU
1	A	120	TRP
1	A	153	LYS
1	A	156	ASN
1	A	199	ASN
1	A	244	LEU
1	A	272	LEU
1	A	286	HIS
1	A	305	GLU
1	A	320	VAL
1	A	334	ARG
1	A	340	LYS
1	A	350	PHE
1	A	360	LEU
1	A	371	ARG
1	A	398	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	9	GLN
1	A	25	ASN
1	A	93	GLN
1	A	105	HIS
1	A	114	GLN

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Mol	Chain	Res	Type
1	A	140	ASN
1	A	147	GLN
1	A	156	ASN
1	A	168	GLN
1	A	175	GLN
1	A	179	ASN
1	A	191	HIS
1	A	200	GLN
1	A	286	HIS
1	A	329	HIS
1	A	353	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 carbohydrates 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).

Mol	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z  > 2	Counts	RMSZ	# Z  > 2
2	FMN	A	401	-	32,33,33	1.49	6 (18%)	34,50,50	2.29	9 (26%)

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	FMN	A	401	-	-	0/18/18/18	0/3/3/3

All (6) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	401	FMN	P-O2P	-2.85	1.44	1.54
2	A	401	FMN	C2-N1	-2.69	1.32	1.38
2	A	401	FMN	P-O1P	-2.18	1.43	1.50
2	A	401	FMN	C4A-N5	2.01	1.36	1.33
2	A	401	FMN	C4-C4A	2.93	1.47	1.41
2	A	401	FMN	C4-N3	3.72	1.39	1.33

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
2	A	401	FMN	C4A-C4-N3	-5.46	116.39	123.52
2	A	401	FMN	N3-C2-N1	-3.90	121.12	127.69
2	A	401	FMN	C4-C4A-C10	-3.41	117.76	119.94
2	A	401	FMN	C4A-C10-N10	-3.13	118.25	120.52
2	A	401	FMN	O3P-P-O5'	-2.22	100.25	106.72
2	A	401	FMN	C4-C4A-N5	2.20	121.37	118.70
2	A	401	FMN	O2P-P-O1P	2.77	119.68	110.63
2	A	401	FMN	C4A-N5-C5A	3.52	120.86	116.72
2	A	401	FMN	C4-N3-C2	8.19	121.99	115.16

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
2	A	401	FMN	2	0

## 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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.