



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 28, 2014 – 07:46 AM GMT

PDB ID : 1YUZ  
Title : Partially Reduced State of Nigerythrin  
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Deposited on : 2005-02-14  
Resolution : 1.40 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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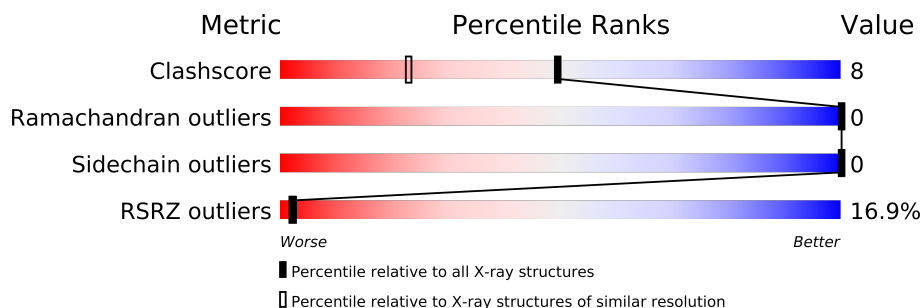
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.40 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	79885	1246 (1.42-1.38)
Ramachandran outliers	78287	1206 (1.42-1.38)
Sidechain outliers	78261	1205 (1.42-1.38)
RSRZ outliers	66119	1097 (1.42-1.38)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	202	
1	B	202	

## 2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 3822 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Nigerythrin.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	200	Total	C	N	O	S	0	5	0
			1580	997	265	308	10			
1	B	202	Total	C	N	O	S	0	10	0
			1636	1032	278	316	10			

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

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

- Molecule 3 is FE (II) ION (three-letter code: FE2) (formula: Fe).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	B	1	Total	Fe	0	0
			1	1		
3	A	1	Total	Fe	0	0
			1	1		

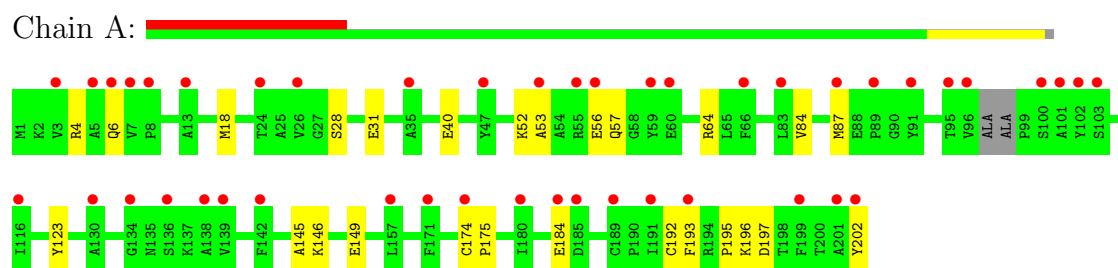
- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	282	Total	O	0	0
			282	282		
4	B	318	Total	O	0	0
			318	318		

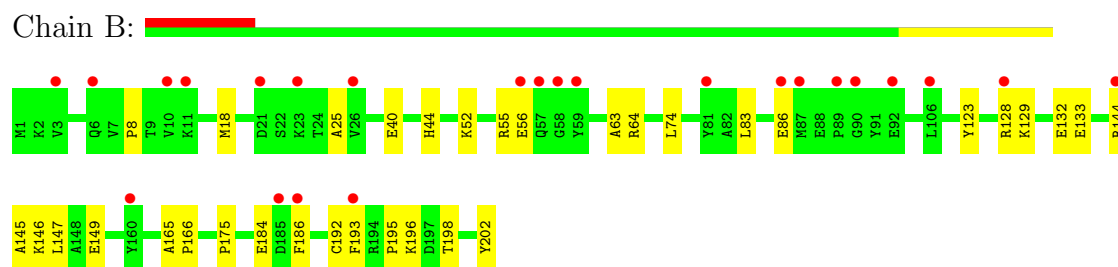
### 3 Residue-property plots

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: Nigerythrin



#### • Molecule 1: Nigerythrin



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	46.67Å 72.31Å 119.02Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	34.60 – 1.40 72.31 – 2.18	Depositor EDS
% Data completeness (in resolution range)	99.7 (34.60-1.40) 92.9 (72.31-2.18)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.07 (at 2.18Å)	Xtriage
Refinement program	CNS 1.0	Depositor
R, $R_{free}$	0.192 , 0.208 0.222 , (Not available)	Depositor DCC
$R_{free}$ test set	No test flags present.	DCC
Wilson B-factor (Å <sup>2</sup> )	8.6	Xtriage
Anisotropy	0.303	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 29.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 39991 reflections	Xtriage
$F_o, F_c$ correlation	0.88	EDS
Total number of atoms	3822	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	15.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.85% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: FE2, FE

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	0.28	0/1613	0.49	0/2180
1	B	0.30	0/1670	0.50	0/2258
All	All	0.29	0/3283	0.50	0/4438

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no planarity outliers.

### 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	1580	0	1524	20	0
1	B	1636	0	1580	28	0
2	A	2	0	0	0	0
2	B	2	0	0	0	0
3	A	1	0	0	0	0
3	B	1	0	0	0	0
4	A	282	0	0	4	0
4	B	318	0	0	3	0
All	All	3822	0	3104	48	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 8.

All (48) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:18:MET:HB3	1:B:128[B]:ARG:NH1	1.73	1.04
1:B:166:PRO:HG3	4:B:327:HOH:O	1.82	0.79
1:B:175:PRO:HG3	1:B:198:THR:O	1.87	0.75
1:A:64:ARG:HD2	1:A:202:TYR:OXT	1.87	0.75
1:B:44:HIS:CD2	1:B:74[B]:LEU:HG	2.21	0.74
1:A:84:VAL:HA	1:A:87:MET:CE	2.26	0.66
1:B:145:ALA:O	1:B:149:GLU:HG2	1.97	0.65
1:A:145:ALA:O	1:A:149:GLU:HG2	2.01	0.60
1:A:195:PRO:HB2	1:A:197:ASP:OD1	2.03	0.58
1:B:25:ALA:HB2	4:B:592:HOH:O	2.05	0.57
1:B:144[A]:ARG:HG2	1:B:144[A]:ARG:HH11	1.71	0.56
1:A:28:SER:OG	1:A:31:GLU:HG3	2.04	0.56
1:B:195:PRO:HD3	4:B:527:HOH:O	2.06	0.54
1:B:165:ALA:N	1:B:166:PRO:HD3	2.23	0.54
1:A:84:VAL:HA	1:A:87:MET:HE3	1.88	0.54
1:A:84:VAL:HA	1:A:87:MET:HE2	1.90	0.53
1:A:52:LYS:O	1:A:56:GLU:HG3	2.08	0.53
1:B:18:MET:HB3	1:B:128[B]:ARG:HH12	1.69	0.52
1:A:53:ALA:O	1:A:57:GLN:HG3	2.10	0.52
1:B:52:LYS:O	1:B:56[A]:GLU:HG3	2.11	0.51
1:B:128[B]:ARG:O	1:B:132:GLU:HG3	2.10	0.51
1:A:64:ARG:NH2	4:A:404:HOH:O	2.31	0.50
1:A:18[B]:MET:HE3	4:A:416:HOH:O	2.12	0.50
1:A:174:CYS:SG	1:A:175:PRO:HD2	2.51	0.50
1:B:55:ARG:HG3	1:B:63:ALA:CB	2.43	0.49
1:B:186:PHE:CZ	1:B:196[A]:LYS:HG3	2.48	0.48
1:B:128[A]:ARG:O	1:B:132:GLU:HG3	2.13	0.47
1:A:18[B]:MET:HE1	4:A:465:HOH:O	2.14	0.47
1:A:123:TYR:HB3	1:A:146:LYS:HB2	1.97	0.47
1:B:144[A]:ARG:NH1	1:B:144[A]:ARG:HG2	2.30	0.46
1:A:40:GLU:OE1	1:A:40:GLU:HA	2.16	0.46
1:B:83:LEU:O	1:B:86[B]:GLU:HG2	2.15	0.46
1:B:129:LYS:HE3	1:B:133:GLU:OE2	2.16	0.46
1:B:18:MET:H	1:B:128[B]:ARG:NH2	2.14	0.45
1:A:192:CYS:O	1:A:193:PHE:HB2	2.15	0.45
1:B:8:PRO:HG2	1:B:147:LEU:HD13	1.97	0.45
1:B:64:ARG:HD2	1:B:202:TYR:OXT	2.16	0.45
1:B:55:ARG:HG3	1:B:63:ALA:HB1	1.98	0.45
1:A:184[A]:GLU:HB3	1:A:196:LYS:NZ	2.32	0.44

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Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:B:40:GLU:HA	1:B:40:GLU:OE1	2.18	0.44
1:B:123:TYR:HB3	1:B:146:LYS:HB2	2.00	0.43
1:B:192:CYS:O	1:B:193:PHE:HB2	2.18	0.43
1:A:87:MET:HE3	1:A:87:MET:HB2	1.70	0.43
1:A:195:PRO:HD3	4:A:501:HOH:O	2.19	0.42
1:B:184:GLU:HB3	1:B:196[A]:LYS:NZ	2.34	0.42
1:B:184:GLU:HB3	1:B:196[A]:LYS:HZ1	1.83	0.41
1:A:4:ARG:HB3	1:A:6:GLN:HE21	1.85	0.41
1:B:44:HIS:HD2	1:B:74[B]:LEU:HG	1.81	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	201/202 (100%)	194 (96%)	7 (4%)	0	100	100
1	B	210/202 (104%)	204 (97%)	6 (3%)	0	100	100
All	All	411/404 (102%)	398 (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	161/157 (102%)	161 (100%)	0	100	100
1	B	165/157 (105%)	165 (100%)	0	100	100
All	All	326/314 (104%)	326 (100%)	0	100	100



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

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

Mol	Chain	Res	Type
1	A	6	GLN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 6 ligands modelled in this entry, 6 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.

## 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	200/202 (99%)	1.48	45 (22%)	1 1	8, 13, 20, 22	2 (1%)
1	B	202/202 (100%)	1.23	24 (11%)	5 4	7, 11, 18, 22	2 (0%)
All	All	402/404 (99%)	1.35	69 (17%)	2 2	7, 12, 19, 22	4 (0%)

All (69) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	100	SER	7.0
1	A	96	VAL	5.4
1	B	87[A]	MET	4.7
1	A	7	VAL	4.5
1	B	3	VAL	4.1
1	A	202	TYR	4.1
1	A	185[A]	ASP	3.9
1	A	95	THR	3.7
1	B	81	TYR	3.5
1	A	102	TYR	3.5
1	A	103[A]	SER	3.5
1	A	191	ILE	3.4
1	A	184[A]	GLU	3.3
1	B	86[A]	GLU	3.3
1	B	92	GLU	3.3
1	B	23	LYS	3.3
1	A	5	ALA	3.2
1	B	58	GLY	3.1
1	A	53	ALA	3.1
1	B	56[A]	GLU	2.9
1	A	60	GLU	2.8
1	B	128[A]	ARG	2.8
1	B	21[A]	ASP	2.8
1	A	83	LEU	2.8

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Mol	Chain	Res	Type	RSRZ
1	A	139	VAL	2.8
1	B	90	GLY	2.7
1	B	89	PRO	2.7
1	A	24	THR	2.7
1	A	55	ARG	2.7
1	A	101	ALA	2.7
1	A	6	GLN	2.6
1	A	89	PRO	2.6
1	A	47	TYR	2.6
1	A	193	PHE	2.6
1	A	8	PRO	2.6
1	A	199	PHE	2.5
1	A	174	CYS	2.5
1	A	130	ALA	2.5
1	A	157	LEU	2.5
1	A	171	PHE	2.4
1	B	57	GLN	2.4
1	B	10[A]	VAL	2.4
1	A	35	ALA	2.4
1	B	6	GLN	2.3
1	A	66	PHE	2.3
1	A	142	PHE	2.3
1	A	59	TYR	2.3
1	A	189	CYS	2.2
1	B	186	PHE	2.2
1	B	59	TYR	2.2
1	B	11[A]	LYS	2.2
1	B	185	ASP	2.2
1	A	134	GLY	2.2
1	A	91	TYR	2.2
1	B	144[A]	ARG	2.1
1	B	160	TYR	2.1
1	A	26	VAL	2.1
1	A	180	ILE	2.1
1	A	87	MET	2.1
1	A	13	ALA	2.1
1	A	3	VAL	2.0
1	B	26	VAL	2.0
1	A	136	SER	2.0
1	B	193	PHE	2.0
1	A	138	ALA	2.0
1	B	106	LEU	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	116	ILE	2.0
1	A	56	GLU	2.0
1	A	201	ALA	2.0

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

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

## 6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

## 6.4 Ligands ⓘ

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	RSR	LLDF	B-factors( $\text{\AA}^2$ )	Q<0.9
2	FE	B	303	1/1	0.08	-1.71	8,8,8,8	0
2	FE	A	301	1/1	0.11	-1.89	11,11,11,11	0
2	FE	A	303	1/1	0.07	-2.71	11,11,11,11	0
3	FE2	B	302	1/1	0.07	-3.68	8,8,8,8	0
2	FE	B	301	1/1	0.04	-4.36	10,10,10,10	0
3	FE2	A	302	1/1	0.06	-4.45	8,8,8,8	0

## 6.5 Other polymers ⓘ

There are no such residues in this entry.