

wwPDB X-ray Structure Validation Summary Report (i)

Dec 5, 2023 - 03:57 am GMT

PDB ID : 2BWD

Title : Atomic Resolution Structure of Achromobacter cycloclastes Cu Nitrite Reduc-

tase with Endogenously bound Nitrite and NO

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Deposited on : 2005-07-13

Resolution : 1.15 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 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:

 $Mol Probity \quad : \quad 4.02b\text{-}467$

Mogul : 1.8.4, CSD as541be (2020)

Xtriage (Phenix) : 1.13 EDS : 2.36

Percentile statistics : 20191225.v01 (using entries in the PDB archive December 25th 2019)

Refmac : 5.8.0158

CCP4 : 7.0.044 (Gargrove)

Ideal geometry (proteins) : Engh & Huber (2001) Ideal geometry (DNA, RNA) : Parkinson et al. (1996)

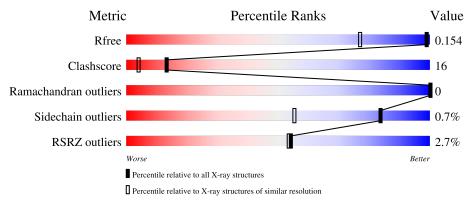
Validation Pipeline (wwPDB-VP) : 2.36

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 1.15 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\AA)}) \end{array}$
R_{free}	130704	1492 (1.18-1.10)
Clashscore	141614	1537 (1.18-1.10)
Ramachandran outliers	138981	1483 (1.18-1.10)
Sidechain outliers	138945	1480 (1.18-1.10)
RSRZ outliers	127900	1464 (1.18-1.10)

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					
			3%				
1	A	340	84%	14%	• •		

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ACT	A	504	-	-	X	-





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Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
4	ACT	A	508	-	X	=	-



2 Entry composition (i)

There are 7 unique types of molecules in this entry. The entry contains 3416 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.

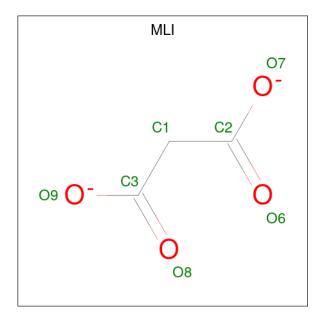
• Molecule 1 is a protein called COPPER-CONTAINING NITRITE REDUCTASE.

Mol	Chain	Residues		At	oms			ZeroOcc	AltConf	Trace
1	Λ	334	Total	С	N	О	S	0	38	0
1	A	334	2822	1810	470	530	12	U	30	

• Molecule 2 is COPPER (II) ION (three-letter code: CU) (formula: Cu).

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

• Molecule 3 is MALONATE ION (three-letter code: MLI) (formula: $C_3H_2O_4$).



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

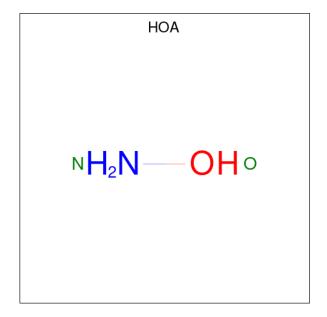
• Molecule 4 is ACETATE ION (three-letter code: ACT) (formula: $C_2H_3O_2$).





Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	1	Total C O 4 2 2	0	0
4	A	1	Total C O 8 4 4	0	1
4	A	1	Total C O 4 2 2	0	0

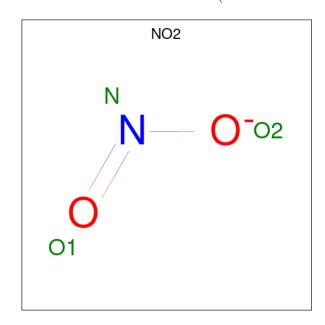
 \bullet Molecule 5 is HYDROXYAMINE (three-letter code: HOA) (formula: ${\rm H_3NO}).$



Mo	l Chai	n	Residues	Atoms			ZeroOcc	AltConf
5	A		1	Total 2	N 1	O 1	0	0



• Molecule 6 is NITRITE ION (three-letter code: NO2) (formula: NO₂).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
6	A	1	Total 3	N 1	O 2	0	0

• Molecule 7 is water.

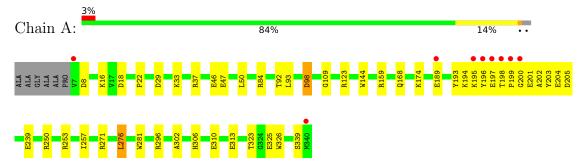
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
7	A	564	Total O 564 564	1	0



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: COPPER-CONTAINING NITRITE REDUCTASE





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 3	Depositor
Cell constants	96.24Å 96.24Å 96.24Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	30.00 - 1.15	Depositor
resolution (A)	15.03 - 1.15	EDS
% Data completeness	97.1 (30.00-1.15)	Depositor
(in resolution range)	97.2 (15.03-1.15)	EDS
R_{merge}	0.06	Depositor
R_{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	1.71 (at 1.15Å)	Xtriage
Refinement program	REFMAC 5.1.9999	Depositor
Ρ. Р.	0.113 , 0.144	Depositor
R, R_{free}	0.130 , 0.154	DCC
R_{free} test set	5072 reflections $(4.97%)$	wwPDB-VP
Wilson B-factor (Å ²)	8.2	Xtriage
Anisotropy	0.000	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.34, 53.6	EDS
L-test for twinning ²	$< L > = 0.51, < L^2> = 0.35$	Xtriage
Estimated twinning fraction	0.028 for l,-k,h	Xtriage
F_o, F_c correlation	0.98	EDS
Total number of atoms	3416	wwPDB-VP
Average B, all atoms (Å ²)	12.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.90% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical values of <|L|>, $<L^2>$ for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



¹Intensities 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: MLI, NO2, CU, ACT, HOA

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	Mol Chain		nd lengths	Bond angles		
Moi Cha	Chain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	A	0.92	$1/2964 \ (0.0\%)$	1.03	17/4035 (0.4%)	

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 maintain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	1

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(A)	$\operatorname{Ideal}(\text{\AA})$
1	A	310	GLU	CD-OE2	6.31	1.32	1.25

The worst 5 of 17 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$Observed(^o)$	$\operatorname{Ideal}({}^{o})$
1	A	37	ARG	NE-CZ-NH2	10.38	125.49	120.30
1	A	8	ASP	CB-CG-OD2	8.84	126.25	118.30
1	A	98[A]	ASP	CB-CG-OD1	8.54	125.98	118.30
1	A	98[B]	ASP	CB-CG-OD1	8.54	125.98	118.30
1	A	37	ARG	NE-CZ-NH1	-8.17	116.22	120.30

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	306	HIS	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	2822	0	2779	89	0
2	A	2	0	0	0	0
3	A	7	0	2	0	0
4	A	16	0	12	2	0
5	A	2	0	0	0	0
6	A	3	0	0	1	0
7	A	564	0	0	80	2
All	All	3416	0	2793	92	2

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 16.

The worst 5 of 92 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:A:202[B]:ALA:HB3	7:A:2379:HOH:O	1.22	1.30
1:A:200[A]:GLY:HA2	7:A:2393:HOH:O	1.25	1.29
1:A:16[B]:LYS:CE	7:A:2028:HOH:O	1.64	1.27
1:A:325[B]:GLU:CG	7:A:2528:HOH:O	1.71	1.26
1:A:325[A]:GLU:HG2	7:A:2528:HOH:O	1.34	1.23

All (2) 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	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	Clash overlap (Å)
7:A:2110:HOH:O	7:A:2440:HOH:O[3_555]	1.80	0.40
7:A:2391:HOH:O	7:A:2508:HOH:O[9_555]	2.14	0.06



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	Analysed Favoured Allowed		Outliers	Percentiles
1	A	368/340 (108%)	360 (98%)	8 (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	Analysed Rotameric		Percentiles	
1	A	306/273 (112%)	303 (99%)	3 (1%)	76	42

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	276[A]	LEU
1	A	276[B]	LEU
1	A	339	SER

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

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 9 ligands modelled in this entry, 2 are 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	Tuna	Chain	Peg	Res Link		ond len	gths	I	Bond an	gles
IVIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2
3	MLI	A	503	-	6,6,6	2.08	1 (16%)	7,7,7	1.38	1 (14%)
6	NO2	A	507	2,5	1,2,2	5.64	1 (100%)	0,1,1	_	-
4	ACT	A	505[A]	-	3,3,3	0.90	0	3,3,3	0.08	0
4	ACT	A	505[B]	-	3,3,3	0.83	0	3,3,3	0.79	0
4	ACT	A	504	-	3,3,3	0.38	0	3,3,3	2.42	2 (66%)
5	HOA	A	506	2,6	0,1,1	-	-	-		
4	ACT	A	508	-	3,3,3	2.71	1 (33%)	3,3,3	4.27	3 (100%)

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.

\mathbf{Mol}	Type	Chain	Res	Link	Chirals	Torsions	Rings
3	MLI	A	503	-	-	0/4/4/4	-

All (3) bond length outliers are listed below:

Mo	ol	Chain	Res	Type	Atoms	\mathbf{Z}	Observed(A)	$\operatorname{Ideal}(ext{\AA})$
6		A	507	NO2	O1-N	5.64	1.50	1.22
4		A	508	ACT	O-C	4.53	1.43	1.22

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Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\operatorname{Observed}(\text{\AA})$	Ideal(A)
3	A	503	MLI	C1-C3	-4.49	1.44	1.51

The worst 5 of 6 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^{o})$	$\operatorname{Ideal}({}^{o})$
4	A	508	ACT	OXT-C-O	6.22	144.98	122.05
4	A	504	ACT	O-C-CH3	-3.36	109.26	122.33
4	A	508	ACT	OXT-C-CH3	-3.18	102.06	115.18
4	A	508	ACT	O-C-CH3	-2.44	112.83	122.33
3	A	503	MLI	O9-C3-C1	2.37	122.11	114.54

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
6	A	507	NO2	1	0
4	A	504	ACT	2	0

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	$\langle { m RSRZ} \rangle$	# RSRZ > 2		$OWAB(A^2)$	Q < 0.9
1	A	334/340 (98%)	-0.19	9 (2%) 54	53	5, 8, 16, 28	20 (5%)

The worst 5 of 9 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	197[A]	GLU	5.6
1	A	195[A]	LYS	4.3
1	A	340[A]	MET	4.1
1	A	196[A]	TYR	3.8
1	A	7[A]	VAL	3.8

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

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

6.3 Carbohydrates (i)

There are no monosaccharides in this entry.

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	$oxed{f B-factors({ m \AA}^2)}$	Q<0.9
4	ACT	A	504	4/4	0.90	0.15	9,13,14,16	4
4	ACT	A	505[A]	4/4	0.94	0.11	9,9,10,10	4

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Mol	Type	Chain	Res	Atoms	RSCC	RSR	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
4	ACT	A	505[B]	4/4	0.94	0.11	8,10,11,13	4
6	NO2	A	507	3/3	0.96	0.10	8,8,12,13	3
5	HOA	A	506	2/2	0.97	0.11	6,6,6,10	2
4	ACT	A	508	4/4	0.97	0.15	7,14,16,18	0
3	MLI	A	503	7/7	0.99	0.05	6,6,7,9	7
2	CU	A	502	1/1	1.00	0.02	6,6,6,6	1
2	CU	A	501	1/1	1.00	0.02	7,7,7,7	0

6.5 Other polymers (i)

There are no such residues in this entry.

