

# Full wwPDB X-ray Structure Validation Report (i)

#### Mar 10, 2024 – 04:30 AM EDT

PDB ID	:	3X2Z
Title	:	Crystal structure of metallo-beta-lactamase in complex with nickel from
		Thermotoga maritima
Authors	:	Choi, H.J.; Kim, H.J.; Matsuura, A.; Mikami, B.; Yoon, H.J.; Lee, H.H.
Deposited on		
Resolution	:	2.33  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 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:

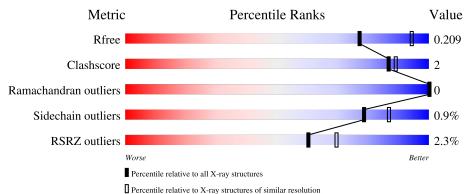
MolProbity	:	4.02b-467
Xtriage (Phenix)	:	1.13
EDS	:	2.36
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
CCP4	:	7.0.044 (Gargrove)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.36

# 1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure:  $X\text{-}RAY \, DIFFRACTION$ 

The reported resolution of this entry is 2.33 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\textbf{Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries,\ resolution\ range}({ m \AA}))$
R <sub>free</sub>	130704	2096 (2.36-2.32)
Clashscore	141614	2193 (2.36-2.32)
Ramachandran outliers	138981	2159 (2.36-2.32)
Sidechain outliers	138945	2160 (2.36-2.32)
RSRZ outliers	127900	2067 (2.36-2.32)

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	Quality of chain	
1	А	227	2% 98%	•
1	В	227	3% 93%	7%
1	С	227	2% 98%	•



# 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 5534 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.

Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	Trace	
1	Δ	227	Total	С	Ν	0	$\mathbf{S}$	0	0	0
	1 A	221	1735	1121	284	319	11			
1	1 B	227	Total	С	Ν	0	S	0	0	0
		221	1735	1121	284	319	11			
1	C	227	Total	С	Ν	0	S	0	0	0
	227	1735	1121	284	319	11	0	0	U	

• Molecule 1 is a protein called UPF0173 metal-dependent hydrolase TM\_1162.

There are 3 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP Q9X0P5
В	0	GLY	-	expression tag	UNP Q9X0P5
С	0	GLY	-	expression tag	UNP Q9X0P5

• Molecule 2 is NICKEL (II) ION (three-letter code: NI) (formula: Ni).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	А	2	Total Ni 2 2	0	0
2	В	2	Total Ni 2 2	0	0
2	С	2	Total Ni 2 2	0	0

• Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	108	Total O 108 108	0	0
3	В	112	Total         O           112         112	0	0

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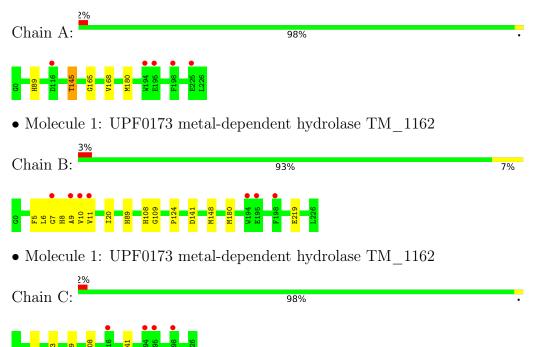
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	С	103	Total         O           103         103	0	0



# 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 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: UPF0173 metal-dependent hydrolase TM\_1162





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 41 21 2	Depositor
Cell constants	$\begin{array}{cccccccccccccccccccccccccccccccccccc$	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	$\frac{90.00}{48.07} - 2.33$	Dan asitan
Resolution (Å)	48.07 - 2.33 35.16 - 2.33	Depositor EDS
% Data completeness	96.7 (48.07-2.33)	Depositor
(in resolution range)	96.7 (35.16-2.33)	EDS
R <sub>merge</sub>	(Not available)	Depositor
R <sub>sym</sub>	(Not available)	Depositor
$\frac{R_{sym}}{\langle I/\sigma(I) \rangle^{-1}}$	$8.65 (at 2.34 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.8.0049	Depositor
D D .	0.191 , $0.205$	Depositor
$R, R_{free}$	0.194 , $0.209$	DCC
$R_{free}$ test set	2009 reflections $(3.11\%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	31.6	Xtriage
Anisotropy	0.005	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.35 , $36.4$	EDS
L-test for twinning <sup>2</sup>	$< L >=0.49, < L^2>=0.33$	Xtriage
Estimated twinning fraction	0.015 for -h,l,k	Xtriage
	0.012 for -l,-k,-h	Ŭ
$F_o, F_c$ correlation	0.95	EDS
Total number of atoms	5534	wwPDB-VP
Average B, all atoms $(Å^2)$	34.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 2.59% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical values of  $\langle |L| \rangle$ ,  $\langle L^2 \rangle$  for acentric reflections are 0.5, 0.333 respectively for untwinned datasets, and 0.375, 0.2 for perfectly twinned datasets.



<sup>&</sup>lt;sup>1</sup>Intensities 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: NI

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	А	0.27	0/1772	0.47	0/2393	
1	В	0.27	0/1772	0.47	0/2393	
1	С	0.27	0/1772	0.47	0/2393	
All	All	0.27	0/5316	0.47	0/7179	

There are no bond length outliers.

There are no bond angle outliers.

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	А	1735	0	1764	2	0
1	В	1735	0	1764	18	0
1	С	1735	0	1764	2	0
2	А	2	0	0	0	0
2	В	2	0	0	0	0
2	С	2	0	0	0	0
3	А	108	0	0	0	0
3	В	112	0	0	1	0
3	С	103	0	0	0	0
All	All	5534	0	5292	22	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 2.

All $(22)$ close contacts	within the	same	asymmetric	unit	are	listed	below,	sorted	by '	their	$\operatorname{clash}$
magnitude.											

Atom-1	Atom-2	Interatomic	Clash
Atom-1	Atom-2	distance (Å)	overlap (Å)
1:B:6:LEU:HB2	1:B:10:VAL:CG1	2.21	0.70
1:A:145:THR:O	1:A:180:MET:HE1	1.97	0.64
1:B:9:ALA:HB3	3:B:407:HOH:O	2.02	0.59
1:B:6:LEU:HD12	1:B:10:VAL:HG11	1.84	0.59
1:B:11:VAL:O	1:B:20:ILE:HG13	2.03	0.58
1:B:8:HIS:CD2	1:B:9:ALA:H	2.22	0.58
1:B:5:PHE:CE2	1:B:7:GLY:HA2	2.42	0.55
1:B:148:MET:HB2	1:B:180:MET:CE	2.37	0.54
1:B:108:HIS:CE1	1:B:141:ASP:HB3	2.44	0.53
1:B:108:HIS:HB3	1:B:124:PRO:HG3	1.92	0.50
1:B:5:PHE:CZ	1:B:7:GLY:HA2	2.47	0.50
1:B:6:LEU:HB2	1:B:10:VAL:HB	1.95	0.48
1:B:180:MET:HB3	1:B:180:MET:HE2	1.88	0.46
1:B:6:LEU:HD12	1:B:10:VAL:CG1	2.45	0.45
1:B:6:LEU:HB2	1:B:10:VAL:CB	2.45	0.45
1:B:8:HIS:CG	1:B:9:ALA:H	2.36	0.44
1:B:108:HIS:HB3	1:B:124:PRO:CG	2.48	0.44
1:B:6:LEU:HB2	1:B:10:VAL:HG12	1.98	0.43
1:C:9:ALA:CB	1:C:53:HIS:CE1	3.03	0.42
1:A:165:GLY:HA2	1:A:168:VAL:O	2.19	0.42
1:B:108:HIS:CG	1:B:109:GLY:H	2.37	0.42
1:C:108:HIS:CE1	1:C:141:ASP:HB3	2.55	0.41

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	Perce	ntiles
1	А	225/227~(99%)	214 (95%)	11 (5%)	0	100	100
1	В	225/227~(99%)	216 (96%)	9~(4%)	0	100	100
1	С	225/227~(99%)	214 (95%)	11 (5%)	0	100	100
All	All	675/681~(99%)	644 (95%)	31 (5%)	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 side chain 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	А	187/187~(100%)	185~(99%)	2(1%)	73 83
1	В	187/187~(100%)	185~(99%)	2(1%)	73 83
1	С	$187/187\ (100\%)$	186 (100%)	1 (0%)	88 93
All	All	561/561~(100%)	556~(99%)	5 (1%)	78 87

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	А	89	HIS
1	А	145	THR
1	В	89	HIS
1	В	219	GLU
1	С	89	HIS

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

No monomer is involved in short contacts.

# 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	<RSRZ $>$	# RSRZ > 2	$OWAB(Å^2)$	Q<0.9
1	А	227/227~(100%)	-0.19	5 (2%) 62 71	20, 31, 47, 59	0
1	В	$227/227 \ (100\%)$	-0.08	7 (3%) 49 59	21, 32, 49, 59	0
1	С	$227/227 \ (100\%)$	-0.17	4 (1%) 68 76	22, 35, 51, 62	0
All	All	681/681~(100%)	-0.15	16 (2%) 60 69	20, 33, 50, 62	0

All (16) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	11	VAL	4.3
1	А	198	PHE	3.8
1	В	198	PHE	3.7
1	С	195	GLU	3.0
1	А	194	TRP	2.8
1	А	195	GLU	2.6
1	С	198	PHE	2.4
1	В	7	GLY	2.4
1	А	225	GLU	2.4
1	В	194	TRP	2.3
1	А	116	ASP	2.2
1	С	116	ASP	2.1
1	С	194	TRP	2.1
1	В	10	VAL	2.1
1	В	9	ALA	2.0
1	В	195	GLU	2.0

# 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	B-factors(Å <sup>2</sup> )	Q<0.9
2	NI	В	302	1/1	0.73	0.13	43,43,43,43	0
2	NI	С	302	1/1	0.85	0.11	49,49,49,49	0
2	NI	А	302	1/1	0.88	0.14	44,44,44,44	0
2	NI	А	301	1/1	0.96	0.17	28,28,28,28	0
2	NI	С	301	1/1	0.97	0.17	28,28,28,28	0
2	NI	В	301	1/1	0.98	0.19	27,27,27,27	0

### 6.5 Other polymers (i)

There are no such residues in this entry.

