

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

#### May 28, 2020 – 09:58 pm BST

PDB ID : 2GZX

Title: Crystal Structure of the TatD deoxyribonuclease MW0446 from Staphylococ-

cus aureus. Northeast Structural Genomics Consortium Target ZR237.

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nomics Consortium (NESG)

Deposited on : 2006-05-12

Resolution : 2.20 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) : 1.13 EDS : 2.11

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)

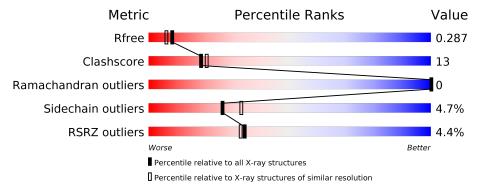
Validation Pipeline (wwPDB-VP) : 2.11

## 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 2.20 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	$\begin{array}{c} \textbf{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
$R_{free}$	130704	4898 (2.20-2.20)
Clashscore	141614	5594 (2.20-2.20)
Ramachandran outliers	138981	5503 (2.20-2.20)
Sidechain outliers	138945	5504 (2.20-2.20)
RSRZ outliers	127900	4800 (2.20-2.20)

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 >=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	A	265	73%	21%	• 5%		
1	В	265	66%	24%	5% 5%		



## 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4104 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 putative TatD related DNase.

Mol	Chain	Residues		Atoms				ZeroOcc	AltConf	Trace	
1	A	253		C 1283					0	0	0
1	В	253	Total 2004	C 1276					0	0	0

There are 28 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	1	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
A	33	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
A	47	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
A	93	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
A	152	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
A	196	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
A	258	LEU	-	EXPRESSION TAG	UNP Q1Y6Z0
A	259	GLU	-	EXPRESSION TAG	UNP Q1Y6Z0
A	260	HIS	-	EXPRESSION TAG	UNP Q1Y6Z0
A	261	HIS	=	EXPRESSION TAG	UNP Q1Y6Z0
A	262	HIS	-	EXPRESSION TAG	UNP Q1Y6Z0
A	263	HIS	-	EXPRESSION TAG	UNP Q1Y6Z0
A	264	HIS	=	EXPRESSION TAG	UNP Q1Y6Z0
A	265	HIS	=	EXPRESSION TAG	UNP Q1Y6Z0
В	1	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
В	33	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
В	47	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
В	93	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
В	152	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
В	196	MSE	MET	MODIFIED RESIDUE	UNP Q1Y6Z0
В	258	LEU	-	EXPRESSION TAG	UNP Q1Y6Z0
В	259	GLU	-	EXPRESSION TAG	UNP Q1Y6Z0
В	260	HIS	-	EXPRESSION TAG	UNP Q1Y6Z0
В	261	HIS	ı	EXPRESSION TAG	UNP Q1Y6Z0
В	262	HIS	-	EXPRESSION TAG	UNP Q1Y6Z0

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Chain	Residue	Modelled	Actual	${f Comment}$	Reference
В	263	HIS	_	EXPRESSION TAG	UNP Q1Y6Z0
В	264	HIS	_	EXPRESSION TAG	UNP Q1Y6Z0
В	265	HIS	-	EXPRESSION TAG	UNP Q1Y6Z0

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

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	В	2	Total Ni 2 2	0	0
2	A	2	Total Ni 2 2	0	0

• Molecule 3 is water.

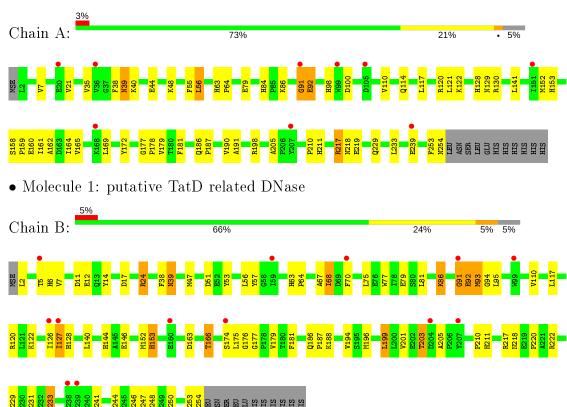
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	41	Total O 41 41	0	0
3	В	40	Total O 40 40	0	0



# 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 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: putative TatD related DNase





# 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants	46.68Å 77.32Å 76.13Å	Depositor
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 98.01° 90.00°	Depositor
Resolution (Å)	26.79 - 2.20	Depositor
resolution (A)	29.66 - 2.20	EDS
% Data completeness	95.9 (26.79-2.20)	Depositor
(in resolution range)	98.3 (29.66-2.20)	EDS
$R_{merge}$	0.10	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	3.06  (at  2.20Å)	Xtriage
Refinement program	CNS 1.1	Depositor
$R, R_{free}$	0.258 , $0.286$	Depositor
it, it free	0.262 , $0.287$	DCC
$R_{free}$ test set	2103 reflections $(3.93\%)$	wwPDB-VP
Wilson B-factor $(\mathring{A}^2)$	24.4	Xtriage
Anisotropy	0.294	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.33 , 34.0	EDS
L-test for twinning <sup>2</sup>	$ < L >=0.49, < L^2>=0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o$ , $F_c$ correlation	0.91	EDS
Total number of atoms	4104	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	29.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 6.95% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical 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.



<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:

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	Chain	Boı	nd lengths	Bond angles		
Mol Chain		RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.59	$2/2054 \ (0.1\%)$	0.77	7/2774~(0.3%)	
1	В	0.52	$2/2043 \ (0.1\%)$	0.74	7/2762~(0.3%)	
All	All	0.56	4/4097 (0.1%)	0.75	$14/5536 \ (0.3\%)$	

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	1
1	В	0	1
All	All	0	2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	${ m Observed}({ m \AA})$	$\mathbf{Ideal}(\mathbf{\AA})$
1	A	92	GLU	CB-CG	-15.30	1.23	1.52
1	В	92	GLU	CB-CG	-9.33	1.34	1.52
1	В	93	MSE	SE-CE	-6.25	1.58	1.95
1	A	91	GLY	CA-C	-5.98	1.42	1.51

The worst 5 of 14 bond angle outliers are listed below:

Mol	Chain	Res	Type	${f Atoms}$	$\mathbf{Z}$	$Observed(^o)$	$\operatorname{Ideal}({}^o)$
1	A	92	GLU	CA-CB-CG	-11.83	87.38	113.40
1	A	91	GLY	CA-C-N	-10.38	94.38	117.20
1	В	91	GLY	CA-C-N	-10.17	94.83	117.20
1	A	92	GLU	N-CA-CB	9.59	127.86	110.60
1	В	92	GLU	CA-CB-CG	-8.90	93.82	113.40



There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	91	GLY	Mainchain
1	В	91	GLY	Mainchain

### 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	2015	0	1977	43	0
1	В	2004	0	1953	63	0
2	A	2	0	0	0	0
2	В	2	0	0	0	0
3	A	41	0	0	3	0
3	В	40	0	0	2	0
All	All	4104	0	3930	105	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 13.

The worst 5 of 105 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$egin{aligned}  ext{Interatomic} \  ext{distance} \ ( ext{Å}) \end{aligned}$	Clash overlap (Å)	
1:B:24:ARG:HG2	1:B:24:ARG:HH11	1.32	0.92	
1:B:47:MSE:HE2	1:B:81:LEU:HD12	1.55	0.89	
1:A:229:GLN:HE21	1:A:233:LEU:HD13	1.44	0.80	
1:B:5:THR:HG22	3:B:370:HOH:O	1.85	0.77	
1:B:38:PHE:HB2	1:B:63:HIS:HB2	1.69	0.75	

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	${f ntiles}$
1	A	251/265~(95%)	243 (97%)	8 (3%)	0	100	100
1	В	251/265~(95%)	241 (96%)	10 (4%)	0	100	100
All	All	$502/530 \; (95\%)$	484 (96%)	18 (4%)	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   Outliers		Percentiles		
1	A	$216/227 \ (95\%)$	210 (97%)	6 (3%)	43 56		
1	В	213/227 (94%)	199 (93%)	14 (7%)	16 19		
All	All	429/454 (94%)	409 (95%)	20 (5%)	26 33		

5 of 20 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	В	51	ASP
1	В	68	ILE
1	В	175	LEU
1	В	24	ARG
1	В	39	ASN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 14 such sidechains are listed below:



Mol	Chain	Res	Type
1	A	248	ASN
1	В	39	ASN
1	В	211	HIS
1	A	244	GLN
1	В	183	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)

Of 4 ligands modelled in this entry, 4 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	$\langle { m RSRZ} \rangle$	$\#\mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q<0.9
1	A	248/265~(93%)	0.44	9 (3%) 42 41	12, 28, 39, 47	0
1	В	248/265~(93%)	0.57	13 (5%) 27 26	16, 30, 42, 49	0
All	All	$496/530 \ (93\%)$	0.50	22 (4%) 34 32	12, 29, 42, 49	0

The worst 5 of 22 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	70	PHE	5.1
1	A	20	GLU	3.4
1	В	91	GLY	3.2
1	В	127	ILE	3.1
1	В	160	GLU	3.0

## 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 carbohydrates 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	${f B-factors}({f A}^2)$	Q<0.9
2	NI	В	298	1/1	0.09	0.37	30,30,30,30	0
2	NI	A	299	1/1	0.43	0.28	22,22,22,22	0
2	NI	A	300	1/1	0.98	0.13	15,15,15,15	0
2	NI	В	297	1/1	0.99	0.13	21,21,21,21	0

# 6.5 Other polymers (i)

There are no such residues in this entry.

