



# wwPDB X-ray Structure Validation Summary Report ⓘ

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PDB ID : 3TOI  
Title : Tailoring Enzyme Stability and Exploiting Stability-Trait Linkage by Iterative Truncation and Optimization  
Authors : Tam, H.K.; Einsle, O.  
Deposited on : 2011-09-05  
Resolution : 1.90 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](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
Xtriage (Phenix)	:	1.9-1692
EDS	:	trunk28620
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc28949

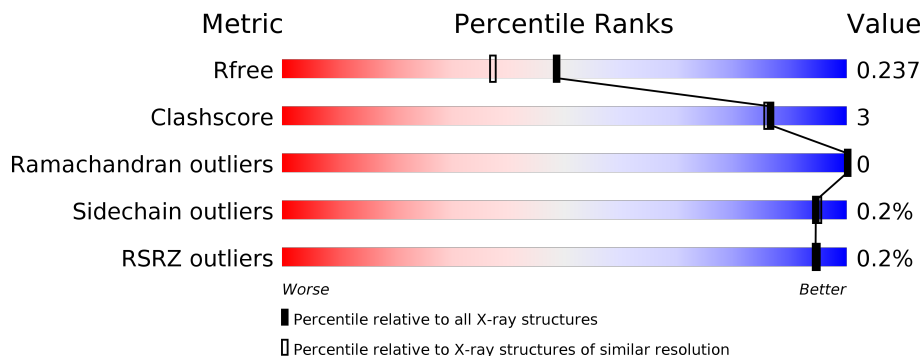
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

## *X-RAY DIFFRACTION*

The reported resolution of this entry is 1.90 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))
$R_{free}$	100719	5047 (1.90-1.90)
Clashscore	112137	5731 (1.90-1.90)
Ramachandran outliers	110173	5669 (1.90-1.90)
Sidechain outliers	110143	5670 (1.90-1.90)
RSRZ outliers	101464	5100 (1.90-1.90)

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.

Mol	Chain	Length	Quality of chain
1	A	254	
1	B	254	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3998 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 Ampicillin resistance protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	0	0	0
			1880	1171	332	365	12			
1	B	245	Total	C	N	O	S	0	0	0
			1868	1165	329	362	12			

There are 34 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	39	ASP	-	EXPRESSION TAG	UNP Q79DR3
A	40	GLY	-	EXPRESSION TAG	UNP Q79DR3
A	56	VAL	ILE	ENGINEERED MUTATION	UNP Q79DR3
A	120	GLY	ARG	ENGINEERED MUTATION	UNP Q79DR3
A	182	THR	MET	ENGINEERED MUTATION	UNP Q79DR3
A	195	SER	THR	ENGINEERED MUTATION	UNP Q79DR3
A	208	MET	ILE	ENGINEERED MUTATION	UNP Q79DR3
A	224	VAL	ALA	ENGINEERED MUTATION	UNP Q79DR3
A	241	HIS	ARG	ENGINEERED MUTATION	UNP Q79DR3
A	265	MET	THR	ENGINEERED MUTATION	UNP Q79DR3
A	288	GLY	-	EXPRESSION TAG	UNP Q79DR3
A	289	GLY	-	EXPRESSION TAG	UNP Q79DR3
A	290	HIS	-	EXPRESSION TAG	UNP Q79DR3
A	291	HIS	-	EXPRESSION TAG	UNP Q79DR3
A	292	HIS	-	EXPRESSION TAG	UNP Q79DR3
A	293	HIS	-	EXPRESSION TAG	UNP Q79DR3
A	294	HIS	-	EXPRESSION TAG	UNP Q79DR3
B	39	ASP	-	EXPRESSION TAG	UNP Q79DR3
B	40	GLY	-	EXPRESSION TAG	UNP Q79DR3
B	56	VAL	ILE	ENGINEERED MUTATION	UNP Q79DR3
B	120	GLY	ARG	ENGINEERED MUTATION	UNP Q79DR3
B	182	THR	MET	ENGINEERED MUTATION	UNP Q79DR3
B	195	SER	THR	ENGINEERED MUTATION	UNP Q79DR3
B	208	MET	ILE	ENGINEERED MUTATION	UNP Q79DR3
B	224	VAL	ALA	ENGINEERED MUTATION	UNP Q79DR3

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Chain	Residue	Modelled	Actual	Comment	Reference
B	241	HIS	ARG	ENGINEERED MUTATION	UNP Q79DR3
B	265	MET	THR	ENGINEERED MUTATION	UNP Q79DR3
B	288	GLY	-	EXPRESSION TAG	UNP Q79DR3
B	289	GLY	-	EXPRESSION TAG	UNP Q79DR3
B	290	HIS	-	EXPRESSION TAG	UNP Q79DR3
B	291	HIS	-	EXPRESSION TAG	UNP Q79DR3
B	292	HIS	-	EXPRESSION TAG	UNP Q79DR3
B	293	HIS	-	EXPRESSION TAG	UNP Q79DR3
B	294	HIS	-	EXPRESSION TAG	UNP Q79DR3

- Molecule 2 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	A	132	Total O 132 132	0	0
2	B	118	Total O 118 118	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: Ampicillin resistance protein

Chain A:  89% 9% .



- Molecule 1: Ampicillin resistance protein

Chain B:  91% 5% .



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	42.00Å 47.52Å 258.83Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	46.74 – 1.90 46.74 – 1.90	Depositor EDS
% Data completeness (in resolution range)	99.5 (46.74-1.90) 99.5 (46.74-1.90)	Depositor EDS
$R_{merge}$	0.08	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	3.57 (at 1.90Å)	Xtriage
Refinement program	REFMAC 5.6.0113	Depositor
R, $R_{free}$	0.199 , 0.237 0.199 , 0.237	Depositor DCC
$R_{free}$ test set	2115 reflections (5.32%)	DCC
Wilson B-factor (Å <sup>2</sup> )	22.7	Xtriage
Anisotropy	0.402	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.38 , 36.1	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.40$ , $\langle L^2 \rangle = 0.22$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.94	EDS
Total number of atoms	3998	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.19% of the height of the origin peak. No significant pseudotranslation is detected.*

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

<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.

## 5 Model quality [i](#)

### 5.1 Standard geometry [i](#)

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	0.46	0/1910	0.58	0/2585
1	B	0.46	2/1898 (0.1%)	0.57	0/2570
All	All	0.46	2/3808 (0.1%)	0.57	0/5155

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	210	TRP	CD2-CE2	5.02	1.47	1.41
1	B	229	TRP	CD2-CE2	5.00	1.47	1.41

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	A	1880	0	1881	12	0
1	B	1868	0	1872	7	0
2	A	132	0	0	2	0
2	B	118	0	0	1	0
All	All	3998	0	3753	19	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 3.

The worst 5 of 19 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:106:SER:HB3	1:A:109:THR:OG1	2.01	0.60
1:B:106:SER:HB3	1:B:109:THR:OG1	2.09	0.52
1:B:83:ARG:HD3	1:B:142:ILE:HG22	1.91	0.52
1:A:171:GLU:OE1	1:A:173:ILE:HD11	2.13	0.47
1:A:175:ASN:ND2	2:A:318:HOH:O	2.48	0.46

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	Percentiles	
1	A	246/254 (97%)	239 (97%)	7 (3%)	0	100	100
1	B	243/254 (96%)	235 (97%)	8 (3%)	0	100	100
All	All	489/508 (96%)	474 (97%)	15 (3%)	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	Rotameric	Outliers	Percentiles	
1	A	201/207 (97%)	200 (100%)	1 (0%)	91	91
1	B	201/207 (97%)	201 (100%)	0	100	100

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Mol	Chain	Analysed	Rotameric	Outliers	Percentiles	
All	All	402/414 (97%)	401 (100%)	1 (0%)	94	95

All (1) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	130	SER

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	154	ASN
1	A	158	HIS
1	A	175	ASN
1	B	100	ASN
1	B	206	GLN

### 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](#)

There are no ligands in this entry.

## 5.7 Other polymers [i](#)

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 [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<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	248/254 (97%)	-0.05	1 (0%) 92 93	14, 22, 31, 40	0
1	B	245/254 (96%)	-0.01	0 100 100	15, 23, 36, 45	0
All	All	493/508 (97%)	-0.03	1 (0%) 94 95	14, 22, 33, 45	0

All (1) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	197	GLU	2.2

### 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](#)

There are no ligands in this entry.

### 6.5 Other polymers [i](#)

There are no such residues in this entry.