



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 17, 2024 – 07:14 PM EST

PDB ID : 3UE3  
Title : Crystal structure of Acinetobacter baumannii PBP3  
Authors : Han, S.  
Deposited on : 2011-10-28  
Resolution : 2.30 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](mailto: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 ⓘ symbol.

The types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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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) : 1.13  
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

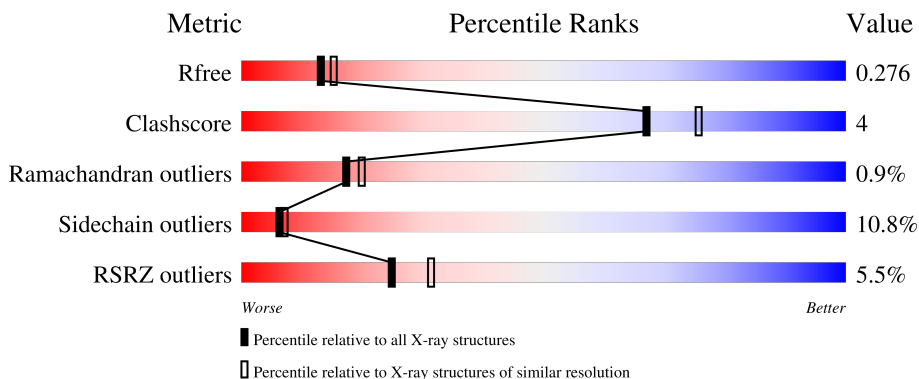
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.30 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))
$R_{free}$	130704	5042 (2.30-2.30)
Clashscore	141614	5643 (2.30-2.30)
Ramachandran outliers	138981	5575 (2.30-2.30)
Sidechain outliers	138945	5575 (2.30-2.30)
RSRZ outliers	127900	4938 (2.30-2.30)

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  $\geq 3$ , 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  $\leq 5\%$ . 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	554	

## 2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 3470 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 Septum formation, penicillin binding protein 3, peptidoglycan synthetase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
			Total	C	N	O	S			
1	A	436	3305	2081	580	628	16	0	0	0

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	56	MET	-	expression tag	UNP C0VN42
A	57	SER	-	expression tag	UNP C0VN42
A	58	HIS	-	expression tag	UNP C0VN42
A	59	HIS	-	expression tag	UNP C0VN42
A	60	HIS	-	expression tag	UNP C0VN42
A	61	HIS	-	expression tag	UNP C0VN42
A	62	HIS	-	expression tag	UNP C0VN42
A	63	HIS	-	expression tag	UNP C0VN42

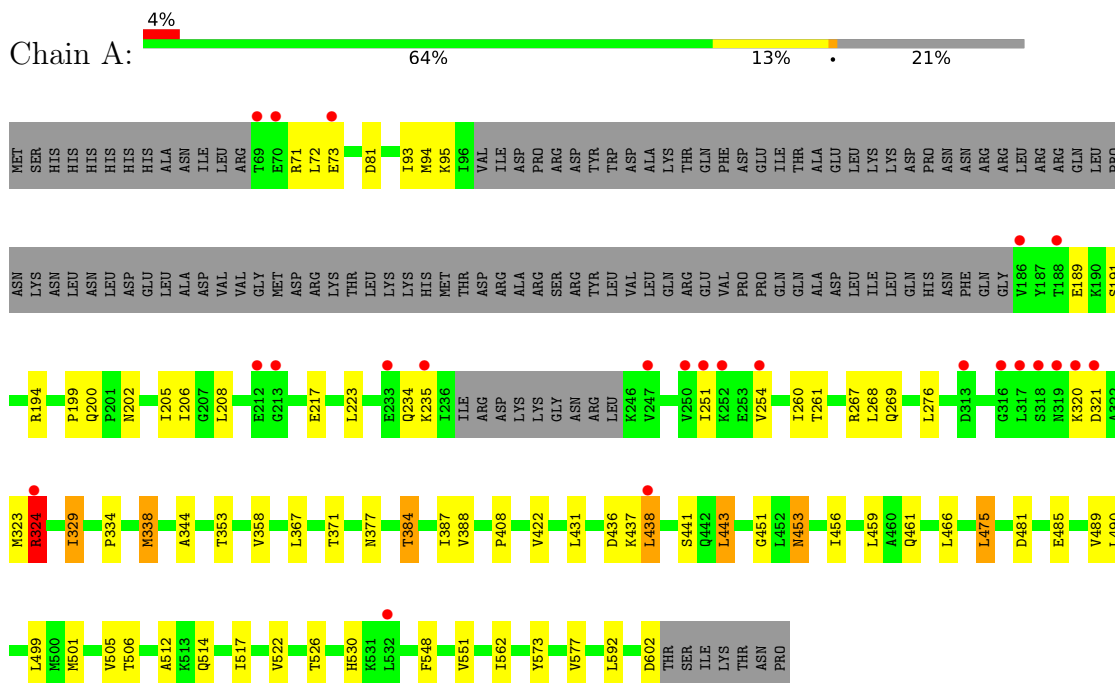
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
			Total	O		
2	A	165	165	165	0	0

### 3 Residue-property plots

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: Septum formation, penicillin binding protein 3, peptidoglycan synthetase



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	39.66Å 89.71Å 211.92Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	38.32 – 2.30 38.32 – 2.30	Depositor EDS
% Data completeness (in resolution range)	97.8 (38.32-2.30) 97.9 (38.32-2.30)	Depositor EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.74 (at 2.29Å)	Xtrriage
Refinement program	BUSTER 2.9.6	Depositor
R, $R_{free}$	0.217 , 0.250 0.236 , 0.276	Depositor DCC
$R_{free}$ test set	1721 reflections (5.07%)	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	25.7	Xtrriage
Anisotropy	0.949	Xtrriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 59.0	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.44$ , $\langle L^2 \rangle = 0.27$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
$F_o, F_c$ correlation	0.89	EDS
Total number of atoms	3470	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	34.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.95% 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	0.50	0/3363	0.78	3/4552 (0.1%)

There are no bond length outliers.

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
1	A	323	MET	C-N-CA	8.32	142.50	121.70
1	A	323	MET	N-CA-C	5.68	126.33	111.00
1	A	437	LYS	C-N-CA	5.40	135.21	121.70

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	3305	0	3353	26	0
2	A	165	0	0	1	0
All	All	3470	0	3353	26	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 4.

All (26) 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:202:ASN:HB3	1:A:206:ILE:HD12	1.83	0.59
1:A:517:ILE:HD12	1:A:522:VAL:HG21	1.83	0.59
1:A:324:ARG:HE	1:A:329:ILE:HD11	1.68	0.58
1:A:384:THR:HB	1:A:501:MET:HG2	1.87	0.56
1:A:408:PRO:HD3	1:A:443:LEU:HD11	1.88	0.55
1:A:344:ALA:HB1	1:A:490:LEU:HD11	1.88	0.55
1:A:431:LEU:HB2	1:A:453:ASN:HB2	1.91	0.53
1:A:436:ASP:C	1:A:438:LEU:H	2.13	0.52
1:A:334:PRO:CG	1:A:338:MET:HG2	2.39	0.52
1:A:235:LYS:HG2	1:A:251:ILE:HD11	1.92	0.50
1:A:367:LEU:HB3	1:A:441:SER:HB2	1.94	0.50
1:A:456:ILE:HG23	1:A:562:ILE:HB	1.95	0.47
1:A:388:VAL:HA	1:A:505:VAL:HG22	1.96	0.47
1:A:223:LEU:HD13	1:A:260:ILE:HD13	1.96	0.47
1:A:526:THR:HG22	1:A:548:PHE:HD2	1.78	0.46
1:A:72:LEU:HD11	1:A:234:GLN:HB2	1.98	0.45
1:A:506:THR:HG21	1:A:522:VAL:HG12	1.98	0.45
1:A:422:VAL:HG23	1:A:475:LEU:HB3	1.99	0.45
1:A:267:ARG:HB3	1:A:592:LEU:HD21	1.99	0.45
1:A:451:GLY:HA3	2:A:722:HOH:O	2.17	0.44
1:A:95:LYS:HB3	1:A:189:GLU:HB2	1.99	0.43
1:A:202:ASN:OD1	1:A:269:GLN:HG2	2.19	0.42
1:A:530:HIS:HE1	1:A:573:TYR:OH	2.01	0.42
1:A:334:PRO:HG2	1:A:338:MET:HG2	2.02	0.41
1:A:505:VAL:HG12	1:A:512:ALA:HB3	2.02	0.40
1:A:387:ILE:HD12	1:A:501:MET:SD	2.62	0.40

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	430/554 (78%)	413 (96%)	13 (3%)	4 (1%)	<b>17</b> <b>20</b>

All (4) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	324	ARG
1	A	320	LYS
1	A	438	LEU
1	A	199	PRO

### 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	352/461 (76%)	314 (89%)	38 (11%)	<b>6</b> <b>7</b>

All (38) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	71	ARG
1	A	73	GLU
1	A	81	ASP
1	A	93	ILE
1	A	94	MET
1	A	191	SER
1	A	194	ARG
1	A	200	GLN
1	A	205	ILE
1	A	208	LEU
1	A	217	GLU
1	A	254	VAL
1	A	261	THR
1	A	268	LEU
1	A	276	LEU
1	A	321	ASP
1	A	324	ARG
1	A	329	ILE
1	A	338	MET
1	A	353	THR
1	A	358	VAL

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Mol	Chain	Res	Type
1	A	371	THR
1	A	377	ASN
1	A	384	THR
1	A	443	LEU
1	A	453	ASN
1	A	459	LEU
1	A	461	GLN
1	A	466	LEU
1	A	475	LEU
1	A	481	ASP
1	A	485	GLU
1	A	489	VAL
1	A	499	LEU
1	A	514	GLN
1	A	551	VAL
1	A	577	VAL
1	A	602	ASP

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (6) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	210	ASN
1	A	234	GLN
1	A	377	ASN
1	A	461	GLN
1	A	514	GLN
1	A	587	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 monosaccharides 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 [i](#)

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	436/554 (78%)	0.10	24 (5%) <span style="border: 1px solid red; padding: 2px;">25</span> <span style="border: 1px solid red; padding: 2px;">31</span>	14, 28, 74, 91	0

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	316	GLY	6.5
1	A	233	GLU	5.5
1	A	532	LEU	5.5
1	A	250	VAL	4.6
1	A	251	ILE	4.4
1	A	438	LEU	3.8
1	A	247	VAL	3.1
1	A	320	LYS	3.1
1	A	252	LYS	3.0
1	A	186	VAL	2.8
1	A	318	SER	2.6
1	A	212	GLU	2.5
1	A	188	THR	2.5
1	A	235	LYS	2.4
1	A	73	GLU	2.4
1	A	70	GLU	2.3
1	A	317	LEU	2.3
1	A	254	VAL	2.3
1	A	69	THR	2.3
1	A	321	ASP	2.3
1	A	319	ASN	2.2
1	A	213	GLY	2.1
1	A	313	ASP	2.0
1	A	324	ARG	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](#)

There are no ligands in this entry.

## 6.5 Other polymers [i](#)

There are no such residues in this entry.