

Full wwPDB X-ray Structure Validation Report (i)

May 21, 2020 – 11:00 pm BST

PDB ID	:	6I1E
Title	:	Crystal structure of Pseudomonas aeruginosa Penicillin-Binding Protein 3 in
		complex with amoxicillin
Authors	:	Bellini, D.; Dowson, C.G.
Deposited on		
Resolution	:	1.64 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 following versions of software and data (see references (1)) were used in the production of this report:

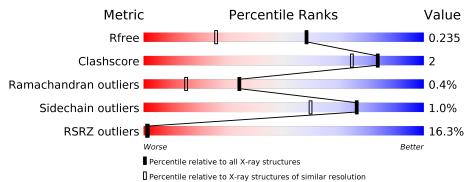
MolProbity		4.02b-467 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix)		1.13
EDS	:	2.11
buster-report	:	1.1.7 (2018)
Percentile statistics	:	20191225.v01 (using entries in the PDB archive December 25th 2019)
Refmac	:	5.8.0158
$\operatorname{CCP4}$:	$7.0.044 (\mathrm{Gargrove})$
Ideal geometry (proteins)	:	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 1.64 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	$egin{array}{c} {f Whole archive}\ (\#{f Entries}) \end{array}$	${f Similar\ resolution}\ (\#{ m Entries},{ m resolution\ range}({ m \AA}))$
R_{free}	130704	3122(1.66-1.62)
Clashscore	141614	3268 (1.66-1.62)
Ramachandran outliers	138981	3215 (1.66-1.62)
Sidechain outliers	138945	3215 (1.66-1.62)
RSRZ outliers	127900	3079 (1.66-1.62)

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		
			16%		
1	А	537	90%	6%	•



2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 4120 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.

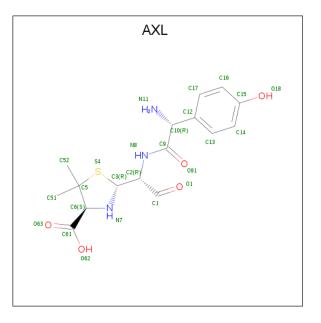
• Molecule 1 is a protein called Peptidoglycan D,D-transpeptidase FtsI.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	515	Total	С	Ν	Ο	\mathbf{S}	0	0	0
	л	515	3927	2478	713	724	12	0	0	0

There are 7 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	43	GLY	-	expression tag	UNP Q51504
A	44	PRO	-	expression tag	UNP Q51504
A	45	GLY	-	expression tag	UNP Q51504
A	46	TYR	-	expression tag	UNP Q51504
A	47	GLN	-	expression tag	UNP Q51504
А	48	ASP	-	expression tag	UNP Q51504
А	49	PRO	-	expression tag	UNP Q51504

• Molecule 2 is 2-{1-[2-AMINO-2-(4-HYDROXY-PHENYL)-ACETYLAMINO]-2-OXO-ETH YL}-5,5-DIMETHYL-THIAZOLIDINE-4-CARBOXYLIC ACID (three-letter code: AXL) (formula: C₁₆H₂₁N₃O₅S) (labeled as "Ligand of Interest" by author).





Mol	Chain	Residues	Atoms				ZeroOcc	AltConf	
2	А	1	Total	С	Ν	Ο	S	0	0
_		_	25	16	- 3	5	1	, i i i i i i i i i i i i i i i i i i i	0

• Molecule 3 is water.

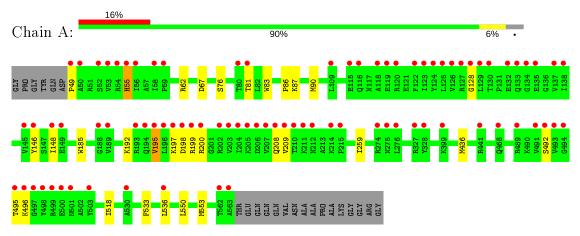
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	А	168	Total O 168 168	0	0



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: Peptidoglycan D,D-transpeptidase FtsI





4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	68.44Å 82.78 Å 89.44 Å	Depositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	60.75 - 1.64	Depositor
Resolution (A)	52.75 - 1.64	EDS
% Data completeness	$100.0 \ (60.75 - 1.64)$	Depositor
(in resolution range)	$100.0\ (52.75 - 1.64)$	EDS
R _{merge}	(Not available)	Depositor
R _{sym}	(Not available)	Depositor
$< I/\sigma(I) > 1$	$1.58 (at 1.64 \text{\AA})$	Xtriage
Refinement program	REFMAC 5.6.0117	Depositor
D D	0.219 , 0.236	Depositor
R, R_{free}	0.215 , 0.235	DCC
R_{free} test set	3096 reflections $(4.92%)$	wwPDB-VP
Wilson B-factor $(Å^2)$	26.9	Xtriage
Anisotropy	0.033	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.35, 37.6	EDS
L-test for twinning ²	$ L > = 0.49, < L^2 > = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.96	EDS
Total number of atoms	4120	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 5.31% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



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: AXL

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	Bo	nd lengths	Bond angles		
	Unain	RMSZ	# Z > 5	RMSZ	# Z > 5	
1	А	0.44	1/4005~(0.0%)	0.55	0/5434	

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
1	А	185	TRP	CD2-CE2	5.17	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	А	3927	0	3989	19	0
2	А	25	0	19	0	0
3	А	168	0	0	0	2
All	All	4120	0	4008	19	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 2.

All (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:495:THR:OG1	1:A:496:LYS:HA	1.72	0.88
1:A:81:THR:HG22	1:A:128:GLY:H	1.44	0.81
1:A:495:THR:N	1:A:496:LYS:HB2	2.02	0.75
1:A:495:THR:H	1:A:496:LYS:HE3	1.59	0.67
1:A:86:PRO:O	1:A:90:MET:HG2	2.00	0.62
1:A:49:PRO:N	1:A:199:ARG:HE	2.03	0.57
1:A:195:VAL:HG11	1:A:197:LYS:HE2	1.89	0.55
1:A:81:THR:HG22	1:A:128:GLY:N	2.20	0.50
1:A:259:ILE:HD12	1:A:436:MET:HG2	1.95	0.49
1:A:62:ARG:HB3	1:A:76:SER:OG	2.14	0.48
1:A:533:PHE:HB2	1:A:536:LEU:HD12	1.97	0.47
1:A:492:SER:HB3	1:A:496:LYS:HB3	1.98	0.45
1:A:55:HIS:HB3	1:A:192:LYS:HD2	1.98	0.45
1:A:83:TRP:CE2	1:A:146:TYR:HB2	2.51	0.45
1:A:81:THR:OG1	1:A:148:ILE:HG13	2.19	0.43
1:A:495:THR:H	1:A:496:LYS:HB2	1.82	0.42
1:A:87:LYS:HD2	1:A:200:ARG:HA	2.02	0.41
1:A:518:ILE:HD11	1:A:553:MET:SD	2.60	0.41
1:A:518:ILE:HD12	1:A:550:LEU:HD23	2.04	0.40

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	Interatomic distance (Å)	Clash overlap (Å)
3:A:811:HOH:O	3:A:816:HOH:O[4_455]	1.77	0.43
3:A:859:HOH:O	3:A:955:HOH:O[3_544]	1.92	0.28

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	А	513/537~(96%)	498 (97%)	13~(2%)	2~(0%)	34 15



All (2) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	А	208	GLN
1	А	209	VAL

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	А	410/424~(97%)	406~(99%)	4 (1%)	76 59	

All (4) residues with a non-rotameric sidechain are listed below:

Mol	Chain	\mathbf{Res}	Type
1	А	55	HIS
1	А	67	ASP
1	А	195	VAL
1	А	198	ASP

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (1) such sidechains are listed below:

Mol	Chain	\mathbf{Res}	Type
1	А	336	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)

1 ligand is modelled in this entry.

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).

Mal	Tuno	Chain	Dog	Link	Bo	ond leng	ths	B	ond ang	les
	Mol Type Chain	nes		Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2	
2	AXL	А	701	1	17,26,26	1.73	4 (23%)	22,38,38	1.48	<mark>3 (13%)</mark>

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.

Mol	Type	Chain	Res	Link	Chirals	Torsions	Rings
2	AXL	А	701	1	-	0/12/37/37	0/2/2/2

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\operatorname{\AA})$	Ideal(Å)
2	А	701	AXL	C5-S4	-4.47	1.76	1.85
2	А	701	AXL	C12-C10	-3.54	1.48	1.52
2	А	701	AXL	C3-S4	-2.87	1.77	1.84
2	А	701	AXL	C6-N7	2.03	1.49	1.46

All (3) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\mathbf{Observed}(^{o})$	$Ideal(^{o})$
2	А	701	AXL	C2-N8-C9	-5.03	117.62	123.12
2	А	701	AXL	C5-S4-C3	2.65	99.66	93.99
2	А	701	AXL	C3-C2-N8	2.21	114.78	109.98

There are no chirality outliers.

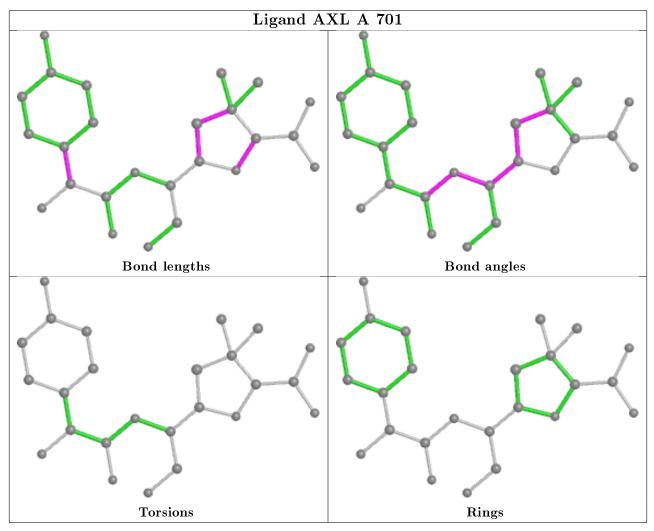


There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

The following is a two-dimensional graphical depiction of Mogul quality analysis of bond lengths, bond angles, torsion angles, and ring geometry for all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the validation Tables will also be included. For torsion angles, if less then 5% of the Mogul distribution of torsion angles is within 10 degrees of the torsion angle in question, then that torsion angle is considered an outlier. Any bond that is central to one or more torsion angles identified as an outlier by Mogul will be highlighted in the graph. For rings, the root-mean-square deviation (RMSD) between the ring in question and similar rings identified by Mogul is calculated over all ring torsion angles. If the average RMSD is greater than 60 degrees and the minimal RMSD between the ring in question and any Mogul-identified rings is also greater than 60 degrees, then that ring is considered an outlier. The outliers are highlighted in purple. The color gray indicates Mogul did not find sufficient equivalents in the CSD to analyse the geometry.





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, 95th 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	$\mathbf{OWAB}(\mathbf{A}^2)$	Q<0.9
1	А	515/537~(95%)	0.93	84 (16%) 1 1	18, 27, 79, 114	0

All (84) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	А	195	VAL	16.7
1	А	204	ILE	16.1
1	А	493	VAL	14.6
1	А	207	VAL	14.2
1	А	498	TYR	12.6
1	А	276	LEU	11.5
1	А	495	THR	10.3
1	А	49	PRO	9.9
1	А	206	ASP	9.9
1	А	203	VAL	9.5
1	А	497	GLY	9.5
1	А	196	LEU	9.3
1	А	496	LYS	9.3
1	А	53	VAL	8.5
1	А	208	GLN	7.6
1	А	55	HIS	7.4
1	А	563	ALA	7.0
1	А	200	ARG	7.0
1	А	494	GLY	6.9
1	А	209	VAL	6.9
1	А	54	ARG	6.7
1	А	500	GLU	6.4
1	А	562	THR	6.2
1	А	492	SER	6.1
1	А	194	GLN	5.7
1	А	275	ASN	5.6
1	A	501	ASN	5.5

Continued on next page...



Mol	Chain	Res	Type	RSRZ
1	А	491	VAL	5.5
1	А	199	ARG	5.4
1	А	274	ARG	5.0
1	А	119	GLU	4.5
1	А	123	ILE	4.4
1	А	210	THR	4.2
1	А	124	TYR	4.2
1	А	135	GLU	4.2
1	А	122	PHE	4.2
1	А	202	ARG	4.1
1	А	212	ASN	3.8
1	А	211	LYS	3.8
1	А	148	ILE	3.8
1	А	189	VAL	3.8
1	А	134	GLY	3.7
1	А	50	ALA	3.7
1	А	205	LYS	3.6
1	А	59	PRO	3.6
1	А	188	GLY	3.6
1	А	58	ILE	3.4
1	А	118	ALA	3.4
1	А	116	GLN	3.4
1	А	193	ARG	3.3
1	А	198	ASP	3.2
1	А	115	GLU	3.2
1	А	530	ALA	3.2
1	А	52	SER	3.2
1	А	56	ILE	3.0
1	А	109	LEU	3.0
1	А	489	ARG	2.9
1	А	214	LYS	2.8
1	А	215	PRO	2.8
1	А	80	THR	2.8
1	А	81	THR	2.8
1	А	129	LEU	2.8
1	А	137	VAL	2.7
1	А	126	VAL	2.7
1	А	327	ARG	2.6
1	А	499	ARG	2.6
1	А	149	GLU	2.6
1	А	130	THR	2.5
		1.05	TTTT	2 4

Continued from previous page...

LEU Continued on next page...

2.5

125

А

1



Mol	Chain	Res	Type	RSRZ
1	А	146	TYR	2.4
1	А	441	ARG	2.4
1	А	120	ARG	2.3
1	А	145	VAL	2.3
1	А	536	LEU	2.3
1	А	138	ILE	2.3
1	А	197	LYS	2.3
1	А	468	GLN	2.2
1	А	127	ARG	2.2
1	А	133	GLN	2.2
1	А	328	TYR	2.1
1	А	399	LYS	2.1
1	А	503	TYR	2.1
1	А	128	GLY	2.1
1	А	132	GLU	2.0

Continued from previous page...

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.

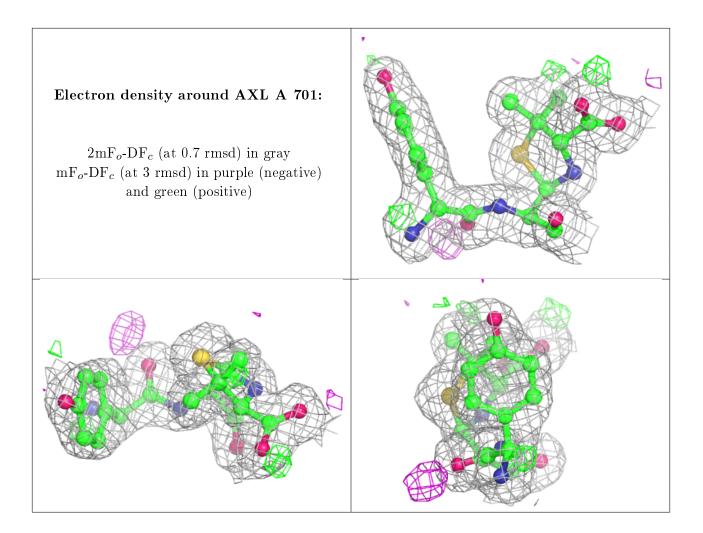
6.4 Ligands (i)

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	$\mathbf{B} ext{-factors}(\mathbf{A}^2)$	Q<0.9
2	AXL	А	701	25/25	0.94	0.09	$20,\!23,\!28,\!30$	0

The following is a graphical depiction of the model fit to experimental electron density of all instances of the Ligand of Interest. In addition, ligands with molecular weight > 250 and outliers as shown on the geometry validation Tables will also be included. Each fit is shown from different orientation to approximate a three-dimensional view.





6.5 Other polymers (i)

There are no such residues in this entry.

