



wwPDB X-ray Structure Validation Summary Report

Feb 27, 2014 – 01:51 PM GMT

PDB ID : 2B5R
Title : 1B Lactamase / B Lactamase Inhibitor
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Deposited on : 2005-09-29
Resolution : 1.65 Å(reported)

This is a wwPDB validation summary report for a publicly released PDB entry.

We welcome your comments at validation@mail.wwpdb.org

A user guide is available at <http://wwpdb.org/ValidationPDFNotes.html>

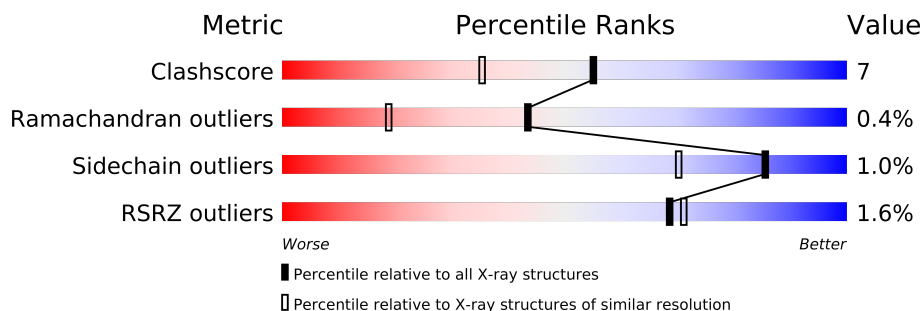
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity	:	4.02b-467
Mogul	:	1.15 2013
Xtriage (Phenix)	:	dev-1323
EDS	:	stable22639
Percentile statistics	:	21963
Refmac	:	5.8.0049
CCP4	:	6.3.0 (Settle)
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et. al. (1996)
Validation Pipeline (wwPDB-VP)	:	stable22683

1 Overall quality at a glance

The reported resolution of this entry is 1.65 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))
Clashscore	79885	1001 (1.66-1.66)
Ramachandran outliers	78287	1581 (1.68-1.64)
Sidechain outliers	78261	1580 (1.68-1.64)
RSRZ outliers	66119	1404 (1.68-1.64)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	263	
1	B	263	
2	C	165	
2	D	165	

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 7047 atoms, of which 0 are hydrogen and 0 are deuterium.

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 Beta-lactamase TEM.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	263	Total	C	N	O	S	0	0	0
			2026	1265	361	389	11			
1	B	263	Total	C	N	O	S	0	0	0
			2026	1265	361	389	11			

There are 6 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	84	ILE	VAL	ENGINEERED	UNP P62593
A	104	TYR	GLU	ENGINEERED	UNP P62593
A	105	ASN	TYR	ENGINEERED	UNP P62593
B	84	ILE	VAL	ENGINEERED	UNP P62593
B	104	TYR	GLU	ENGINEERED	UNP P62593
B	105	ASN	TYR	ENGINEERED	UNP P62593

- Molecule 2 is a protein called Beta-lactamase inhibitory protein.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
2	C	165	Total	C	N	O	S	0	0	0
			1235	778	207	243	7			
2	D	165	Total	C	N	O	S	0	0	0
			1235	778	207	243	7			

- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	116	Total	O	0	0
			116	116		
3	B	136	Total	O	0	0
			136	136		
3	C	119	Total	O	0	0
			119	119		

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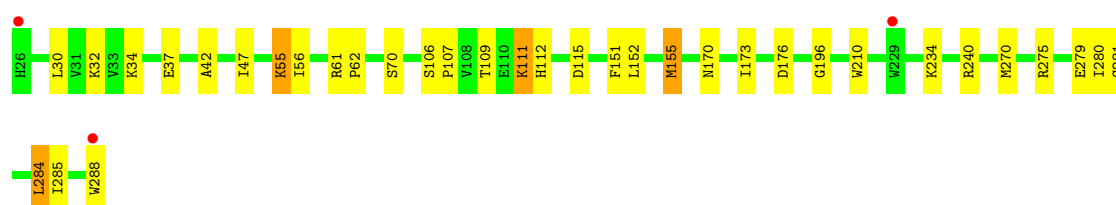
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	D	154	Total 154	O 154	0	0

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 errors 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: Beta-lactamase TEM

Chain A:



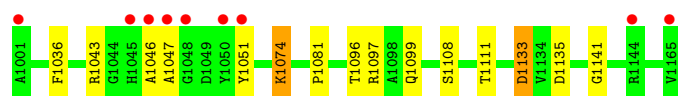
- Molecule 1: Beta-lactamase TEM

Chain B:



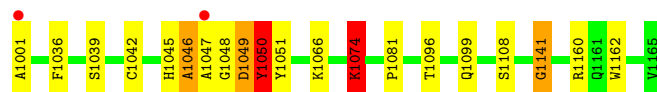
- Molecule 2: Beta-lactamase inhibitory protein

Chain C:



- Molecule 2: Beta-lactamase inhibitory protein

Chain D:



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	45.80Å 124.47Å 157.26Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	48.30 – 1.65 48.31 – 1.65	Depositor EDS
% Data completeness (in resolution range)	99.7 (48.30-1.65) 98.0 (48.31-1.65)	Depositor EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.09 (at 1.65Å)	Xtriage
Refinement program	CNS 0.9	Depositor
R, R_{free}	0.190 , 0.222 0.196 , (Not available)	Depositor DCC
R_{free} test set	No test flags present.	DCC
Wilson B-factor (Å ²)	15.4	Xtriage
Anisotropy	0.229	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.37 , 36.7	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle L \rangle = 0.48$, $\langle L^2 \rangle = 0.32$	Xtriage
Outliers	0 of 109217 reflections	Xtriage
F_o, F_c correlation	0.95	EDS
Total number of atoms	7047	wwPDB-VP
Average B, all atoms (Å ²)	18.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.01% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

5 Model quality

5.1 Standard geometry

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	0.79	1/2060 (0.0%)	0.94	2/2790 (0.1%)
1	B	0.86	0/2060	0.98	7/2790 (0.3%)
2	C	0.84	0/1264	0.89	3/1715 (0.2%)
2	D	0.89	1/1264 (0.1%)	1.01	3/1715 (0.2%)
All	All	0.84	2/6648 (0.0%)	0.96	15/9010 (0.2%)

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
2	D	0	1

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	D	1042	CYS	CB-SG	5.57	1.91	1.82
1	A	210	TRP	CB-CG	5.06	1.59	1.50

The worst 5 of 15 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	288	TRP	CA-C-O	-14.65	89.34	120.10
1	B	288	TRP	N-CA-C	-11.88	78.91	111.00
2	D	1001	ALA	N-CA-C	-9.24	86.04	111.00
1	B	288	TRP	CA-CB-CG	7.26	127.50	113.70
1	B	288	TRP	CB-CA-C	-6.18	98.05	110.40

There are no chirality outliers.

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
2	D	1141	GLY	Peptide

5.2 Close contacts ⓘ

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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	2026	0	2036	35	0
1	B	2026	0	2036	17	0
2	C	1235	0	1170	16	0
2	D	1235	0	1168	23	0
3	A	116	0	0	2	0
3	B	136	0	0	3	0
3	C	119	0	0	6	0
3	D	154	0	0	5	0
All	All	7047	0	6410	86	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 7.

The worst 5 of 86 close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:196:GLY:HA3	3:A:2452:HOH:O	1.36	1.22
1:A:173:ILE:HD13	1:A:240:ARG:NH1	1.58	1.18
1:A:152:LEU:HD23	1:A:155:MET:CE	1.83	1.09
1:A:173:ILE:HD13	1:A:240:ARG:HH12	0.92	1.07
1:A:152:LEU:HD23	1:A:155:MET:HE3	1.31	1.06

There are no symmetry-related clashes.

5.3 Torsion angles

5.3.1 Protein backbone ⓘ

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	261/263 (99%)	257 (98%)	4 (2%)	0	100	100
1	B	261/263 (99%)	257 (98%)	4 (2%)	0	100	100
2	C	163/165 (99%)	156 (96%)	7 (4%)	0	100	100
2	D	163/165 (99%)	157 (96%)	3 (2%)	3 (2%)	13	1
All	All	848/856 (99%)	827 (98%)	18 (2%)	3 (0%)	43	19

All (3) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
2	D	1046	ALA
2	D	1049	ASP
2	D	1050	TYR

5.3.2 Protein sidechains ⓘ

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	217/217 (100%)	213 (98%)	4 (2%)	71	46
1	B	217/217 (100%)	215 (99%)	2 (1%)	87	74
2	C	126/126 (100%)	126 (100%)	0	100	100
2	D	126/126 (100%)	125 (99%)	1 (1%)	89	78
All	All	686/686 (100%)	679 (99%)	7 (1%)	85	70

5 of 7 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	284	LEU
2	D	1074	LYS
1	B	30	LEU
1	A	55	LYS
1	B	284	LEU

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. 5 of 6 such

sidechains are listed below:

Mol	Chain	Res	Type
1	B	132	ASN
2	D	1099	GLN
2	D	1045	HIS
2	C	1099	GLN
2	D	1089	ASN

5.3.3 RNA ⓘ

There are no RNA chains in this entry.

5.4 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

5.5 Carbohydrates ⓘ

There are no carbohydrates in this entry.

5.6 Ligand geometry ⓘ

There are no ligands in this entry.

5.7 Other polymers ⓘ

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 ⓘ

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, 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	OWAB(Å ²)	Q<0.9
1	A	263/263 (100%)	-0.22	3 (1%) 77 81	8, 17, 32, 45	0
1	B	263/263 (100%)	-0.43	0 100 100	9, 14, 29, 36	0
2	C	165/165 (100%)	-0.06	9 (5%) 24 22	11, 18, 32, 50	0
2	D	165/165 (100%)	-0.34	2 (1%) 75 79	9, 14, 26, 41	0
All	All	856/856 (100%)	-0.27	14 (1%) 68 71	8, 16, 31, 50	0

The worst 5 of 14 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	C	1046	ALA	5.9
2	C	1047	ALA	5.9
2	C	1001	ALA	4.6
2	C	1050	TYR	3.7
2	C	1165	VAL	3.7

6.2 Non-standard residues in protein, DNA, RNA chains ⓘ

There are no non-standard protein/DNA/RNA residues in this entry.

6.3 Carbohydrates ⓘ

There are no carbohydrates in this entry.

6.4 Ligands ⓘ

There are no ligands in this entry.

6.5 Other polymers ⓘ

There are no such residues in this entry.