

wwPDB X-ray Structure Validation Summary Report (i)

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PDB ID 6WY4

> Title Crystal Structure of Wild Type Class D beta-lactamase from Clostridium dif-

> > ficile 630

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Resolution 1.80 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 types of validation reports are described at http://www.wwpdb.org/validation/2017/FAQs#types.

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) Mogul

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)

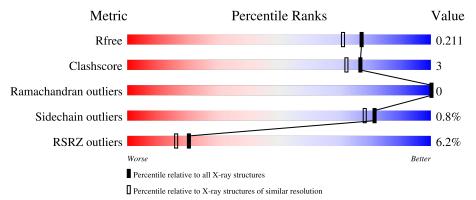
Validation Pipeline (wwPDB-VP) : 2.36

1 Overall quality at a glance (i)

The following experimental techniques were used to determine the structure: X- $RAY\ DIFFRACTION$

The reported resolution of this entry is 1.80 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 $(\# \mathrm{Entries})$	$\begin{array}{c} {\rm Similar\ resolution} \\ (\#{\rm Entries,\ resolution\ range(\mathring{\rm A})}) \end{array}$
R_{free}	130704	5950 (1.80-1.80)
Clashscore	141614	6793 (1.80-1.80)
Ramachandran outliers	138981	6697 (1.80-1.80)
Sidechain outliers	138945	6696 (1.80-1.80)
RSRZ outliers	127900	5850 (1.80-1.80)

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 >=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	255	92%	5% •
1	В	255	7% 88%	9% •
1	С	255	7% 92%	7%
1	D	255	7% 90%	9%



2 Entry composition (i)

There are 4 unique types of molecules in this entry. The entry contains 9261 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 Beta-lactamase.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	A	248	Total	С	N	О	S	0	6	0
1	A	240	2062	1312	334	407	9	0	0	
1	В	248	Total	С	N	О	S	0	7	0
1	Ъ	240	2067	1312	338	408	9	0	'	
1	С	254	Total	С	N	О	S	0	2	0
1		204	2070	1313	339	410	8	0	<u> </u>	
1	D	254	Total	С	N	О	S	0	1	0
1	ע	204	2086	1323	343	412	8		4	

There are 16 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	291	GLY	-	expression tag	UNP Q188Q3
A	292	ALA	-	expression tag	UNP Q188Q3
A	293	ALA	-	expression tag	UNP Q188Q3
A	294	LEU	-	expression tag	UNP Q188Q3
В	291	GLY	-	expression tag	UNP Q188Q3
В	292	ALA	-	expression tag	UNP Q188Q3
В	293	ALA	-	expression tag	UNP Q188Q3
В	294	LEU	-	expression tag	UNP Q188Q3
С	291	GLY	-	expression tag	UNP Q188Q3
С	292	ALA	-	expression tag	UNP Q188Q3
С	293	ALA	-	expression tag	UNP Q188Q3
С	294	LEU	-	expression tag	UNP Q188Q3
D	291	GLY	-	expression tag	UNP Q188Q3
D	292	ALA	-	expression tag	UNP Q188Q3
D	293	ALA	-	expression tag	UNP Q188Q3
D	294	LEU	-	expression tag	UNP Q188Q3

• Molecule 2 is DI(HYDROXYETHYL)ETHER (three-letter code: PEG) (formula: C₄H₁₀O₃).





Mol	Chain	Residues	Atoms		ZeroOcc	AltConf	
2	В	1	Total 7	C 4	O 3	0	0

 \bullet Molecule 3 is SODIUM ION (three-letter code: NA) (formula: Na).

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	D	1	Total Na 1 1	0	0

• Molecule 4 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
4	A	286	Total O 309 309	0	25
4	В	250	Total O 265 265	0	16
4	С	184	Total O 201 201	0	21
4	D	181	Total O 193 193	0	14



3 Residue-property plots (i)

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: Beta-lactamase Chain A: 92% • Molecule 1: Beta-lactamase Chain B: 88% • Molecule 1: Beta-lactamase Chain C: 92% • Molecule 1: Beta-lactamase Chain D: 90%

4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants	85.33Å 93.26Å 138.05Å	Donositor
a, b, c, α , β , γ	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	29.66 - 1.80	Depositor
Resolution (A)	29.66 - 1.80	EDS
% Data completeness	100.0 (29.66-1.80)	Depositor
(in resolution range)	100.0 (29.66-1.80)	EDS
R_{merge}	0.07	Depositor
R_{sym}	0.07	Depositor
$< I/\sigma(I) > 1$	2.52 (at 1.80Å)	Xtriage
Refinement program	REFMAC 5.8.0258	Depositor
D D.	0.176 , 0.209	Depositor
R, R_{free}	0.182 , 0.211	DCC
R_{free} test set	5203 reflections (5.08%)	wwPDB-VP
Wilson B-factor (Å ²)	23.7	Xtriage
Anisotropy	0.112	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$, $B_{sol}(Å^2)$	0.30 , 48.9	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	9261	wwPDB-VP
Average B, all atoms (Å ²)	33.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 3.94% of the height of the origin peak. No significant pseudotranslation is detected.

²Theoretical 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.



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: KCX, NA, PEG

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	Mol Chain		lengths	Bond	angles
IVIOI			RMSZ # Z > 5		# Z > 5
1	A	0.65	0/2089	0.72	0/2808
1	В	0.66	0/2093	0.72	0/2812
1	С	0.66	0/2096	0.70	0/2816
1	D	0.66	0/2112	0.70	0/2838
All	All	0.66	0/8390	0.71	0/11274

There are no bond length outliers.

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	2062	0	2012	7	0
1	В	2067	0	2021	20	0
1	С	2070	0	2023	10	0
1	D	2086	0	2040	13	0
2	В	7	0	10	0	0
3	D	1	0	0	0	0
4	A	309	0	0	1	0
4	В	265	0	0	2	0
4	С	201	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
4	D	193	0	0	0	0
All	All	9261	0	8106	50	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 50 close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

Atom-1	Atom-2	$\begin{array}{c} {\rm Interatomic} \\ {\rm distance} \ ({\rm \AA}) \end{array}$	$\begin{array}{c} \text{Clash} \\ \text{overlap } (\text{\AA}) \end{array}$
1:B:73[B]:GLU:OE1	4:B:401:HOH:O	2.05	0.74
1:B:193[B]:MET:SD	1:B:246:VAL:HG12	2.32	0.69
1:D:84:ILE:HG23	1:D:211:LEU:HD21	1.83	0.61
1:D:99:LYS:NZ	1:D:100:GLU:OE2	2.35	0.60
1:D:217:ILE:HB	1:D:226:VAL:HG13	1.82	0.59

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	ntiles
1	A	251/255~(98%)	243 (97%)	8 (3%)	0	100	100
1	В	$252/255\ (99\%)$	244 (97%)	8 (3%)	0	100	100
1	\mathbf{C}	253/255~(99%)	248 (98%)	5 (2%)	0	100	100
1	D	255/255 (100%)	246 (96%)	9 (4%)	0	100	100
All	All	1011/1020 (99%)	981 (97%)	30 (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	$227/225 \ (101\%)$	225 (99%)	2 (1%)		78	75
1	В	228/225 (101%)	227 (100%)	1 (0%)		91	89
1	С	226/225 (100%)	224 (99%)	2 (1%)		78	75
1	D	228/225 (101%)	226 (99%)	2 (1%)		78	75
All	All	909/900 (101%)	902 (99%)	7 (1%)		81	78

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

Mol	Chain	Res	Type
1	С	248	MET
1	С	251	ARG
1	D	248	MET
1	D	113	LYS
1	В	248	MET

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. There are no such sidechains identified.

5.3.3 RNA (i)

There are no RNA molecules in this entry.

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

4 non-standard protein/DNA/RNA residues are 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).



Mol Ty	Tuno	Chain	Res	Link	Bond lengths				Bond angles		
IVIOI	Type	Chain	nes	LIIIK	Counts	RMSZ	# Z >2	Counts	RMSZ	# Z > 2	
1	KCX	A	83	1	9,11,12	0.48	0	5,12,14	0.43	0	
1	KCX	В	83	1	9,11,12	1.85	1 (11%)	5,12,14	3.36	1 (20%)	
1	KCX	D	83	1	9,11,12	0.40	0	5,12,14	0.50	0	
1	KCX	С	83	1	9,11,12	1.88	1 (11%)	5,12,14	3.24	1 (20%)	

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
1	KCX	A	83	1	-	0/9/10/12	-
1	KCX	В	83	1	-	0/9/10/12	-
1	KCX	D	83	1	-	0/9/10/12	-
1	KCX	С	83	1	-	0/9/10/12	-

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	$\operatorname{Observed}(\text{\AA})$	Ideal(Å)
1	С	83	KCX	OQ1-CX	5.49	1.31	1.21
1	В	83	KCX	OQ1-CX	5.47	1.31	1.21

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	\mathbf{Z}	$\mathbf{Observed}(^o)$	$\operatorname{Ideal}({}^o)$
1	В	83	KCX	OQ1-CX-NZ	-7.52	113.31	124.96
1	С	83	KCX	OQ1-CX-NZ	-7.24	113.74	124.96

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

2 monomers are involved in 3 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
1	A	83	KCX	2	0
1	В	83	KCX	1	0



5.5 Carbohydrates (i)

There are no monosaccharides in this entry.

5.6 Ligand geometry (i)

Of 2 ligands modelled in this entry, 1 is monoatomic - leaving 1 for Mogul analysis.

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).

Mol Type	Chain	Pos	Res Link	В	Bond lengths			Bond angles		
	туре	Chain	nes	LIIIK	Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
2	PEG	В	301	-	6,6,6	0.13	0	5,5,5	0.07	0

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	PEG	В	301	-	-	1/4/4/4	-

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

All (1) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
2	В	301	PEG	O2-C3-C4-O4

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	<rsrz></rsrz>	$\#\mathrm{RSRZ}{>}2$	$OWAB(A^2)$	Q < 0.9
1	A	247/255 (96%)	-0.07	7 (2%) 53 47	13, 22, 41, 70	0
1	В	247/255 (96%)	0.18	18 (7%) 15 11	14, 25, 62, 86	0
1	С	253/255~(99%)	0.33	18 (7%) 16 12	15, 34, 67, 75	0
1	D	253/255 (99%)	0.38	19 (7%) 14 11	22, 38, 65, 79	0
All	All	1000/1020 (98%)	0.21	62 (6%) 20 16	13, 30, 65, 86	0

The worst 5 of 62 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	В	252	GLU	5.3
1	A	222	LYS	4.9
1	D	289	ARG	4.4
1	В	220	ASN	4.2
1	D	166	LEU	4.1

6.2 Non-standard residues in protein, DNA, RNA chains (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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
1	KCX	D	83	12/13	0.90	0.18	24,28,39,41	0
1	KCX	В	83	12/13	0.92	0.17	13,15,19,19	0
1	KCX	С	83	12/13	0.95	0.14	18,21,27,29	0
1	KCX	A	83	12/13	0.97	0.15	14,16,17,19	0



6.3 Carbohydrates (i)

There are no monosaccharides 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	$\mathbf{B} ext{-}\mathbf{factors}(\mathbf{\mathring{A}}^2)$	Q<0.9
2	PEG	В	301	7/7	0.81	0.19	58,59,60,64	0
3	NA	D	301	1/1	0.95	0.19	39,39,39,39	0

6.5 Other polymers (i)

There are no such residues in this entry.

