

# Full wwPDB X-ray Structure Validation Report (i)

May 26, 2020 – 02:56 pm BST

PDB ID : 5J4A

Title: CdiA-CT toxin from Burkholderia pseudomallei E479 in complex with cognate

CdiI immunity protein

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Deposited on : 2016-03-31

Resolution : 2.00 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:

 $\begin{array}{ccc} Mol Probity & : & 4.02b\text{-}467 \\ Xtriage & (Phenix) & : & 1.13 \end{array}$ 

EDS : 2.11

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)

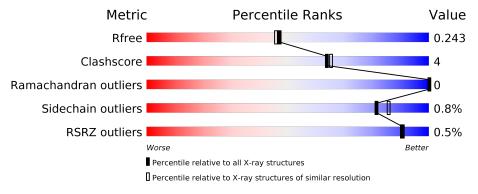
Validation Pipeline (wwPDB-VP) : 2.11

## 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 2.00 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	$\begin{array}{c} \text{Whole archive} \\ (\#\text{Entries}) \end{array}$	$\begin{array}{c} {\rm Similar \; resolution} \\ (\#{\rm Entries, \; resolution \; range(\AA)}) \end{array}$
$R_{free}$	130704	8085 (2.00-2.00)
Clashscore	141614	9178 (2.00-2.00)
Ramachandran outliers	138981	9054 (2.00-2.00)
Sidechain outliers	138945	9053 (2.00-2.00)
RSRZ outliers	127900	7900 (2.00-2.00)

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				
1	A	161	67% 5%	28%			
1	С	161	64% 7% •	28%			
2	В	120	82%	5%	13%		
2	D	120	73%	14%	13%		



## 2 Entry composition (i)

There are 3 unique types of molecules in this entry. The entry contains 3537 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 tRNA nuclease CdiA.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
1	Λ	116	Total	С	N	О	S	0	0	0
1	A	110	873	554	151	166	2	0	U	0
1	С	116	Total	С	N	О	S	0	0	0
1		110	873	554	151	166	2		U	U

There are 4 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	156	MET	-	initiating methionine	UNP H9T8G6
A	285	ALA	ASP	engineered mutation	UNP H9T8G6
С	156	MET	=	initiating methionine	UNP H9T8G6
С	285	ALA	ASP	engineered mutation	UNP H9T8G6

• Molecule 2 is a protein called Immunity protein CdiI.

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf	Trace		
2	D	104	Total	С	N	О	S	0	0	0
	Б	104	818	516	137	161	4	0	U	U
9	D	104	Total	С	N	О	S	0	0	0
2	ש	104	818	516	137	161	4	U	U	U

There are 22 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
В	0	LYS	-	expression tag	UNP H9T8G7
В	110	THR	=	expression tag	UNP H9T8G7
В	111	SER	-	expression tag	UNP H9T8G7
В	112	LEU	-	expression tag	UNP H9T8G7
В	113	GLU	-	expression tag	UNP H9T8G7
В	114	HIS	-	expression tag	UNP H9T8G7
В	115	HIS	_	expression tag	UNP H9T8G7
В	116	HIS	-	expression tag	UNP H9T8G7

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Chain	Residue	Modelled	Actual	Comment	Reference
В	117	HIS	-	expression tag	UNP H9T8G7
В	118	HIS	-	expression tag	UNP H9T8G7
В	119	HIS	_	expression tag	UNP H9T8G7
D	0	LYS	_	expression tag	UNP H9T8G7
D	110	THR	-	expression tag	UNP H9T8G7
D	111	SER	_	expression tag	UNP H9T8G7
D	112	LEU	_	expression tag	UNP H9T8G7
D	113	GLU	_	expression tag	UNP H9T8G7
D	114	HIS	_	expression tag	UNP H9T8G7
D	115	HIS	-	expression tag	UNP H9T8G7
D	116	HIS	_	expression tag	UNP H9T8G7
D	117	HIS	-	expression tag	UNP H9T8G7
D	118	HIS	_	expression tag	UNP H9T8G7
D	119	HIS	-	expression tag	UNP H9T8G7

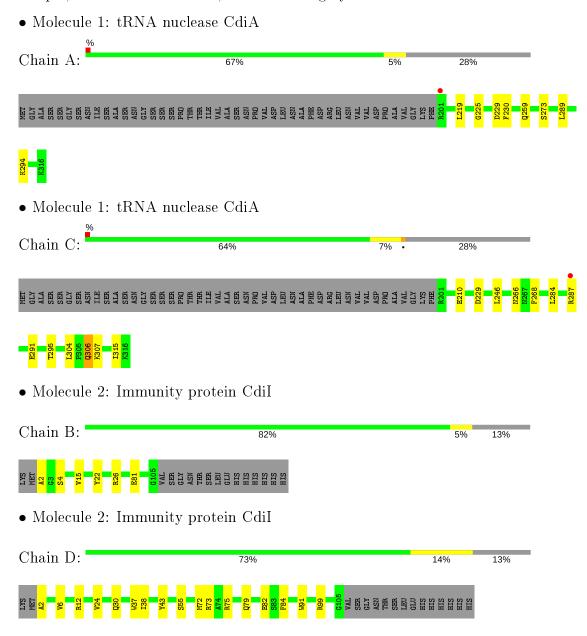
#### • Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	40	Total O 40 40	0	0
3	В	51	Total O 51 51	0	0
3	С	35	Total O 35 35	0	0
3	D	29	Total O 29 29	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.





## 4 Data and refinement statistics (i)

Property	Value	Source
Space group	P 2 21 21	Depositor
Cell constants	54.47Å 73.26Å 110.03Å	Danagitan
a, b, c, $\alpha$ , $\beta$ , $\gamma$	90.00° 90.00° 90.00°	Depositor
Resolution (Å)	43.71 - 2.00	Depositor
Resolution (A)	43.71 - 2.00	EDS
% Data completeness	99.2 (43.71-2.00)	Depositor
(in resolution range)	95.6 (43.71-2.00)	EDS
$R_{merge}$	(Not available)	Depositor
$R_{sym}$	(Not available)	Depositor
$< I/\sigma(I) > 1$	5.56 (at 2.00Å)	Xtriage
Refinement program	PHENIX phenix.refine (1.8.2_1309)	Depositor
D D.	0.193 , 0.237	Depositor
$R, R_{free}$	0.201 , $0.243$	DCC
$R_{free}$ test set	2000 reflections $(6.64\%)$	wwPDB-VP
Wilson B-factor (Å <sup>2</sup> )	34.8	Xtriage
Anisotropy	0.762	Xtriage
Bulk solvent $k_{sol}(e/Å^3)$ , $B_{sol}(Å^2)$	0.37, 39.5	EDS
L-test for twinning <sup>2</sup>	$< L > = 0.49, < L^2> = 0.33$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3537	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: The largest off-origin peak in the Patterson function is 7.48% of the height of the origin peak. No significant pseudotranslation is detected.

<sup>&</sup>lt;sup>2</sup>Theoretical 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.



<sup>&</sup>lt;sup>1</sup>Intensities estimated from amplitudes.

## 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	Mol Chain		lengths	Bond angles		
WIOI	Chain	RMSZ	# Z  > 5	RMSZ	# Z  > 5	
1	A	0.38	0/888	0.53	0/1202	
1	С	0.35	0/888	0.55	0/1202	
2	В	0.41	0/833	0.54	0/1125	
2	D	0.38	0/833	0.51	0/1125	
All	All	0.38	0/3442	0.53	0/4654	

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	873	0	879	5	0
1	С	873	0	879	11	0
2	В	818	0	799	4	0
2	D	818	0	799	11	0
3	A	40	0	0	0	0
3	В	51	0	0	1	0
3	С	35	0	0	3	0
3	D	29	0	0	0	0
All	All	3537	0	3356	28	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 (28) close contacts within the same asymmetric unit are listed below, sorted by their clash magnitude.

A , 4	A	Interatomic	Clash
Atom-1	Atom-2	${f distance} ({f \AA})$	overlap (Å)
1:A:229:ASP:OD2	2:B:2:ALA:HB2	1.84	0.77
1:C:229:ASP:OD2	2:D:2:ALA:HB2	1.86	0.76
1:C:246:LEU:HD13	1:C:284:LEU:HB3	1.71	0.73
1:C:246:LEU:HD11	1:C:268:PHE:CZ	2.30	0.66
1:A:289:LEU:O	1:A:294:LYS:NZ	2.30	0.65
2:D:82:GLU:OE1	2:D:82:GLU:N	2.32	0.63
1:A:225:GLY:HA2	3:B:202:HOH:O	1.99	0.62
2:B:26:ARG:NH1	2:B:81:GLU:OE1	2.34	0.61
1:C:246:LEU:HD11	1:C:268:PHE:HZ	1.65	0.60
1:A:219:LEU:HD22	1:A:230:PHE:HB3	1.84	0.58
1:C:287:ARG:HG3	3:C:401:HOH:O	2.04	0.58
2:B:22:TYR:O	2:B:26:ARG:HG2	2.06	0.56
2:D:6:VAL:HB	2:D:55:SER:HB2	1.89	0.54
1:C:306:GLN:HG3	1:C:307:LYS:N	2.23	0.53
2:D:6:VAL:HG22	2:D:12:ARG:HG2	1.90	0.52
2:B:4:SER:HA	2:B:15:VAL:HG22	1.92	0.52
2:D:30:GLN:O	2:D:73:ARG:NH2	2.40	0.48
1:A:259:GLN:OE1	1:C:266:ASN:HB2	2.15	0.46
1:C:210:GLU:HG2	1:C:315:ILE:HG23	1.96	0.46
2:D:79:GLN:HG3	2:D:84:PHE:CE2	2.52	0.45
2:D:24:VAL:HG13	2:D:43:TYR:CE2	2.53	0.44
1:C:287:ARG:NH2	3:C:403:HOH:O	2.51	0.43
2:D:37:TRP:CE2	2:D:38:ILE:HG13	2.54	0.42
1:C:291:GLU:O	1:C:295:THR:OG1	2.29	0.42
3:C:417:HOH:O	2:D:12:ARG:HD3	2.20	0.41
1:C:210:GLU:HG2	1:C:315:ILE:CG2	2.50	0.41
2:D:72:MET:HE1	2:D:99:ARG:HH21	1.86	0.40
2:D:75:ARG:HD3	2:D:91:TRP:HB3	2.03	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 r	number	of	residues	for	which	the	backbone	conformation	was
analysed, and the total numb	er of	residues								

Mol	Chain	Analysed	Favoured	Allowed	Outliers	Perce	entiles
1	A	114/161 (71%)	113 (99%)	1 (1%)	0	100	100
1	С	114/161 (71%)	112 (98%)	2 (2%)	0	100	100
2	В	102/120~(85%)	101 (99%)	1 (1%)	0	100	100
2	D	102/120~(85%)	101 (99%)	1 (1%)	0	100	100
All	All	432/562 (77%)	427 (99%)	5 (1%)	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	92/128~(72%)	91 (99%)	1 (1%)	73 78
1	С	92/128 (72%)	90 (98%)	2 (2%)	52 55
2	В	91/106 (86%)	91 (100%)	0	100 100
2	D	91/106 (86%)	91 (100%)	0	100 100
All	All	366/468 (78%)	363 (99%)	3 (1%)	81 86

All (3) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	273	SER
1	С	304	LEU
1	С	306	GLN

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

$\mathbf{Mol}$	Chain	Res	Type
1	С	220	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 (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	$\langle { m RSRZ} \rangle$	$\#\mathrm{RSRZ}{>}2$	$\mathbf{OWAB}(\mathrm{\AA}^2)$	Q < 0.9
1	A	116/161 (72%)	-0.17	1 (0%) 84 83	26, 34, 47, 57	0
1	С	116/161 (72%)	0.14	1 (0%) 84 83	28, 41, 54, 64	0
2	В	104/120 (86%)	-0.09	0 100 100	24, 31, 42, 55	0
2	D	104/120 (86%)	-0.03	0 100 100	26, 36, 47, 58	0
All	All	440/562 (78%)	-0.03	2 (0%) 91 90	24, 36, 50, 64	0

#### All (2) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	201	ARG	2.4
1	С	287	ARG	2.3

### 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.

