



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 25, 2017 – 10:45 am GMT

PDB ID : 5M07  
Title : Crystal structure of Mycobacterium tuberculosis PknI kinase domain, C20A mutant  
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Deposited on : 2016-10-03  
Resolution : 2.50 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

<http://wwpdb.org/validation/2016/XrayValidationReportHelp>

with specific help available everywhere you see the ⓘ symbol.

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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)	:	1.9-1692
EDS	:	recalc29047
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	recalc29047

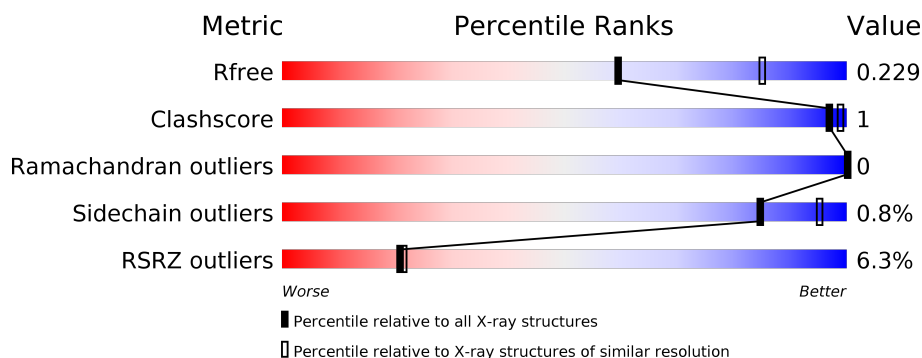
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 2.50 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))
$R_{free}$	100719	3846 (2.50-2.50)
Clashscore	112137	4554 (2.50-2.50)
Ramachandran outliers	110173	4463 (2.50-2.50)
Sidechain outliers	110143	4465 (2.50-2.50)
RSRZ outliers	101464	3876 (2.50-2.50)

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  $\geq 3$ , 2, 1 and 0 types of geometric quality criteria. 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	276	<div> <div>6%</div> <div>88%</div> <div>10%</div> </div>
1	B	276	<div> <div>5%</div> <div>86%</div> <div>11%</div> </div>

The following table lists non-polymeric compounds, carbohydrate monomers and non-standard residues in protein, DNA, RNA chains that are outliers for geometric or electron-density-fit criteria:

Mol	Type	Chain	Res	Chirality	Geometry	Clashes	Electron density
2	NA	A	301	-	-	-	X

## 2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 3870 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 Serine/threonine-protein kinase PknI.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	248	Total	C	N	O	S	0	0	0
			1862	1180	333	342	7			
1	B	247	Total	C	N	O	S	0	0	0
			1852	1174	333	339	6			

There are 42 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	-19	MET	-	initiating methionine	UNP P9WI69
A	-18	GLY	-	expression tag	UNP P9WI69
A	-17	SER	-	expression tag	UNP P9WI69
A	-16	SER	-	expression tag	UNP P9WI69
A	-15	HIS	-	expression tag	UNP P9WI69
A	-14	HIS	-	expression tag	UNP P9WI69
A	-13	HIS	-	expression tag	UNP P9WI69
A	-12	HIS	-	expression tag	UNP P9WI69
A	-11	HIS	-	expression tag	UNP P9WI69
A	-10	HIS	-	expression tag	UNP P9WI69
A	-9	SER	-	expression tag	UNP P9WI69
A	-8	SER	-	expression tag	UNP P9WI69
A	-7	GLY	-	expression tag	UNP P9WI69
A	-6	LEU	-	expression tag	UNP P9WI69
A	-5	VAL	-	expression tag	UNP P9WI69
A	-4	PRO	-	expression tag	UNP P9WI69
A	-3	ARG	-	expression tag	UNP P9WI69
A	-2	GLY	-	expression tag	UNP P9WI69
A	-1	SER	-	expression tag	UNP P9WI69
A	0	HIS	-	expression tag	UNP P9WI69
A	20	ALA	CYS	engineered mutation	UNP P9WI69
B	-19	MET	-	initiating methionine	UNP P9WI69
B	-18	GLY	-	expression tag	UNP P9WI69
B	-17	SER	-	expression tag	UNP P9WI69
B	-16	SER	-	expression tag	UNP P9WI69

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Chain	Residue	Modelled	Actual	Comment	Reference
B	-15	HIS	-	expression tag	UNP P9WI69
B	-14	HIS	-	expression tag	UNP P9WI69
B	-13	HIS	-	expression tag	UNP P9WI69
B	-12	HIS	-	expression tag	UNP P9WI69
B	-11	HIS	-	expression tag	UNP P9WI69
B	-10	HIS	-	expression tag	UNP P9WI69
B	-9	SER	-	expression tag	UNP P9WI69
B	-8	SER	-	expression tag	UNP P9WI69
B	-7	GLY	-	expression tag	UNP P9WI69
B	-6	LEU	-	expression tag	UNP P9WI69
B	-5	VAL	-	expression tag	UNP P9WI69
B	-4	PRO	-	expression tag	UNP P9WI69
B	-3	ARG	-	expression tag	UNP P9WI69
B	-2	GLY	-	expression tag	UNP P9WI69
B	-1	SER	-	expression tag	UNP P9WI69
B	0	HIS	-	expression tag	UNP P9WI69
B	20	ALA	CYS	engineered mutation	UNP P9WI69

- Molecule 2 is SODIUM ION (three-letter code: NA) (formula: Na).

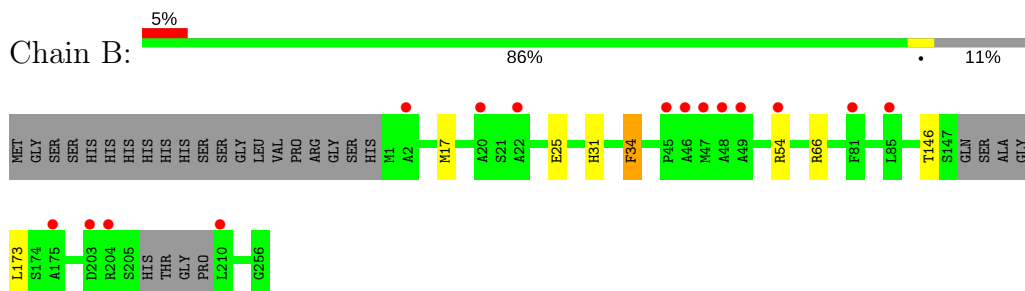
Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
2	B	1	Total Na 1 1	0	0
2	A	1	Total Na 1 1	0	0

- Molecule 3 is water.

Mol	Chain	Residues	Atoms	ZeroOcc	AltConf
3	A	83	Total O 83 83	0	0
3	B	71	Total O 71 71	0	0



- Molecule 1: Serine/threonine-protein kinase PknI



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 43 2 2	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	108.30Å 108.30Å 185.47Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	36.14 – 2.50 46.76 – 2.50	Depositor EDS
% Data completeness (in resolution range)	100.0 (36.14-2.50) 100.0 (46.76-2.50)	Depositor EDS
$R_{merge}$	0.05	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	1.87 (at 2.51Å)	Xtriage
Refinement program	BUSTER 2.10.2	Depositor
R, $R_{free}$	0.208 , 0.235 0.208 , 0.229	Depositor DCC
$R_{free}$ test set	1958 reflections (5.30%)	DCC
Wilson B-factor (Å <sup>2</sup> )	74.4	Xtriage
Anisotropy	0.330	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.34 , 68.2	EDS
L-test for twinning <sup>2</sup>	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	3870	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	85.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.72% 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

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: NA

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	0.49	0/1905	0.64	0/2595
1	B	0.50	0/1895	0.62	0/2582
All	All	0.50	0/3800	0.63	0/5177

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

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	1862	0	1804	4	0
1	B	1852	0	1794	5	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0
3	A	83	0	0	0	0
3	B	71	0	0	0	0
All	All	3870	0	3598	8	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 1.

All (8) 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:B:31:HIS:HB3	1:B:34:PHE:CD2	2.38	0.59
1:A:67:LEU:HD11	1:A:133:LEU:HD23	1.96	0.47
1:B:146:THR:HB	1:B:154:ARG:H	1.81	0.46
1:A:15:VAL:HG12	1:A:16:ARG:HG2	1.99	0.45
1:A:15:VAL:HG21	1:A:37:TRP:HZ3	1.82	0.44
1:B:17:MET:HE3	1:B:25:GLU:HB3	2.00	0.43
1:B:172:GLU:HG2	1:B:173:LEU:N	2.34	0.42
1:A:16:ARG:NH2	1:B:66:ARG:HA	2.36	0.41

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	242/276 (88%)	236 (98%)	6 (2%)	0	100	100
1	B	241/276 (87%)	235 (98%)	6 (2%)	0	100	100
All	All	483/552 (88%)	471 (98%)	12 (2%)	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	182/210 (87%)	181 (100%)	1 (0%)	91	97
1	B	180/210 (86%)	178 (99%)	2 (1%)	78	92
All	All	362/420 (86%)	359 (99%)	3 (1%)	85	95

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

Mol	Chain	Res	Type
1	A	174	SER
1	B	34	PHE
1	B	54	ARG

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

Mol	Chain	Res	Type
1	A	76	HIS

### 5.3.3 RNA ⓘ

There are no RNA molecules 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 ⓘ

Of 2 ligands modelled in this entry, 2 are monoatomic - leaving 0 for Mogul analysis.

There are no bond length outliers.

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

No monomer is involved in short contacts.

## 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, 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	248/276 (89%)	0.32	16 (6%) 20 20	61, 79, 124, 151	0
1	B	247/276 (89%)	0.43	15 (6%) 22 22	61, 81, 136, 151	0
All	All	495/552 (89%)	0.38	31 (6%) 21 21	61, 80, 129, 151	0

All (31) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	47	MET	7.4
1	A	47	MET	7.0
1	B	210	LEU	6.6
1	A	48	ALA	6.2
1	B	49	ALA	5.8
1	B	48	ALA	4.9
1	A	49	ALA	4.8
1	B	46	ALA	4.0
1	A	46	ALA	3.5
1	B	54	ARG	3.5
1	A	175	ALA	3.5
1	A	174	SER	3.3
1	B	45	PRO	3.3
1	A	204	ARG	3.2
1	B	2	ALA	3.2
1	A	57	PHE	2.9
1	B	22	ALA	2.9
1	A	22	ALA	2.8
1	A	3	LEU	2.8
1	B	85	LEU	2.6
1	B	20	ALA	2.5
1	B	175	ALA	2.5
1	B	204	ARG	2.5
1	A	85	LEU	2.3

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Mol	Chain	Res	Type	RSRZ
1	A	43	LEU	2.3
1	A	34	PHE	2.3
1	A	54	ARG	2.2
1	B	81	PHE	2.1
1	A	9	PHE	2.1
1	B	203	ASP	2.0
1	A	173	LEU	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 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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. The B-factors column lists the minimum, median, 95<sup>th</sup> 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	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	NA	A	301	1/1	0.44	0.29	2.63	96,96,96,96	0
2	NA	B	301	1/1	0.96	0.12	-3.42	66,66,66,66	0

## 6.5 Other polymers [i](#)

There are no such residues in this entry.