



Full wwPDB X-ray Structure Validation Report ⓘ

May 15, 2020 – 08:33 pm BST

PDB ID : 4X3F
Title : Crystal structure of the intracellular domain of the M. tuberculosis Ser/Thr kinase PknA
Authors : Wagner, T.; Wehenkel, A.; Alzari, P.M.; Bellinzoni, M.
Deposited on : 2014-11-28
Resolution : 2.90 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 ⓘ symbol.

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

MolProbity : 4.02b-467
Mogul : 1.8.5 (274361), CSD as541be (2020)
Xtriage (Phenix) : 1.13
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

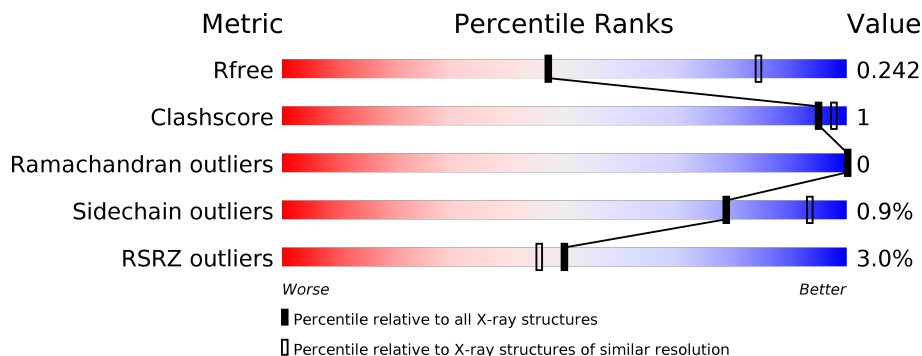
1 Overall quality at a glance i

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.90 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}	130704	1957 (2.90-2.90)
Clashscore	141614	2172 (2.90-2.90)
Ramachandran outliers	138981	2115 (2.90-2.90)
Sidechain outliers	138945	2117 (2.90-2.90)
RSRZ outliers	127900	1906 (2.90-2.90)

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 $\leq 5\%$. 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	337	<div style="display: flex; align-items: center;"> <div style="width: 4%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 79%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 15%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">4% 79% • 18%</p>
2	B	337	<div style="display: flex; align-items: center;"> <div style="width: 3%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 80%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 16%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">3% 80% • 17%</p>
3	C	337	<div style="display: flex; align-items: center;"> <div style="width: 0%; height: 10px; background-color: red; margin-right: 5px;"></div> <div style="width: 78%; height: 10px; background-color: green; margin-right: 5px;"></div> <div style="width: 1%; height: 10px; background-color: yellow; margin-right: 5px;"></div> <div style="width: 21%; height: 10px; background-color: grey;"></div> </div> <p style="text-align: center;">% 78% • 19%</p>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 6027 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 PknA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
1	A	277	1995	1255	361	369	3	7	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	0	GLY	-	expression tag	UNP P9WI83

- Molecule 2 is a protein called Serine/threonine-protein kinase PknA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
2	B	280	2036	1280	371	374	2	9	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
B	0	GLY	-	expression tag	UNP P9WI83

- Molecule 3 is a protein called Serine/threonine-protein kinase PknA.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
			Total	C	N	O	P	S			
3	C	274	1986	1253	362	364	1	6	0	0	0

There is a discrepancy between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
C	0	GLY	-	expression tag	UNP P9WI83

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	3	Total 3	O 3	0	0
4	B	2	Total 2	O 2	0	0
4	C	5	Total 5	O 5	0	0

4 Data and refinement statistics

Property	Value	Source
Space group	C 1 2 1	Depositor
Cell constants a, b, c, α , β , γ	102.31Å 150.18Å 96.54Å 90.00° 96.98° 90.00°	Depositor
Resolution (Å)	43.56 – 2.90 43.56 – 2.90	Depositor EDS
% Data completeness (in resolution range)	99.6 (43.56-2.90) 99.6 (43.56-2.90)	Depositor EDS
R_{merge}	0.11	Depositor
R_{sym}	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ ¹	1.36 (at 2.90Å)	Xtrriage
Refinement program	REFMAC 5.8.0049	Depositor
R, R_{free}	0.219 , 0.243 0.219 , 0.242	Depositor DCC
R_{free} test set	1606 reflections (5.03%)	wwPDB-VP
Wilson B-factor (Å ²)	80.7	Xtrriage
Anisotropy	0.497	Xtrriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 70.2	EDS
L-test for twinning ²	$\langle L \rangle = 0.46$, $\langle L^2 \rangle = 0.29$	Xtrriage
Estimated twinning fraction	No twinning to report.	Xtrriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	6027	wwPDB-VP
Average B, all atoms (Å ²)	95.0	wwPDB-VP

Xtrriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 4.51% of the height of the origin peak. No significant pseudotranslation is detected.*

¹Intensities estimated from amplitudes.

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.

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: TPO, SEP

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	0.44	0/2005	0.55	0/2735
2	B	0.46	0/2057	0.57	0/2804
3	C	0.46	0/2018	0.59	2/2756 (0.1%)
All	All	0.45	0/6080	0.57	2/8295 (0.0%)

There are no bond length outliers.

All (2) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	C	260	ARG	NE-CZ-NH1	6.67	123.64	120.30
3	C	260	ARG	NE-CZ-NH2	-5.08	117.76	120.30

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	1995	0	1902	5	0
2	B	2036	0	1978	5	0
3	C	1986	0	1923	5	0
4	A	3	0	0	0	0
4	B	2	0	0	0	0
4	C	5	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	6027	0	5803	14	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 (14) 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)
2:B:2:SER:N	2:B:3:PRO:HD3	2.20	0.56
3:C:68:LEU:HD23	3:C:73:ILE:HD13	1.97	0.47
2:B:68:LEU:HD23	2:B:73:ILE:HD13	1.97	0.46
1:A:270:ALA:HA	1:A:281:PRO:HD2	1.98	0.46
2:B:8:THR:HG22	2:B:14:ARG:HB2	1.99	0.45
3:C:270:ALA:HA	3:C:281:PRO:HD2	1.98	0.44
3:C:8:THR:HG22	3:C:14:ARG:HB2	1.99	0.43
1:A:140:ARG:NH1	1:A:165:ALA:HB2	2.34	0.43
1:A:34:ARG:CZ	3:C:37:ARG:HB3	2.49	0.43
1:A:68:LEU:HD23	1:A:73:ILE:HD13	2.00	0.42
3:C:288:PRO:HA	3:C:289:PRO:HD3	1.97	0.42
2:B:270:ALA:HA	2:B:281:PRO:HD2	2.02	0.41
1:A:288:PRO:HA	1:A:289:PRO:HD3	1.96	0.41
2:B:288:PRO:HA	2:B:289:PRO:HD3	1.98	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	270/337 (80%)	263 (97%)	7 (3%)	0	100	100
2	B	274/337 (81%)	265 (97%)	9 (3%)	0	100	100

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Mol	Chain	Analysed	Favoured	Allowed	Outliers	Percentiles	
3	C	267/337 (79%)	260 (97%)	7 (3%)	0	100	100
All	All	811/1011 (80%)	788 (97%)	23 (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	183/256 (72%)	182 (100%)	1 (0%)	88	96
2	B	195/257 (76%)	193 (99%)	2 (1%)	76	92
3	C	191/258 (74%)	189 (99%)	2 (1%)	76	92
All	All	569/771 (74%)	564 (99%)	5 (1%)	78	93

All (5) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	121	MET
2	B	95	MET
2	B	121	MET
3	C	81	GLU
3	C	121	MET

Some 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](#)

6 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	Type	Chain	Res	Link	Bond lengths			Bond angles		
					Counts	RMSZ	# Z > 2	Counts	RMSZ	# Z > 2
1	TPO	A	152	1	8,10,11	1.17	1 (12%)	10,14,16	0.94	0
1	SEP	A	46	1	8,9,10	0.83	0	8,12,14	0.95	0
2	TPO	B	21	2	8,10,11	2.33	2 (25%)	10,14,16	1.05	0
1	TPO	A	224	1	8,10,11	0.97	1 (12%)	10,14,16	1.19	0
2	TPO	B	224	2	8,10,11	0.64	0	10,14,16	1.09	0
3	TPO	C	224	3	8,10,11	0.67	0	10,14,16	1.14	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
1	TPO	A	152	1	-	3/9/11/13	-
1	SEP	A	46	1	-	5/5/8/10	-
2	TPO	B	21	2	-	5/9/11/13	-
1	TPO	A	224	1	-	2/9/11/13	-
2	TPO	B	224	2	-	2/9/11/13	-
3	TPO	C	224	3	-	2/9/11/13	-

All (4) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	B	21	TPO	P-OG1	5.91	1.70	1.59
1	A	224	TPO	P-O1P	2.45	1.58	1.50
1	A	152	TPO	P-O1P	2.39	1.58	1.50
2	B	21	TPO	P-O1P	2.34	1.58	1.50

There are no bond angle outliers.

There are no chirality outliers.

All (19) torsion outliers are listed below:

Mol	Chain	Res	Type	Atoms
1	A	152	TPO	O-C-CA-CB
1	A	152	TPO	CG2-CB-OG1-P
1	A	152	TPO	CB-OG1-P-O3P
1	A	46	SEP	N-CA-CB-OG
1	A	46	SEP	CB-OG-P-O1P
1	A	46	SEP	CB-OG-P-O2P
1	A	46	SEP	CB-OG-P-O3P
2	B	21	TPO	N-CA-CB-CG2
2	B	21	TPO	N-CA-CB-OG1
2	B	21	TPO	C-CA-CB-CG2
2	B	21	TPO	CB-OG1-P-O1P
1	A	46	SEP	CA-CB-OG-P
2	B	21	TPO	CB-OG1-P-O2P
1	A	224	TPO	C-CA-CB-CG2
2	B	224	TPO	C-CA-CB-CG2
3	C	224	TPO	C-CA-CB-CG2
1	A	224	TPO	O-C-CA-CB
2	B	224	TPO	O-C-CA-CB
3	C	224	TPO	O-C-CA-CB

There are no ring outliers.

No monomer is involved in short contacts.

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

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	274/337 (81%)	0.22	13 (4%) 31 28	68, 96, 147, 178	0
2	B	278/337 (82%)	0.13	10 (3%) 42 37	63, 88, 134, 180	0
3	C	273/337 (81%)	-0.00	2 (0%) 87 87	65, 89, 125, 147	0
All	All	825/1011 (81%)	0.12	25 (3%) 50 45	63, 91, 136, 180	0

All (25) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
2	B	86	GLY	5.8
1	A	86	GLY	5.5
1	A	49	SER	5.5
1	A	48	PHE	5.1
1	A	54	PHE	4.9
2	B	289	PRO	3.7
1	A	58	PHE	3.2
3	C	288	PRO	3.1
2	B	50	SER	3.0
2	B	176	MET	2.9
1	A	288	PRO	2.7
1	A	87	GLU	2.7
1	A	51	ASP	2.6
3	C	289	PRO	2.5
2	B	88	GLY	2.5
2	B	163	ALA	2.4
1	A	91	ALA	2.3
1	A	166	VAL	2.3
1	A	50	SER	2.3
1	A	275	ARG	2.2
2	B	288	PRO	2.1
2	B	87	GLU	2.1
2	B	85	ASN	2.0

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Mol	Chain	Res	Type	RSRZ
1	A	274	VAL	2.0
2	B	115	LEU	2.0

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, 95th 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	B-factors(Å ²)	Q<0.9
1	SEP	A	46	10/11	0.82	0.29	147,167,181,181	0
2	TPO	B	21	11/12	0.84	0.21	106,119,159,159	0
1	TPO	A	152	11/12	0.87	0.37	110,122,140,151	0
1	TPO	A	224	11/12	0.93	0.18	96,103,114,116	0
3	TPO	C	224	11/12	0.93	0.15	97,104,120,126	0
2	TPO	B	224	11/12	0.96	0.12	86,95,116,120	0

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.