



Full wwPDB X-ray Structure Validation Report ⓘ

May 28, 2020 – 08:35 pm BST

PDB ID : 1U46
Title : Crystal Structure of the Unphosphorylated Kinase Domain of the Tyrosine Kinase ACK1
Authors : Loughheed, J.C.; Chen, R.H.; Mak, P.; Stout, T.J.
Deposited on : 2004-07-23
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 ⓘ symbol.

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

MolProbity	:	4.02b-467
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

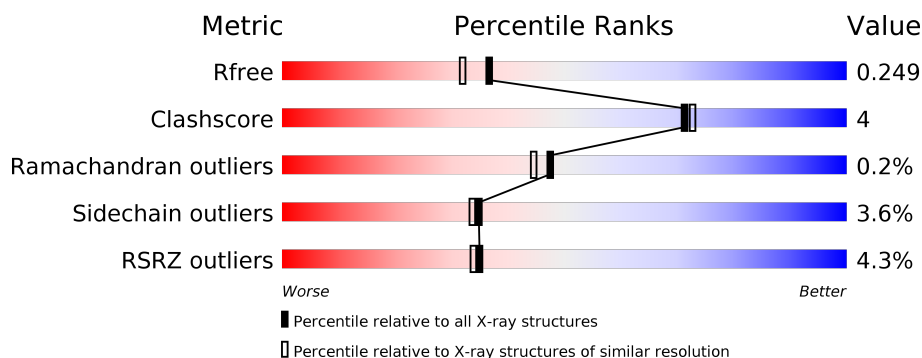
1 Overall quality at a glance

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	Whole archive (#Entries)	Similar resolution (#Entries, resolution range(Å))
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 $\leq 5\%$. 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	291	<div> <div>5%</div> <div> <div></div> <div>73%</div> <div>13%</div> <div>12%</div> </div> </div>
1	B	291	<div> <div>3%</div> <div> <div></div> <div>80%</div> <div>8%</div> <div>12%</div> </div> </div>

2 Entry composition

There are 3 unique types of molecules in this entry. The entry contains 4340 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 Activated CDC42 kinase 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	255	Total	C	N	O	S	0	0	0
			2044	1306	361	362	15			
1	B	257	Total	C	N	O	S	0	0	0
			2065	1315	369	366	15			

There are 8 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	107	GLY	-	CLONING ARTIFACT	UNP Q07912
A	108	SER	-	CLONING ARTIFACT	UNP Q07912
A	396	GLU	-	CLONING ARTIFACT	UNP Q07912
A	397	PHE	-	CLONING ARTIFACT	UNP Q07912
B	107	GLY	-	CLONING ARTIFACT	UNP Q07912
B	108	SER	-	CLONING ARTIFACT	UNP Q07912
B	396	GLU	-	CLONING ARTIFACT	UNP Q07912
B	397	PHE	-	CLONING ARTIFACT	UNP Q07912

- Molecule 2 is CHLORIDE ION (three-letter code: CL) (formula: Cl).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Cl	0	0
			1	1		

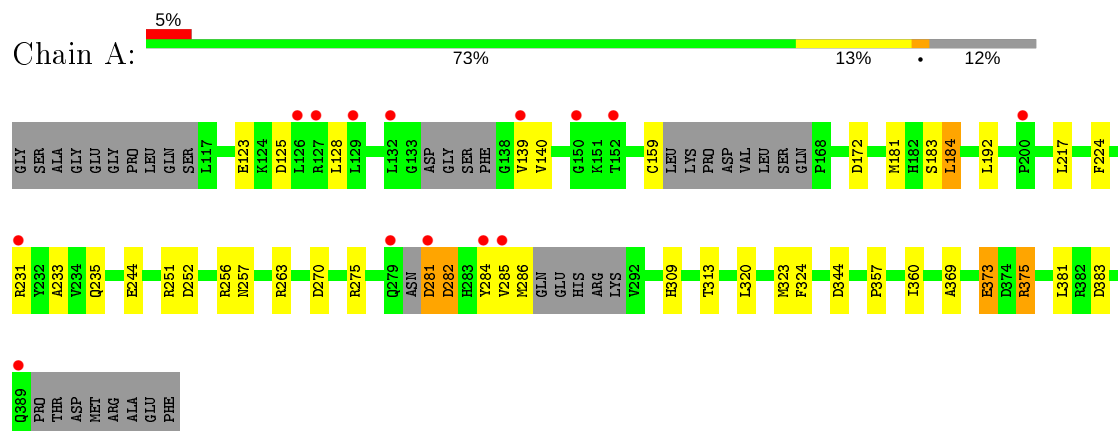
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	93	Total	O	0	0
			93	93		
3	B	137	Total	O	0	0
			137	137		

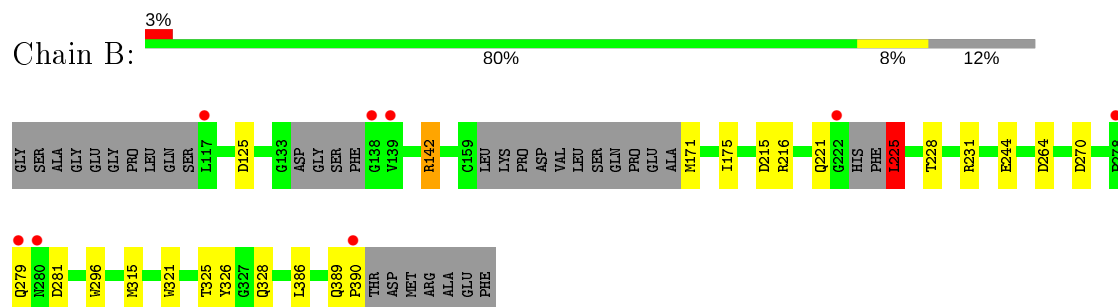
3 Residue-property plots [i](#)

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.

• Molecule 1: Activated CDC42 kinase 1



• Molecule 1: Activated CDC42 kinase 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, α , β , γ	81.18 Å 43.40 Å 84.45 Å 90.00° 111.82° 90.00°	Depositor
Resolution (Å)	30.00 – 2.00 29.90 – 2.00	Depositor EDS
% Data completeness (in resolution range)	100.0 (30.00-2.00) 100.0 (29.90-2.00)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.08	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.46 (at 2.00 Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, R_{free}	0.199 , 0.249 0.208 , 0.249	Depositor DCC
R_{free} test set	1886 reflections (5.04%)	wwPDB-VP
Wilson B-factor (Å ²)	26.6	Xtriage
Anisotropy	0.025	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.34 , 53.3	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	0.019 for l,-k,h	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	4340	wwPDB-VP
Average B, all atoms (Å ²)	32.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 5.78% 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: CL

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	0.84	1/2090 (0.0%)	1.07	13/2823 (0.5%)
1	B	0.91	1/2111 (0.0%)	1.01	8/2852 (0.3%)
All	All	0.87	2/4201 (0.0%)	1.04	21/5675 (0.4%)

All (2) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	244	GLU	CD-OE1	6.64	1.32	1.25
1	A	323	MET	SD-CE	-5.18	1.48	1.77

All (21) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	375	ARG	NE-CZ-NH1	18.95	129.78	120.30
1	A	375	ARG	NE-CZ-NH2	-13.76	113.42	120.30
1	A	270	ASP	CB-CG-OD2	13.05	130.04	118.30
1	B	142	ARG	NE-CZ-NH2	-10.86	114.87	120.30
1	B	231	ARG	NE-CZ-NH2	-6.87	116.87	120.30
1	B	125	ASP	CB-CG-OD2	6.66	124.30	118.30
1	B	142	ARG	NE-CZ-NH1	6.24	123.42	120.30
1	B	215	ASP	CB-CG-OD2	6.24	123.91	118.30
1	B	264	ASP	CB-CG-OD1	6.23	123.91	118.30
1	B	231	ARG	NE-CZ-NH1	6.12	123.36	120.30
1	A	256	ARG	NE-CZ-NH1	6.03	123.31	120.30
1	A	256	ARG	NE-CZ-NH2	-5.99	117.31	120.30
1	A	281	ASP	CB-CG-OD2	5.81	123.53	118.30
1	B	225	LEU	CB-CG-CD1	-5.67	101.36	111.00
1	A	282	ASP	CB-CG-OD2	5.61	123.35	118.30
1	A	383	ASP	CB-CG-OD2	5.48	123.23	118.30
1	A	184	LEU	CA-CB-CG	5.39	127.69	115.30

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	375	ARG	CD-NE-CZ	5.22	130.90	123.60
1	A	125	ASP	CB-CG-OD2	5.20	122.98	118.30
1	A	172	ASP	CB-CG-OD2	5.19	122.97	118.30
1	A	344	ASP	CB-CG-OD2	5.03	122.83	118.30

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	2044	0	2045	18	0
1	B	2065	0	2071	10	0
2	B	1	0	0	0	0
3	A	93	0	0	1	0
3	B	137	0	0	0	0
All	All	4340	0	4116	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.

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:B:225:LEU:HD11	1:B:326:TYR:HE1	1.16	1.09
1:B:225:LEU:HD11	1:B:326:TYR:CE1	1.97	0.98
1:A:231:ARG:HD3	1:A:235:GLN:NE2	2.05	0.69
1:B:225:LEU:CD1	1:B:326:TYR:HE1	2.03	0.61
1:B:225:LEU:CD1	1:B:326:TYR:CE1	2.78	0.59
1:B:386:LEU:HA	1:B:390:PRO:HB3	1.89	0.55
1:A:231:ARG:HD3	1:A:235:GLN:HE21	1.70	0.55
1:A:373:GLU:HG3	3:A:452:HOH:O	2.08	0.54
1:A:369:ALA:O	1:A:375:ARG:HD2	2.09	0.52
1:A:233:ALA:HB1	1:A:320:LEU:HD21	1.92	0.51
1:B:225:LEU:HD12	1:B:225:LEU:N	2.26	0.50

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:357:PRO:HG2	1:A:360:ILE:HD12	1.95	0.49
1:B:171:MET:CE	1:B:175:ILE:HD11	2.43	0.48
1:A:181:MET:HA	1:A:184:LEU:HD22	1.96	0.48
1:A:285:VAL:HG12	1:A:286:MET:O	2.13	0.47
1:A:275:ARG:HG3	1:A:284:TYR:CE1	2.50	0.46
1:A:324:PHE:CZ	1:A:360:ILE:HD13	2.52	0.44
1:A:244:GLU:HA	1:A:309:HIS:CE1	2.52	0.44
1:B:225:LEU:HB2	1:B:228:THR:HB	2.01	0.43
1:A:217:LEU:HD23	1:A:224:PHE:CD2	2.54	0.43
1:A:128:LEU:HD12	1:A:128:LEU:H	1.84	0.43
1:A:313:THR:HG23	1:A:381:LEU:HD11	2.00	0.42
1:A:252:ASP:O	1:A:257:ASN:ND2	2.50	0.42
1:A:251:ARG:NH2	1:A:286:MET:CE	2.84	0.41
1:B:296:TRP:O	1:B:315:MET:HG2	2.20	0.41
1:A:139:VAL:HG12	1:A:140:VAL:N	2.36	0.41
1:B:321:TRP:O	1:B:325:THR:HG23	2.21	0.41
1:A:231:ARG:CD	1:A:235:GLN:NE2	2.80	0.40

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	245/291 (84%)	237 (97%)	8 (3%)	0	100	100
1	B	249/291 (86%)	242 (97%)	6 (2%)	1 (0%)	34	30
All	All	494/582 (85%)	479 (97%)	14 (3%)	1 (0%)	47	44

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	270	ASP

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	220/250 (88%)	212 (96%)	8 (4%)	35	34
1	B	223/250 (89%)	215 (96%)	8 (4%)	35	34
All	All	443/500 (89%)	427 (96%)	16 (4%)	35	34

All (16) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	123	GLU
1	A	159	CYS
1	A	183	SER
1	A	192	LEU
1	A	263	ARG
1	A	281	ASP
1	A	282	ASP
1	A	373	GLU
1	B	142	ARG
1	B	216	ARG
1	B	221	GLN
1	B	225	LEU
1	B	279	GLN
1	B	281	ASP
1	B	328	GLN
1	B	389	GLN

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (5) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	179	ASN
1	A	283	HIS
1	A	328	GLN
1	B	370	HIS
1	B	389	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](#)

Of 1 ligands modelled in this entry, 1 is 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 [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	255/291 (87%)	0.25	14 (5%)	25 24	15, 33, 57, 70	19 (7%)
1	B	257/291 (88%)	-0.01	8 (3%)	49 48	15, 28, 49, 60	13 (5%)
All	All	512/582 (87%)	0.12	22 (4%)	35 34	15, 30, 54, 70	32 (6%)

All (22) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	280	ASN	4.8
1	B	138	GLY	4.7
1	A	200	PRO	4.4
1	A	389	GLN	4.2
1	A	281	ASP	3.6
1	B	390	PRO	3.5
1	B	117	LEU	3.4
1	B	139	VAL	3.2
1	A	152	THR	3.1
1	A	127	ARG	3.0
1	A	129	LEU	2.9
1	A	150	GLY	2.7
1	A	132	LEU	2.7
1	A	284	TYR	2.5
1	A	285	VAL	2.4
1	A	126	LEU	2.4
1	A	139	VAL	2.4
1	B	279	GLN	2.3
1	A	279	GLN	2.3
1	A	231	ARG	2.2
1	B	222	GLY	2.1
1	B	278	PRO	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. 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(\AA^2)	Q<0.9
2	CL	B	1	1/1	0.99	0.04	29,29,29,29	0

6.5 Other polymers [i](#)

There are no such residues in this entry.