



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 12, 2026 – 11:25 AM UTC

PDB ID : 1YHV / pdb_00001yhv
Title : Crystal Structure of PAK1 kinase domain with two point mutations (K299R, T423E)
Authors : Lei, M.; Robinson, M.A.; Harrison, S.C.
Deposited on : 2005-01-10
Resolution : 1.80 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 types of validation reports are described at

<http://www.wwpdb.org/validation/2017/FAQs#types>.

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

MolProbity	:	4-5-2 with Phenix2.0
Xtriage (Phenix)	:	2.0
EDS	:	3.0
Percentile statistics	:	20250101.v01 (using entries in the PDB archive January 1st 2025)
CCP4	:	9.0.010 (Gargrove)
Density-Fitness	:	1.0.12
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	2.49

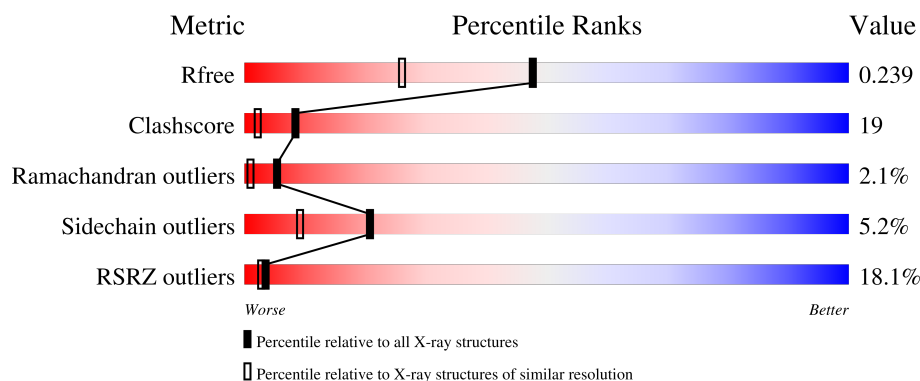
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 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 (#Entries)	Similar resolution (#Entries, resolution range(Å))
R_{free}	180053	7662 (1.80-1.80)
Clashscore	190562	8479 (1.80-1.80)
Ramachandran outliers	187476	8391 (1.80-1.80)
Sidechain outliers	187428	8390 (1.80-1.80)
RSRZ outliers	180081	7663 (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 $\leq 5\%$. 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	297	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2488 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 PAK 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	293	Total	C	N	O	S	0	0	0
			2270	1440	376	438	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	299	ARG	LYS	engineered mutation	UNP Q13153
A	423	GLU	THR	engineered mutation	UNP Q13153

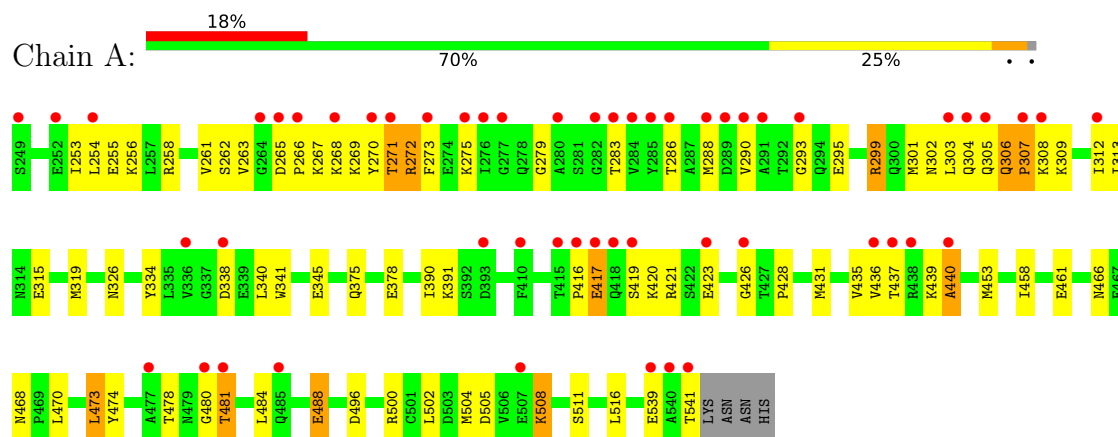
- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	218	Total	O	0	0
			218	218		

3 Residue-property plots

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: Serine/threonine-protein kinase PAK 1



4 Data and refinement statistics

Property	Value	Source
Space group	C 2 2 21	Depositor
Cell constants a, b, c, α , β , γ	51.96Å 103.30Å 122.88Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	50.00 – 1.80 50.00 – 1.80	Depositor EDS
% Data completeness (in resolution range)	(Not available) (50.00-1.80) 98.1 (50.00-1.80)	Depositor EDS
R_{merge}	0.08	Depositor
R_{sym}	0.25	Depositor
$\langle I/\sigma(I) \rangle$ ¹	4.12 (at 1.80Å)	Xtriage
Refinement program	CNS	Depositor
R, R_{free}	0.220 , 0.240 0.223 , 0.239	Depositor DCC
R_{free} test set	1514 reflections (4.96%)	wwPDB-VP
Wilson B-factor (Å ²)	27.5	Xtriage
Anisotropy	0.572	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.38 , 42.1	EDS
L-test for twinning ²	$\langle L \rangle = 0.49$, $\langle L^2 \rangle = 0.32$	Xtriage
Estimated twinning fraction	No twinning to report.	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	2488	wwPDB-VP
Average B, all atoms (Å ²)	36.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.09% 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](#)

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	0.53	0/2308	1.07	12/3127 (0.4%)

There are no bond length outliers.

All (12) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	306	GLN	OE1-CD-NE2	-6.88	115.72	122.60
1	A	263	VAL	N-CA-C	6.72	117.37	110.05
1	A	262	SER	N-CA-C	-6.36	101.96	110.55
1	A	261	VAL	N-CA-C	6.33	117.37	110.21
1	A	345	GLU	N-CA-C	-5.85	102.81	110.53
1	A	417	GLU	CA-C-N	5.63	130.53	122.08
1	A	417	GLU	C-N-CA	5.63	130.53	122.08
1	A	511	SER	N-CA-C	-5.38	102.79	110.59
1	A	338	ASP	N-CA-C	-5.32	106.38	112.92
1	A	481	THR	CA-C-N	5.21	125.12	119.85
1	A	481	THR	C-N-CA	5.21	125.12	119.85
1	A	458	ILE	CB-CA-C	-5.09	105.25	111.92

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	2270	0	2277	87	0

Continued on next page...

Continued from previous page...

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
2	A	218	0	0	10	0
All	All	2488	0	2277	87	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 19.

All (87) 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:A:309:LYS:O	1:A:312:ILE:HG22	1.55	1.04
1:A:271:THR:HG22	1:A:272:ARG:N	1.73	1.00
1:A:391:LYS:HA	1:A:453:MET:HE2	1.51	0.90
1:A:436:VAL:HG21	1:A:473:LEU:HD23	1.53	0.90
1:A:391:LYS:CA	1:A:453:MET:HE2	2.02	0.89
1:A:293:GLY:HA2	2:A:170:HOH:O	1.79	0.82
1:A:481:THR:HG22	1:A:502:LEU:O	1.81	0.81
1:A:266:PRO:HG2	1:A:341:TRP:CE3	2.17	0.79
1:A:474:TYR:O	1:A:478:THR:HG23	1.88	0.74
1:A:436:VAL:CG2	1:A:473:LEU:HD23	2.17	0.73
1:A:271:THR:HG22	1:A:272:ARG:H	1.54	0.72
1:A:271:THR:CG2	1:A:272:ARG:N	2.47	0.71
1:A:481:THR:HG23	1:A:504:MET:HG2	1.72	0.70
1:A:426:GLY:O	1:A:431:MET:HE2	1.92	0.69
1:A:539:GLU:C	1:A:541:THR:H	1.99	0.69
1:A:306:GLN:HG3	2:A:208:HOH:O	1.92	0.69
1:A:436:VAL:HG21	1:A:473:LEU:CD2	2.24	0.67
1:A:307:PRO:HB2	1:A:308:LYS:HD2	1.76	0.66
1:A:275:LYS:HE3	1:A:283:THR:HG23	1.79	0.65
1:A:378:GLU:OE1	1:A:516:LEU:CD1	2.44	0.64
1:A:288:MET:HG3	2:A:116:HOH:O	1.97	0.64
1:A:468:ASN:OD1	1:A:470:LEU:HG	1.98	0.63
1:A:391:LYS:CA	1:A:453:MET:CE	2.75	0.63
1:A:288:MET:HE2	2:A:116:HOH:O	1.98	0.62
1:A:326:ASN:ND2	1:A:375:GLN:HE21	2.01	0.59
1:A:326:ASN:HD21	1:A:375:GLN:HE21	1.49	0.58
1:A:309:LYS:O	1:A:312:ILE:CG2	2.42	0.57
1:A:312:ILE:CD1	1:A:340:LEU:HD23	2.35	0.57
1:A:436:VAL:CG2	1:A:473:LEU:CD2	2.81	0.57
1:A:266:PRO:HG2	1:A:341:TRP:CD2	2.40	0.57
1:A:378:GLU:OE1	1:A:516:LEU:HD13	2.04	0.57
1:A:390:ILE:HG22	1:A:453:MET:CE	2.35	0.56

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:255:GLU:OE1	1:A:258:ARG:NH1	2.38	0.56
1:A:265:ASP:OD2	1:A:267:LYS:HB2	2.06	0.56
1:A:391:LYS:C	1:A:453:MET:HE2	2.32	0.55
1:A:271:THR:O	1:A:273:PHE:HD1	1.90	0.55
1:A:474:TYR:CE1	1:A:478:THR:HG21	2.41	0.54
1:A:428:PRO:N	1:A:431:MET:HE3	2.22	0.54
1:A:391:LYS:N	1:A:453:MET:HE1	2.22	0.53
1:A:312:ILE:HD11	1:A:340:LEU:HD23	1.90	0.53
1:A:269:LYS:HD3	1:A:270:TYR:CZ	2.44	0.52
1:A:428:PRO:HA	1:A:431:MET:HE3	1.92	0.52
1:A:266:PRO:HG3	1:A:334:TYR:CG	2.45	0.51
1:A:253:ILE:HD13	1:A:309:LYS:HB2	1.93	0.50
1:A:488:GLU:H	1:A:488:GLU:CD	2.18	0.50
1:A:390:ILE:HG22	1:A:453:MET:HE3	1.93	0.50
1:A:500:ARG:HD3	2:A:90:HOH:O	2.12	0.49
1:A:254:LEU:O	1:A:258:ARG:HG3	2.13	0.48
1:A:256:LYS:HG2	1:A:313:ILE:HD13	1.95	0.48
1:A:268:LYS:HE3	2:A:168:HOH:O	2.11	0.48
1:A:391:LYS:N	1:A:453:MET:CE	2.76	0.48
1:A:301:MET:HE1	1:A:312:ILE:HD11	1.96	0.47
1:A:326:ASN:HD21	1:A:375:GLN:NE2	2.11	0.47
1:A:267:LYS:HE2	1:A:341:TRP:CH2	2.50	0.47
1:A:309:LYS:C	1:A:312:ILE:HG22	2.34	0.46
1:A:265:ASP:OD2	1:A:267:LYS:CB	2.63	0.46
1:A:428:PRO:CA	1:A:431:MET:HE3	2.45	0.46
1:A:505:ASP:HB3	1:A:508:LYS:HB2	1.98	0.46
1:A:480:GLY:HA2	1:A:504:MET:SD	2.56	0.46
1:A:426:GLY:C	1:A:431:MET:HE2	2.42	0.45
1:A:439:LYS:O	1:A:440:ALA:C	2.59	0.45
1:A:496:ASP:O	1:A:500:ARG:HG2	2.17	0.45
1:A:267:LYS:HE2	1:A:341:TRP:HH2	1.82	0.45
1:A:302:ASN:HB3	1:A:305:GLN:HG2	2.00	0.44
1:A:303:LEU:HD22	1:A:309:LYS:HD3	1.99	0.44
1:A:468:ASN:OD1	1:A:470:LEU:CG	2.65	0.44
1:A:303:LEU:O	1:A:305:GLN:N	2.51	0.44
1:A:391:LYS:C	1:A:453:MET:CE	2.90	0.44
1:A:436:VAL:HG23	2:A:111:HOH:O	2.17	0.44
1:A:481:THR:CG2	1:A:504:MET:HG2	2.46	0.43
1:A:266:PRO:HG3	1:A:334:TYR:CB	2.48	0.43
1:A:306:GLN:HA	1:A:307:PRO:HD2	1.82	0.43
1:A:436:VAL:O	1:A:437:THR:C	2.60	0.43

Continued on next page...

Continued from previous page...

Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:286:THR:HG21	1:A:295:GLU:OE1	2.19	0.43
1:A:315:GLU:O	1:A:319:MET:HG3	2.18	0.43
1:A:312:ILE:HG23	1:A:313:ILE:N	2.34	0.42
1:A:421:ARG:O	1:A:440:ALA:HA	2.20	0.42
1:A:303:LEU:C	1:A:305:GLN:N	2.78	0.42
1:A:315:GLU:HG2	1:A:319:MET:HE2	2.00	0.42
1:A:390:ILE:HG22	1:A:453:MET:HE1	2.01	0.42
1:A:453:MET:HE3	1:A:453:MET:HB2	1.78	0.42
1:A:299:ARG:HD2	1:A:301:MET:SD	2.60	0.42
1:A:312:ILE:HD13	1:A:340:LEU:HD23	2.01	0.42
1:A:481:THR:HG21	2:A:173:HOH:O	2.18	0.42
1:A:390:ILE:CG2	1:A:453:MET:HE1	2.50	0.41
1:A:435:VAL:HB	2:A:111:HOH:O	2.20	0.41
1:A:431:MET:HB3	2:A:111:HOH:O	2.22	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 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	291/297 (98%)	273 (94%)	12 (4%)	6 (2%)	5 1

All (6) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	272	ARG
1	A	440	ALA
1	A	279	GLY
1	A	417	GLU
1	A	304	GLN
1	A	307	PRO

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	248/258 (96%)	235 (95%)	13 (5%)	21 9

All (13) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	271	THR
1	A	290	VAL
1	A	299	ARG
1	A	416	PRO
1	A	419	SER
1	A	420	LYS
1	A	423	GLU
1	A	461	GLU
1	A	466	ASN
1	A	473	LEU
1	A	484	LEU
1	A	488	GLU
1	A	508	LYS

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (3) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	326	ASN
1	A	383	ASN
1	A	418	GLN

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 [i](#)

There are no oligosaccharides 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	293/297 (98%)	0.94	53 (18%) 3 3	19, 33, 60, 66	0

All (53) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	480	GLY	6.0
1	A	417	GLU	5.6
1	A	540	ALA	4.8
1	A	290	VAL	4.6
1	A	437	THR	4.4
1	A	418	GLN	4.4
1	A	541	THR	4.4
1	A	265	ASP	4.3
1	A	440	ALA	4.0
1	A	273	PHE	4.0
1	A	426	GLY	3.7
1	A	419	SER	3.7
1	A	271	THR	3.7
1	A	276	ILE	3.6
1	A	438	ARG	3.5
1	A	275	LYS	3.2
1	A	304	GLN	3.2
1	A	307	PRO	3.1
1	A	481	THR	3.1
1	A	264	GLY	3.0
1	A	280	ALA	3.0
1	A	436	VAL	2.9
1	A	286	THR	2.9
1	A	283	THR	2.8
1	A	305	GLN	2.8
1	A	284	VAL	2.8
1	A	270	TYR	2.8

Continued on next page...

Continued from previous page...

Mol	Chain	Res	Type	RSRZ
1	A	293	GLY	2.8
1	A	308	LYS	2.7
1	A	266	PRO	2.6
1	A	288	MET	2.6
1	A	338	ASP	2.5
1	A	282	GLY	2.5
1	A	485	GLN	2.5
1	A	336	VAL	2.5
1	A	415	THR	2.4
1	A	268	LYS	2.4
1	A	254	LEU	2.3
1	A	539	GLU	2.3
1	A	277	GLY	2.3
1	A	416	PRO	2.3
1	A	410	PHE	2.3
1	A	393	ASP	2.2
1	A	249	SER	2.2
1	A	303	LEU	2.2
1	A	477	ALA	2.2
1	A	423	GLU	2.1
1	A	289	ASP	2.1
1	A	507	GLU	2.1
1	A	252	GLU	2.1
1	A	291	ALA	2.1
1	A	312	ILE	2.0
1	A	285	TYR	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 oligosaccharides 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.