



Full wwPDB X-ray Structure Validation Report ⓘ

Mar 6, 2026 – 01:24 PM UTC

PDB ID : 1L8K / pdb_00001l8k
Title : T Cell Protein-Tyrosine Phosphatase Structure
Authors : Iversen, L.F.
Deposited on : 2002-03-21
Resolution : 2.56 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) : **NOT EXECUTED**
EDS : **NOT EXECUTED**
Percentile statistics : 20250101.v01 (using entries in the PDB archive January 1st 2025)
Ideal geometry (proteins) : Engh & Huber (2001)
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP) : 2.49

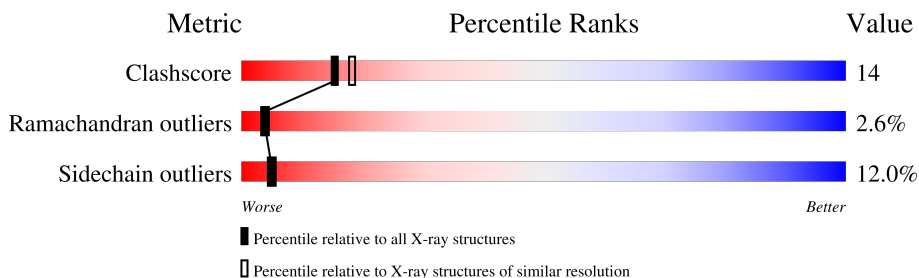
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.56 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))
Clashscore	190562	1897 (2.58-2.54)
Ramachandran outliers	187476	1875 (2.58-2.54)
Sidechain outliers	187428	1875 (2.58-2.54)

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 $\leq 5\%$.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	314	

2 Entry composition

There are 2 unique types of molecules in this entry. The entry contains 2284 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 T-cell protein-tyrosine phosphatase.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	273	Total	C	N	O	S	0	0	0
			2244	1422	392	418	12			

- Molecule 2 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	A	40	Total	O	0	0
			40	40		

4 Data and refinement statistics

Xtriage (Phenix) and EDS were not executed - this section is therefore incomplete.

Property	Value	Source
Space group	P 43 21 2	Depositor
Cell constants a, b, c, α , β , γ	60.49Å 60.49Å 187.63Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	20.00 – 2.56	Depositor
% Data completeness (in resolution range)	(Not available) (20.00-2.56)	Depositor
R_{merge}	(Not available)	Depositor
R_{sym}	(Not available)	Depositor
Refinement program	X-PLOR 3.851	Depositor
R, R_{free}	0.215 , 0.288	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2284	wwPDB-VP
Average B, all atoms (Å ²)	65.0	wwPDB-VP

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	0.47	0/2299	0.99	9/3114 (0.3%)

There are no bond length outliers.

All (9) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	120	SER	N-CA-C	-11.04	99.53	113.23
1	A	236	MET	N-CA-C	-6.56	100.19	110.36
1	A	150	VAL	N-CA-C	6.39	116.92	107.15
1	A	99	LEU	N-CA-C	-6.06	104.67	111.28
1	A	179	THR	N-CA-C	6.04	119.96	112.59
1	A	216	CYS	N-CA-C	-5.92	98.20	110.80
1	A	121	VAL	N-CA-C	-5.84	100.01	108.71
1	A	68	TYR	N-CA-C	5.25	117.79	109.24
1	A	96	HIS	N-CA-C	-5.10	105.80	111.36

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	2244	0	2189	60	0
2	A	40	0	0	1	0
All	All	2284	0	2189	60	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 14.

All (60) 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:109:VAL:HB	1:A:213:VAL:HG22	1.61	0.83
1:A:149:ASP:HB3	1:A:156:VAL:HG13	1.66	0.75
1:A:113:ASN:HD22	1:A:113:ASN:H	1.37	0.70
1:A:81:ARG:HD3	1:A:230:ASP:OD2	1.92	0.70
1:A:236:MET:C	1:A:238:LYS:H	2.01	0.66
1:A:46:ASN:ND2	1:A:70:ASN:HD22	1.95	0.64
1:A:227:SER:HB3	1:A:251:MET:HE1	1.79	0.64
1:A:47:ARG:H	1:A:87:GLN:HE22	1.44	0.64
1:A:46:ASN:HD22	1:A:70:ASN:HD22	1.45	0.62
1:A:236:MET:O	1:A:237:GLU:HB3	2.00	0.61
1:A:242:ILE:HD12	1:A:242:ILE:N	2.16	0.60
1:A:22:TYR:HB2	1:A:265:LEU:HD12	1.85	0.59
1:A:91:PRO:HD3	1:A:122:LYS:HB3	1.85	0.58
1:A:90:LEU:HD22	1:A:122:LYS:HG3	1.86	0.58
1:A:19:GLN:O	1:A:23:LEU:HB2	2.04	0.57
1:A:113:ASN:HD22	1:A:113:ASN:N	2.02	0.56
1:A:19:GLN:N	1:A:20:PRO:HD2	2.21	0.55
1:A:275:GLY:C	1:A:277:LYS:H	2.15	0.55
1:A:106:THR:HG21	1:A:213:VAL:HG13	1.90	0.54
1:A:236:MET:C	1:A:238:LYS:N	2.66	0.54
1:A:247:VAL:O	1:A:251:MET:HG3	2.06	0.54
1:A:229:VAL:HG12	1:A:272:ILE:HG13	1.89	0.54
1:A:230:ASP:O	1:A:234:VAL:HG23	2.07	0.54
1:A:20:PRO:O	1:A:24:GLU:HG3	2.08	0.53
1:A:133:GLU:CD	1:A:133:GLU:H	2.16	0.53
1:A:113:ASN:H	1:A:113:ASN:ND2	2.05	0.53
1:A:5:ILE:HD12	1:A:5:ILE:N	2.24	0.53
1:A:84:ILE:HG13	1:A:211:PRO:HB2	1.91	0.53
1:A:25:ILE:HG21	1:A:248:LEU:HD12	1.91	0.52
1:A:149:ASP:HB3	1:A:156:VAL:CG1	2.37	0.52
1:A:248:LEU:HD21	1:A:259:ILE:HD13	1.91	0.52
1:A:268:SER:O	1:A:272:ILE:HG13	2.12	0.49
1:A:196:LEU:O	1:A:199:VAL:HG22	2.13	0.49
1:A:189:PRO:HD3	1:A:270:MET:HE2	1.95	0.49
1:A:198:LYS:HD3	1:A:198:LYS:O	2.13	0.49
1:A:154:TYR:OH	1:A:198:LYS:HG3	2.13	0.48
1:A:258:LEU:O	1:A:259:ILE:HB	2.13	0.48

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Atom-1	Atom-2	Interatomic distance (Å)	Clash overlap (Å)
1:A:110:VAL:HG22	1:A:214:ILE:HG23	1.94	0.48
1:A:160:GLN:NE2	1:A:169:THR:HG21	2.28	0.47
1:A:110:VAL:HG13	1:A:214:ILE:HD13	1.96	0.47
1:A:95:CYS:SG	1:A:138:GLU:HG3	2.54	0.47
1:A:14:THR:C	1:A:16:ARG:H	2.22	0.46
1:A:105:LYS:HA	1:A:170:ARG:NH1	2.29	0.46
1:A:47:ARG:HH12	1:A:122:LYS:HB2	1.81	0.46
1:A:90:LEU:CD2	1:A:122:LYS:HG3	2.46	0.45
1:A:104:GLN:O	1:A:210:GLY:HA3	2.17	0.45
1:A:5:ILE:HA	2:A:326:HOH:O	2.15	0.44
1:A:269:TYR:O	1:A:273:ILE:HG13	2.17	0.44
1:A:19:GLN:N	1:A:20:PRO:CD	2.81	0.44
1:A:225:THR:O	1:A:229:VAL:HG22	2.17	0.44
1:A:76:ILE:HD11	1:A:251:MET:HE3	2.00	0.44
1:A:32:TYR:OH	1:A:253:LYS:HD2	2.19	0.43
1:A:140:GLY:C	1:A:164:ILE:HG12	2.43	0.43
1:A:189:PRO:HB2	1:A:193:LEU:HD23	2.00	0.43
1:A:189:PRO:O	1:A:193:LEU:HD23	2.19	0.42
1:A:238:LYS:HG2	1:A:239:GLY:N	2.34	0.42
1:A:106:THR:HG23	1:A:212:ALA:HA	2.02	0.42
1:A:16:ARG:HE	1:A:16:ARG:HA	1.84	0.41
1:A:109:VAL:HB	1:A:213:VAL:CG2	2.42	0.41
1:A:99:LEU:HD23	1:A:141:PHE:HE2	1.84	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	271/314 (86%)	242 (89%)	22 (8%)	7 (3%)	4 4

All (7) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	122	LYS
1	A	183	PHE
1	A	276	ALA
1	A	15	GLN
1	A	123	CYS
1	A	259	ILE
1	A	118	LYS

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	250/289 (86%)	220 (88%)	30 (12%)	5 5

All (30) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	21	LEU
1	A	23	LEU
1	A	30	HIS
1	A	50	ASP
1	A	75	ASP
1	A	81	ARG
1	A	83	TYR
1	A	99	LEU
1	A	101	VAL
1	A	109	VAL
1	A	110	VAL
1	A	113	ASN
1	A	115	ILE
1	A	119	GLU
1	A	121	VAL
1	A	133	GLU
1	A	138	GLU
1	A	145	LEU
1	A	156	VAL
1	A	183	PHE

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Mol	Chain	Res	Type
1	A	185	VAL
1	A	193	LEU
1	A	213	VAL
1	A	216	CYS
1	A	222	ARG
1	A	228	LEU
1	A	233	LEU
1	A	248	LEU
1	A	258	LEU
1	A	265	LEU

Sometimes sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (16) such sidechains are listed below:

Mol	Chain	Res	Type
1	A	15	GLN
1	A	19	GLN
1	A	34	HIS
1	A	46	ASN
1	A	63	ASN
1	A	80	GLN
1	A	87	GLN
1	A	113	ASN
1	A	125	GLN
1	A	132	GLN
1	A	157	HIS
1	A	160	GLN
1	A	165	ASN
1	A	194	ASN
1	A	243	ASN
1	A	246	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

EDS was not executed - this section is therefore empty.

6.2 Non-standard residues in protein, DNA, RNA chains

EDS was not executed - this section is therefore empty.

6.3 Carbohydrates

EDS was not executed - this section is therefore empty.

6.4 Ligands

EDS was not executed - this section is therefore empty.

6.5 Other polymers

EDS was not executed - this section is therefore empty.