



# Full wwPDB X-ray Structure Validation Report ⓘ

Jan 31, 2016 – 07:10 PM GMT

PDB ID : 1ECV  
Title : CRYSTAL STRUCTURE OF PROTEIN TYROSINE PHOSPHATASE 1B  
COMPLEXED WITH 5-iodo-2-(oxalyl-amino)-benzoic acid  
Authors : Andersen, H.S.; Iversen, L.F.; Branner, S.; Rasmussen, H.B.; Moller, N.P.H.  
Deposited on : 2000-01-26  
Resolution : 1.95 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.

---

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

MolProbity : 4.02b-467  
Mogul : 1.7 (RC4), CSD as536be (2015)  
Xtriage (Phenix) : **NOT EXECUTED**  
EDS : **NOT EXECUTED**  
Percentile statistics : 20151230.v01 (using entries in the PDB archive December 30th 2015)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et al. (1996)  
Validation Pipeline (wwPDB-VP) : trunk26865

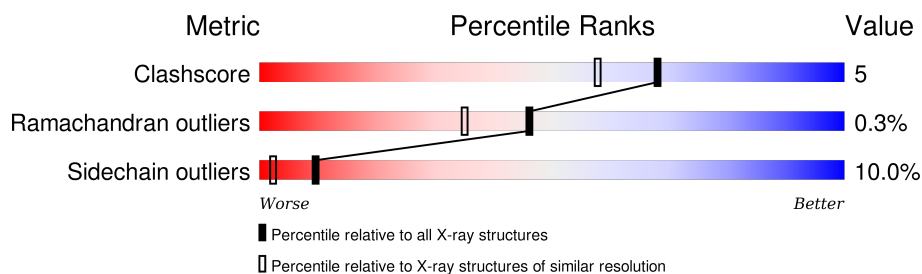
# 1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

*X-RAY DIFFRACTION*

The reported resolution of this entry is 1.95 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	102246	1953 (1.96-1.96)
Ramachandran outliers	100387	1936 (1.96-1.96)
Sidechain outliers	100360	1936 (1.96-1.96)

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.

Note EDS was not executed.

Mol	Chain	Length	Quality of chain
1	A	298	 75% 16% 8% •

## 2 Entry composition [i](#)

There are 4 unique types of molecules in this entry. The entry contains 2724 atoms, of which 28 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.

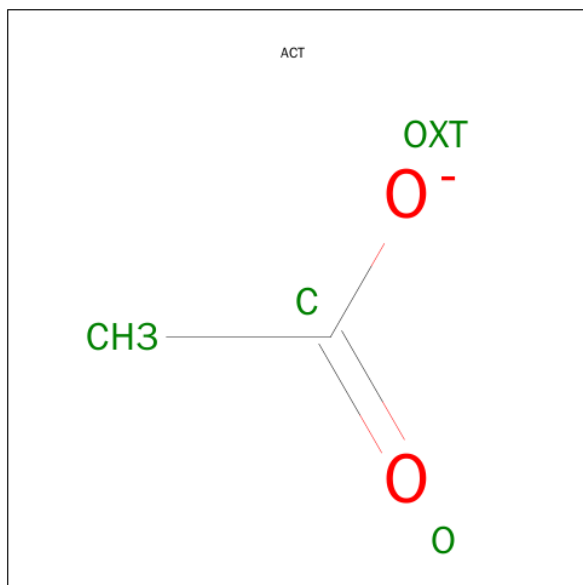
- Molecule 1 is a protein called PROTEIN-TYROSINE PHOSPHATASE 1B.

Mol	Chain	Residues	Atoms						ZeroOcc	AltConf	Trace
1	A	297	Total	C	H	N	O	S	0	0	0
			2455	1535	28	418	458	16			

There are 2 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	151	THR	SER	CONFLICT	UNP P18031
A	252	ASP	GLU	CONFLICT	UNP P18031

- Molecule 2 is ACETATE ION (three-letter code: ACT) (formula:  $C_2H_3O_2$ ).



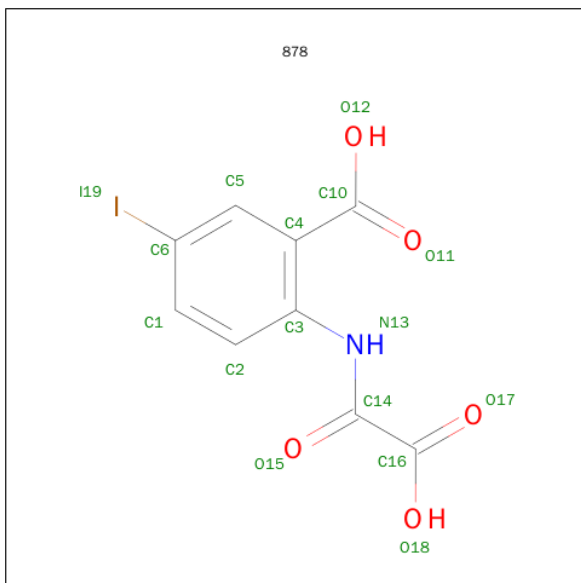
Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		
2	A	1	Total	C	O	0	0
			4	2	2		

*Continued on next page...*

Continued from previous page...

Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	C	O	0	0
			4	2	2		

- Molecule 3 is 5-iodo-2-(oxalyl-amino)-benzoic acid (three-letter code: 878) (formula: C<sub>9</sub>H<sub>6</sub>INO<sub>5</sub>).



Mol	Chain	Residues	Atoms					ZeroOcc	AltConf
3	A	1	Total	C	I	N	O	0	0
			16	9	1	1	5		

- Molecule 4 is water.

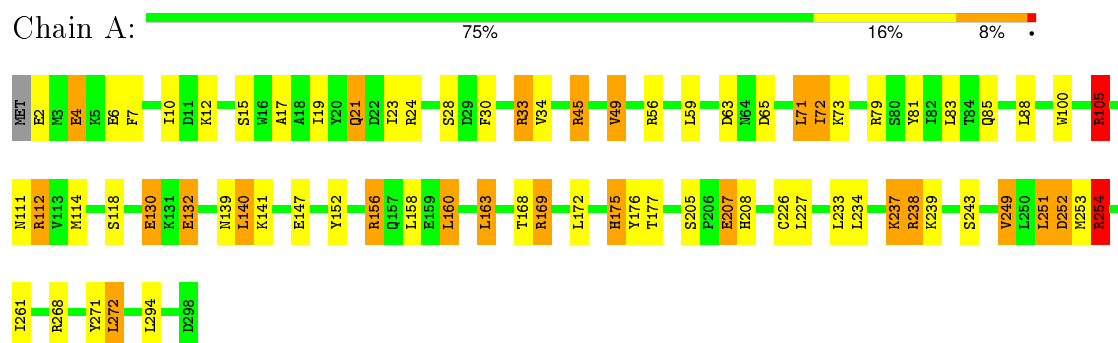
Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	241	Total	O	0	0
			241	241		

### 3 Residue-property plots

These plots are drawn for all protein, RNA and DNA chains in the entry. The first graphic for a chain summarises the proportions of errors 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 ( $\text{RSRZ} > 2$ ). 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.

Note EDS was not executed.

#### • Molecule 1: PROTEIN-TYROSINE PHOSPHATASE 1B



## 4 Data and refinement statistics

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

Property	Value	Source
Space group	P 31 2 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	87.87Å 87.87Å 103.12Å 90.00° 90.00° 120.00°	Depositor
Resolution (Å)	6.00 – 1.95	Depositor
% Data completeness (in resolution range)	99.5 (6.00-1.95)	Depositor
$R_{merge}$	0.06	Depositor
$R_{sym}$	(Not available)	Depositor
Refinement program	X-PLOR	Depositor
R, $R_{free}$	0.194 , 0.249	Depositor
Estimated twinning fraction	No twinning to report.	Xtriage
Total number of atoms	2724	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	21.0	wwPDB-VP

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: 878, ACT

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	1.08	3/2482 (0.1%)	1.72	42/3345 (1.3%)

Chiral center outliers are detected by calculating the chiral volume of a chiral center and verifying if the center is modelled as a planar moiety or with the opposite hand. A planarity outlier is detected by checking planarity of atoms in a peptide group, atoms in a mainchain group or atoms of a sidechain that are expected to be planar.

Mol	Chain	#Chirality outliers	#Planarity outliers
1	A	0	2

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	21	GLN	CG-CD	5.89	1.64	1.51
1	A	252	ASP	CB-CG	5.35	1.62	1.51
1	A	24	ARG	NE-CZ	5.21	1.39	1.33

All (42) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	254	ARG	NE-CZ-NH1	18.15	129.38	120.30
1	A	254	ARG	NE-CZ-NH2	-16.05	112.28	120.30
1	A	33	ARG	NE-CZ-NH2	-12.05	114.27	120.30
1	A	238	ARG	NE-CZ-NH2	-10.97	114.82	120.30
1	A	79	ARG	NE-CZ-NH1	10.68	125.64	120.30
1	A	45	ARG	NE-CZ-NH1	10.10	125.35	120.30
1	A	169	ARG	NE-CZ-NH2	-9.65	115.47	120.30
1	A	56	ARG	NE-CZ-NH1	9.55	125.07	120.30
1	A	238	ARG	NE-CZ-NH1	9.49	125.05	120.30
1	A	268	ARG	NE-CZ-NH2	-8.17	116.22	120.30

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	24	ARG	CD-NE-CZ	7.67	134.34	123.60
1	A	252	ASP	CB-CG-OD2	7.45	125.01	118.30
1	A	254	ARG	CD-NE-CZ	7.42	133.98	123.60
1	A	176	TYR	CB-CG-CD2	-7.36	116.59	121.00
1	A	130	GLU	CB-CA-C	-7.20	96.00	110.40
1	A	71	LEU	CB-CG-CD2	6.93	122.79	111.00
1	A	112	ARG	NE-CZ-NH2	-6.83	116.88	120.30
1	A	254	ARG	CB-CG-CD	6.53	128.59	111.60
1	A	118	SER	CB-CA-C	-6.46	97.83	110.10
1	A	156	ARG	NE-CZ-NH2	-6.39	117.11	120.30
1	A	141	LYS	CB-CG-CD	-6.37	95.04	111.60
1	A	132	GLU	CA-CB-CG	6.36	127.40	113.40
1	A	268	ARG	NE-CZ-NH1	6.31	123.45	120.30
1	A	65	ASP	CB-CG-OD1	6.30	123.97	118.30
1	A	114	MET	CG-SD-CE	-6.29	90.13	100.20
1	A	140	LEU	CB-CG-CD1	6.28	121.67	111.00
1	A	81	TYR	CB-CG-CD1	-6.27	117.24	121.00
1	A	34	VAL	CA-CB-CG2	6.11	120.07	110.90
1	A	207	GLU	CA-CB-CG	6.03	126.65	113.40
1	A	105	ARG	NE-CZ-NH1	5.98	123.29	120.30
1	A	56	ARG	NE-CZ-NH2	-5.95	117.33	120.30
1	A	253	MET	CG-SD-CE	-5.77	90.97	100.20
1	A	271	TYR	CB-CG-CD2	-5.59	117.64	121.00
1	A	72	ILE	CA-CB-CG1	5.44	121.34	111.00
1	A	152	TYR	CB-CG-CD2	-5.39	117.77	121.00
1	A	49	VAL	CA-CB-CG1	5.21	118.72	110.90
1	A	252	ASP	OD1-CG-OD2	-5.17	113.47	123.30
1	A	4	GLU	CA-CB-CG	5.15	124.72	113.40
1	A	33	ARG	NE-CZ-NH1	5.07	122.83	120.30
1	A	33	ARG	CD-NE-CZ	5.03	130.64	123.60
1	A	249	VAL	CA-CB-CG2	5.01	118.42	110.90
1	A	208	HIS	CA-CB-CG	-5.01	105.08	113.60

There are no chirality outliers.

All (2) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	105	ARG	Sidechain
1	A	254	ARG	Sidechain



## 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	2427	28	2381	23	0
2	A	12	0	9	0	0
3	A	16	0	4	0	0
4	A	241	0	0	2	0
All	All	2696	28	2394	23	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 5.

All (23) 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:100:TRP:HE3	1:A:160:LEU:HD13	1.62	0.64
1:A:147:GLU:HB3	1:A:156:ARG:HG2	1.82	0.61
1:A:205:SER:HB3	1:A:207:GLU:OE1	2.00	0.61
1:A:45:ARG:H	1:A:85:GLN:HE22	1.48	0.59
1:A:205:SER:HB3	1:A:207:GLU:CD	2.30	0.52
1:A:238:ARG:HD2	1:A:243:SER:OG	2.10	0.52
1:A:111:ASN:O	1:A:175:HIS:HE1	1.93	0.51
1:A:45:ARG:H	1:A:85:GLN:NE2	2.09	0.51
1:A:239:LYS:HE2	4:A:386:HOH:O	2.09	0.51
1:A:100:TRP:CZ3	1:A:169:ARG:HG3	2.47	0.50
1:A:112:ARG:HD3	1:A:177:THR:O	2.14	0.48
1:A:23:ILE:HG23	1:A:251:LEU:HD13	1.96	0.47
1:A:10:ILE:HD13	1:A:19:ILE:HG13	1.95	0.47
1:A:6:GLU:O	1:A:10:ILE:HG13	2.16	0.45
1:A:139:ASN:O	1:A:163:LEU:HB2	2.16	0.45
1:A:73:LYS:HD3	4:A:530:HOH:O	2.16	0.44
1:A:83:LEU:HD11	1:A:226:CYS:SG	2.58	0.44
1:A:7:PHE:CZ	1:A:272:LEU:HD13	2.53	0.43
1:A:237:LYS:HB3	1:A:237:LYS:HE2	1.69	0.43
1:A:28:SER:HB3	1:A:30:PHE:CE2	2.54	0.42
1:A:2:GLU:HG3	1:A:4:GLU:H	1.85	0.42
1:A:227:LEU:HD12	1:A:249:VAL:HG22	2.03	0.41
1:A:17:ALA:O	1:A:21:GLN:HG3	2.21	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	295/298 (99%)	288 (98%)	6 (2%)	1 (0%)	46 35

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	261	ILE

### 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	270/271 (100%)	243 (90%)	27 (10%)	9 2

All (27) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	12	LYS
1	A	15	SER
1	A	33	ARG
1	A	49	VAL
1	A	59	LEU
1	A	63	ASP
1	A	71	LEU

*Continued on next page...*

*Continued from previous page...*

Mol	Chain	Res	Type
1	A	72	ILE
1	A	88	LEU
1	A	105	ARG
1	A	130	GLU
1	A	132	GLU
1	A	140	LEU
1	A	158	LEU
1	A	160	LEU
1	A	163	LEU
1	A	168	THR
1	A	172	LEU
1	A	175	HIS
1	A	233	LEU
1	A	234	LEU
1	A	237	LYS
1	A	251	LEU
1	A	252	ASP
1	A	254	ARG
1	A	272	LEU
1	A	294	LEU

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	85	GLN
1	A	102	GLN
1	A	111	ASN
1	A	123	GLN
1	A	175	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

4 ligands 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
3	878	A	301	-	10,16,16	1.02	1 (10%)	12,22,22	1.10	1 (8%)
2	ACT	A	302	-	1,3,3	3.22	1 (100%)	0,3,3	0.00	-
2	ACT	A	303	-	1,3,3	1.99	0	0,3,3	0.00	-
2	ACT	A	304	-	1,3,3	3.51	1 (100%)	0,3,3	0.00	-

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
3	878	A	301	-	-	0/4/12/12	0/1/1/1
2	ACT	A	302	-	-	0/0/0/0	0/0/0/0
2	ACT	A	303	-	-	0/0/0/0	0/0/0/0
2	ACT	A	304	-	-	0/0/0/0	0/0/0/0

All (3) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	301	878	C3-N13	-2.89	1.36	1.41
2	A	302	ACT	CH3-C	3.22	1.53	1.48
2	A	304	ACT	CH3-C	3.51	1.53	1.48

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed( $^{\circ}$ )	Ideal( $^{\circ}$ )
3	A	301	878	C5-C4-C3	2.53	121.33	118.97

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 ⓘ

EDS was not executed - this section will therefore be empty.

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

EDS was not executed - this section will therefore be empty.

### 6.3 Carbohydrates ⓘ

EDS was not executed - this section will therefore be empty.

### 6.4 Ligands ⓘ

EDS was not executed - this section will therefore be empty.

### 6.5 Other polymers ⓘ

EDS was not executed - this section will therefore be empty.