



wwPDB X-ray Structure Validation Summary Report ⓘ

Oct 2, 2017 – 11:42 PM EDT

PDB ID : 1Q6T
Title : THE STRUCTURE OF PHOSPHOTYROSINE PHOSPHATASE 1B IN COMPLEX WITH COMPOUND 11
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Deposited on : unknown
Resolution : 2.30 Å(reported)

This is a wwPDB X-ray Structure Validation Summary Report for a publicly released PDB entry.

We welcome your comments at 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	:	1.7.2 (RC1), CSD as538be (2017)
Xtriage (Phenix)	:	1.9-1692
EDS	:	rb-20030345
Percentile statistics	:	20161228.v01 (using entries in the PDB archive December 28th 2016)
Refmac	:	5.8.0135
CCP4	:	6.5.0
Ideal geometry (proteins)	:	Engh & Huber (2001)
Ideal geometry (DNA, RNA)	:	Parkinson et al. (1996)
Validation Pipeline (wwPDB-VP)	:	rb-20030345

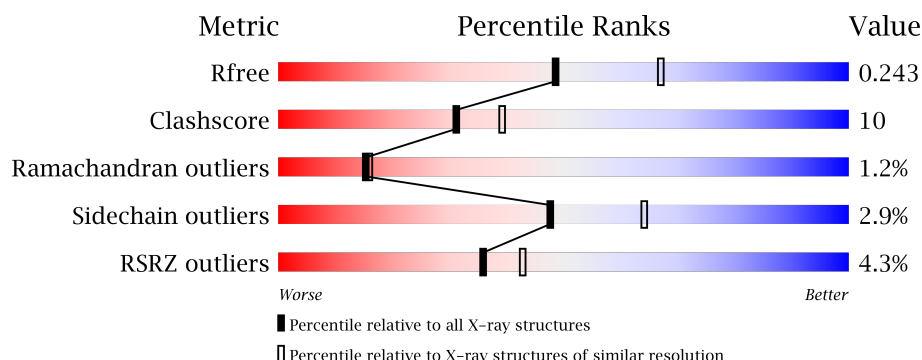
1 Overall quality at a glance

The following experimental techniques were used to determine the structure:

X-RAY DIFFRACTION

The reported resolution of this entry is 2.30 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}	100719	4130 (2.30-2.30)
Clashscore	112137	4751 (2.30-2.30)
Ramachandran outliers	110173	4705 (2.30-2.30)
Sidechain outliers	110143	4704 (2.30-2.30)
RSRZ outliers	101464	4156 (2.30-2.30)

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. 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	310	<div> <div>5%</div> <div> <div></div> <div>74%</div> <div>18%</div> <div>• 7%</div> </div> </div>
1	B	310	<div> <div>4%</div> <div> <div></div> <div>74%</div> <div>17%</div> <div>• 7%</div> </div> </div>

2 Entry composition

There are 4 unique types of molecules in this entry. The entry contains 5018 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 Protein-tyrosine phosphatase, non-receptor type 1.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	289	Total	C	N	O	S	0	0	0
			2349	1490	404	439	16			
1	B	289	Total	C	N	O	S	0	0	0
			2346	1488	404	439	15			

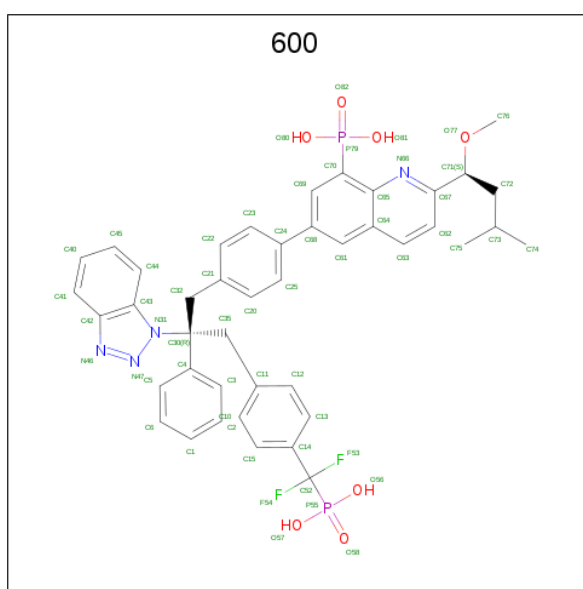
There are 24 discrepancies between the modelled and reference sequences:

Chain	Residue	Modelled	Actual	Comment	Reference
A	489	MET	-	CLONING ARTIFACT	UNP P18031
A	490	ASP	-	CLONING ARTIFACT	UNP P18031
A	491	TYR	-	CLONING ARTIFACT	UNP P18031
A	492	LYS	-	CLONING ARTIFACT	UNP P18031
A	493	ASP	-	CLONING ARTIFACT	UNP P18031
A	494	ASP	-	CLONING ARTIFACT	UNP P18031
A	495	ASP	-	CLONING ARTIFACT	UNP P18031
A	496	ASP	-	CLONING ARTIFACT	UNP P18031
A	497	LYS	-	CLONING ARTIFACT	UNP P18031
A	498	LEU	-	CLONING ARTIFACT	UNP P18031
A	499	GLU	-	CLONING ARTIFACT	UNP P18031
A	500	PHE	-	CLONING ARTIFACT	UNP P18031
B	989	MET	-	CLONING ARTIFACT	UNP P18031
B	990	ASP	-	CLONING ARTIFACT	UNP P18031
B	991	TYR	-	CLONING ARTIFACT	UNP P18031
B	992	LYS	-	CLONING ARTIFACT	UNP P18031
B	993	ASP	-	CLONING ARTIFACT	UNP P18031
B	994	ASP	-	CLONING ARTIFACT	UNP P18031
B	995	ASP	-	CLONING ARTIFACT	UNP P18031
B	996	ASP	-	CLONING ARTIFACT	UNP P18031
B	997	LYS	-	CLONING ARTIFACT	UNP P18031
B	998	LEU	-	CLONING ARTIFACT	UNP P18031
B	999	GLU	-	CLONING ARTIFACT	UNP P18031
B	1000	PHE	-	CLONING ARTIFACT	UNP P18031

- Molecule 2 is MAGNESIUM ION (three-letter code: MG) (formula: Mg).

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
2	B	1	Total	Mg	0	0
			1	1		
2	A	1	Total	Mg	0	0
			1	1		

- Molecule 3 is 6-[4-((2S)-2-(1H-1,2,3-BENZOTRIAZOL-1-YL)-3-{4-[DIFLUORO(PHOSPHONO)METHYL]PHENYL}-2-PHENYLPROPYL)PHENYL]-2-[(1S)-1-METHOXY-3-METHYLBUTYL]QUINOLIN-8-YLPHOSPHONIC ACID (three-letter code: 600) (formula: C₄₃H₄₂F₂N₄O₇P₂).

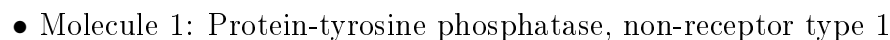


Mol	Chain	Residues	Atoms						ZeroOcc	AltConf
3	A	1	Total	C	F	N	O	P	0	0
			58	43	2	4	7	2		
3	B	1	Total	C	F	N	O	P	0	0
			58	43	2	4	7	2		

- Molecule 4 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
4	A	97	Total	O	0	0
			97	97		
4	B	108	Total	O	0	0
			108	108		

- Molecule 1: Protein-tyrosine phosphatase, non-receptor type 1



4 Data and refinement statistics

Property	Value	Source
Space group	P 21 21 21	Depositor
Cell constants a, b, c, α , β , γ	88.03Å 87.83Å 138.93Å 90.00° 90.00° 90.00°	Depositor
Resolution (Å)	12.00 – 2.30 28.38 – 1.96	Depositor EDS
% Data completeness (in resolution range)	97.4 (12.00-2.30) 91.2 (28.38-1.96)	Depositor EDS
R_{merge}	(Not available)	Depositor
R_{sym}	0.06	Depositor
$\langle I/\sigma(I) \rangle$ ¹	2.25 (at 1.96Å)	Xtriage
Refinement program	CNX	Depositor
R, R_{free}	0.214 , 0.246 0.213 , 0.243	Depositor DCC
R_{free} test set	2378 reflections (5.34%)	DCC
Wilson B-factor (Å ²)	34.9	Xtriage
Anisotropy	0.241	Xtriage
Bulk solvent k_{sol} (e/Å ³), B_{sol} (Å ²)	0.39 , 44.6	EDS
L-test for twinning ²	$\langle L \rangle = 0.43$, $\langle L^2 \rangle = 0.26$	Xtriage
Estimated twinning fraction	0.063 for k,h,-l	Xtriage
F_o, F_c correlation	0.94	EDS
Total number of atoms	5018	wwPDB-VP
Average B, all atoms (Å ²)	40.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 3.65% 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: MG, 600

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	0.49	0/2402	0.63	0/3237
1	B	0.53	1/2399 (0.0%)	0.70	1/3234 (0.0%)
All	All	0.51	1/4801 (0.0%)	0.67	1/6471 (0.0%)

All (1) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	B	1119	LEU	C-N	-5.74	1.20	1.34

All (1) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	B	1063	ASP	CA-CB-CG	-10.34	90.64	113.40

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	2349	0	2306	51	0
1	B	2346	0	2298	53	0
2	A	1	0	0	0	0
2	B	1	0	0	0	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
3	A	58	0	37	2	0
3	B	58	0	37	0	0
4	A	97	0	0	2	0
4	B	108	0	0	2	0
All	All	5018	0	4678	98	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 10.

The worst 5 of 98 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:1161:GLU:HG3	1:B:1168:THR:HG22	1.56	0.88
1:A:500:PHE:HA	1:A:503:MET:HE3	1.61	0.81
1:B:1002:GLU:HG3	1:B:1005:LYS:HE2	1.64	0.78
1:B:1112:ARG:NH1	1:B:1112:ARG:HB2	2.00	0.76
1:B:1011:ASP:O	1:B:1012:LYS:HG3	1.85	0.76

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	287/310 (93%)	270 (94%)	14 (5%)	3 (1%)	18	20
1	B	287/310 (93%)	270 (94%)	13 (4%)	4 (1%)	13	13
All	All	574/620 (93%)	540 (94%)	27 (5%)	7 (1%)	15	16

5 of 7 Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	B	1061	GLN
1	B	1116	LYS
1	A	619	LEU
1	B	1011	ASP
1	A	575	GLU

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	259/283 (92%)	252 (97%)	7 (3%)	50	67
1	B	258/283 (91%)	250 (97%)	8 (3%)	45	61
All	All	517/566 (91%)	502 (97%)	15 (3%)	48	64

5 of 15 residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	782	MET
1	B	1006	GLU
1	B	1260	LEU
1	A	765	ASP
1	B	1197	LYS

Some sidechains can be flipped to improve hydrogen bonding and reduce clashes. All (2) such sidechains are listed below:

Mol	Chain	Res	Type
1	B	1009	GLN
1	B	1139	ASN

5.3.3 RNA ⓘ

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 4 ligands modelled in this entry, 2 are monoatomic - leaving 2 for Mogul analysis.

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	600	A	801	-	56,64,64	3.63	38 (67%)	69,97,97	3.57	13 (18%)
3	600	B	1301	-	56,64,64	3.61	32 (57%)	69,97,97	3.57	16 (23%)

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	600	A	801	-	-	0/46/57/57	0/7/7/7
3	600	B	1301	-	-	0/46/57/57	0/7/7/7

The worst 5 of 70 bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
3	A	801	600	C67-C71	-15.55	1.42	1.51
3	B	1301	600	C67-C71	-15.40	1.42	1.51
3	B	1301	600	P55-O57	-4.45	1.46	1.54
3	A	801	600	O77-C76	-4.30	1.26	1.42
3	B	1301	600	P79-O81	-4.15	1.45	1.54

The worst 5 of 29 bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
3	B	1301	600	C72-C71-C67	-11.23	96.32	111.63
3	A	801	600	C72-C71-C67	-10.25	97.65	111.63
3	B	1301	600	C43-C42-N46	-4.21	103.29	108.59
3	A	801	600	C43-C42-N46	-3.88	103.71	108.59
3	A	801	600	C64-C65-N66	-3.67	117.91	122.69

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

1 monomer is involved in 2 short contacts:

Mol	Chain	Res	Type	Clashes	Symm-Clashes
3	A	801	600	2	0

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 [i](#)

6.1 Protein, DNA and RNA chains [i](#)

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	289/310 (93%)	0.08	14 (4%) 31 38	23, 38, 68, 76	0
1	B	289/310 (93%)	0.13	11 (3%) 41 48	21, 38, 70, 78	0
All	All	578/620 (93%)	0.11	25 (4%) 36 43	21, 39, 70, 78	0

The worst 5 of 25 RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	B	1285	SER	11.3
1	A	618	SER	5.9
1	A	619	LEU	5.5
1	B	1000	PHE	4.9
1	B	1117	GLY	4.8

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. LLDF column lists the quality of electron density of the group with respect to its neighbouring residues in protein, DNA or RNA chains. 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	LLDF	B-factors(\AA^2)	Q<0.9
3	600	B	1301	58/58	0.96	0.14	1.04	19,28,44,47	0
3	600	A	801	58/58	0.96	0.16	0.83	22,28,55,59	0
2	MG	B	5000	1/1	0.86	0.11	-	26,26,26,26	1
2	MG	A	6000	1/1	0.97	0.14	-	32,32,32,32	1

6.5 Other polymers [i](#)

There are no such residues in this entry.