



# Full wwPDB X-ray Structure Validation Report ⓘ

Feb 27, 2014 – 12:58 PM GMT

PDB ID : 1V0R  
Title : TUNGSTATE-INHIBITED PHOSPHOLIPASE D FROM STREPTOMYCES  
SP. STRAIN PMF  
Authors : Leiros, I.; Mcsweeney, S.; Hough, E.  
Deposited on : 2004-04-01  
Resolution : 1.70 Å(reported)

This is a full wwPDB 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/ValidationPDFNotes.html>

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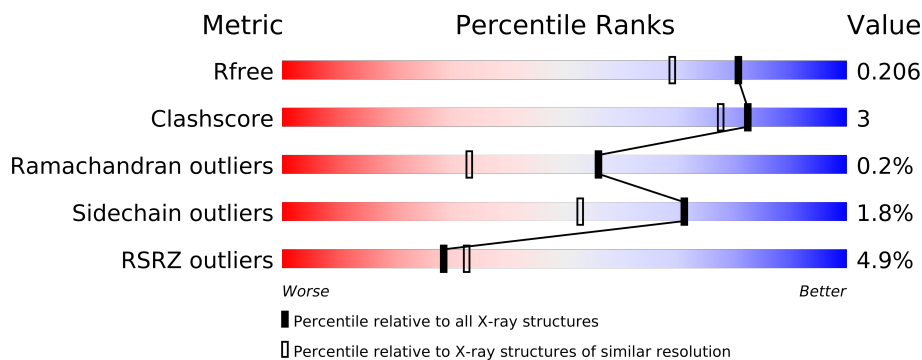
The following versions of software and data (see [references](#)) were used in the production of this report:

MolProbity : 4.02b-467  
Mogul : 1.15 2013  
Xtriage (Phenix) : dev-1323  
EDS : stable22639  
Percentile statistics : 21963  
Refmac : 5.8.0049  
CCP4 : 6.3.0 (Settle)  
Ideal geometry (proteins) : Engh & Huber (2001)  
Ideal geometry (DNA, RNA) : Parkinson et. al. (1996)  
Validation Pipeline (wwPDB-VP) : stable22683

# 1 Overall quality at a glance

The reported resolution of this entry is 1.70 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))
$R_{free}$	66092	2456 (1.70-1.70)
Clashscore	79885	2929 (1.70-1.70)
Ramachandran outliers	78287	2878 (1.70-1.70)
Sidechain outliers	78261	2878 (1.70-1.70)
RSRZ outliers	66119	2456 (1.70-1.70)

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. The upper red bar (where present) indicates the fraction of residues that have poor fit to the electron density.

Mol	Chain	Length	Quality of chain
1	A	506	

## 2 Entry composition i

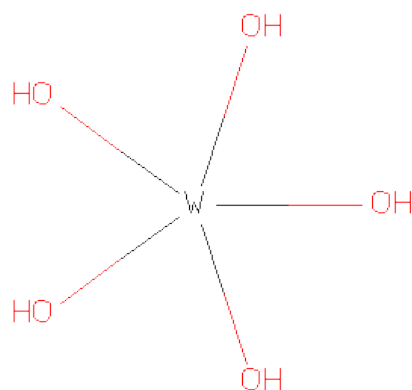
There are 3 unique types of molecules in this entry. The entry contains 4250 atoms, of which 0 are hydrogen and 0 are deuterium.

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 PHOSPHOLIPASE D.

Mol	Chain	Residues	Atoms					ZeroOcc	AltConf	Trace
1	A	493	Total	C	N	O	S	52	0	0
			3700	2308	651	727	14			

- Molecule 2 is TUNGSTATE(VI) ION (three-letter code: WO5) (formula:  $\text{H}_5\text{O}_5\text{W}$ ).



Mol	Chain	Residues	Atoms			ZeroOcc	AltConf
2	A	1	Total	O	W	0	0
			6	5	1		

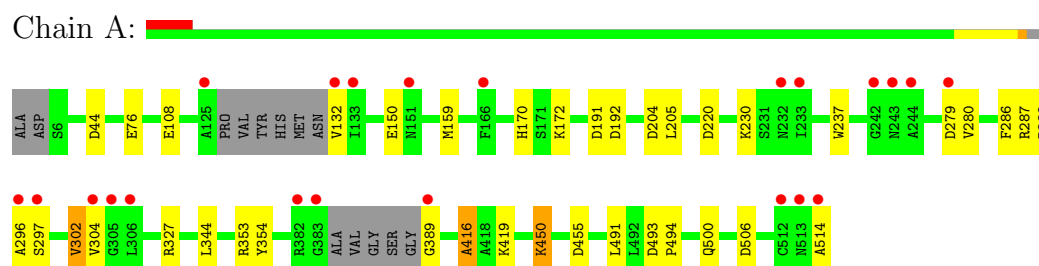
- Molecule 3 is water.

Mol	Chain	Residues	Atoms		ZeroOcc	AltConf
3	A	544	Total	O	0	0
			544	544		

### 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 ( $RSRZ > 2$ ). 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: PHOSPHOLIPASE D



## 4 Data and refinement statistics

Property	Value	Source
Space group	P 1 21 1	Depositor
Cell constants a, b, c, $\alpha$ , $\beta$ , $\gamma$	57.61Å 56.90Å 68.95Å 90.00° 93.67° 90.00°	Depositor
Resolution (Å)	69.01 – 1.70 22.94 – 1.70	Depositor EDS
% Data completeness (in resolution range)	99.6 (69.01-1.70) 99.7 (22.94-1.70)	Depositor EDS
$R_{merge}$	0.04	Depositor
$R_{sym}$	(Not available)	Depositor
$\langle I/\sigma(I) \rangle$ <sup>1</sup>	2.89 (at 1.70Å)	Xtriage
Refinement program	REFMAC 5.1.24	Depositor
R, $R_{free}$	0.155 , 0.198 0.169 , 0.206	Depositor DCC
$R_{free}$ test set	2485 reflections (5.35%)	DCC
Wilson B-factor (Å <sup>2</sup> )	18.4	Xtriage
Anisotropy	0.534	Xtriage
Bulk solvent $k_{sol}$ (e/Å <sup>3</sup> ), $B_{sol}$ (Å <sup>2</sup> )	0.35 , 45.0	EDS
Estimated twinning fraction	No twinning to report.	Xtriage
L-test for twinning	$\langle  L  \rangle = 0.49$ , $\langle L^2 \rangle = 0.33$	Xtriage
Outliers	0 of 48957 reflections	Xtriage
$F_o, F_c$ correlation	0.96	EDS
Total number of atoms	4250	wwPDB-VP
Average B, all atoms (Å <sup>2</sup> )	23.0	wwPDB-VP

Xtriage's analysis on translational NCS is as follows: *The largest off-origin peak in the Patterson function is 6.67% of the height of the origin peak. No significant pseudotranslation is detected.*

<sup>1</sup>Intensities estimated from amplitudes.

## 5 Model quality

### 5.1 Standard geometry

Bond lengths and bond angles in the following residue types are not validated in this section: WO5

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	1.16	8/3775 (0.2%)	1.09	22/5141 (0.4%)

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	1	1

All (8) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
1	A	290	LEU	C-N	32.05	1.95	1.34
1	A	150	GLU	CB-CG	-22.19	1.09	1.52
1	A	295	THR	CA-CB	-21.15	0.98	1.53
1	A	302	VAL	C-N	17.00	1.73	1.34
1	A	416	ALA	C-N	8.46	1.53	1.34
1	A	132	VAL	CA-CB	-6.84	1.40	1.54
1	A	170	HIS	CD2-NE2	-6.59	1.23	1.38
1	A	170	HIS	CG-CD2	-6.34	1.25	1.35

All (22) bond angle outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	295	THR	N-CA-CB	29.42	166.20	110.30
1	A	295	THR	CA-CB-CG2	14.88	133.24	112.40
1	A	295	THR	CB-CA-C	-11.27	81.17	111.60
1	A	150	GLU	CA-CB-CG	10.06	135.53	113.40
1	A	170	HIS	CG-CD2-NE2	9.36	126.98	109.20

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Mol	Chain	Res	Type	Atoms	Z	Observed(°)	Ideal(°)
1	A	327	ARG	NE-CZ-NH1	7.83	124.21	120.30
1	A	220	ASP	CB-CG-OD1	7.49	125.04	118.30
1	A	290	LEU	CA-C-N	-7.20	96.93	117.10
1	A	290	LEU	O-C-N	-6.95	107.90	121.10
1	A	302	VAL	O-C-N	-6.59	112.16	122.70
1	A	237	TRP	CD1-CG-CD2	6.52	111.51	106.30
1	A	192	ASP	CB-CG-OD2	6.48	124.14	118.30
1	A	204	ASP	CB-CG-OD1	6.24	123.92	118.30
1	A	191	ASP	CB-CG-OD2	6.16	123.84	118.30
1	A	506	ASP	CB-CG-OD2	6.13	123.82	118.30
1	A	44	ASP	CB-CG-OD1	5.89	123.60	118.30
1	A	455	ASP	CB-CG-OD2	5.83	123.55	118.30
1	A	170	HIS	ND1-CG-CD2	-5.77	97.92	106.00
1	A	289	ASP	CB-CG-OD2	5.72	123.45	118.30
1	A	327	ARG	NE-CZ-NH2	-5.39	117.60	120.30
1	A	353	ARG	NE-CZ-NH2	-5.22	117.69	120.30
1	A	132	VAL	N-CA-CB	5.21	122.95	111.50

All (1) chirality outliers are listed below:

Mol	Chain	Res	Type	Atom
1	A	295	THR	CA

All (1) planarity outliers are listed below:

Mol	Chain	Res	Type	Group
1	A	416	ALA	Mainchain

## 5.2 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 hydrogens added by MolProbity. The Clashes column lists the number of clashes within the asymmetric unit, and the number in parentheses is this value normalized per 1000 atoms of the molecule in the chain. The Symm-Clashes column gives symmetry related clashes, in the same way as for the Clashes column.

Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
1	A	3700	0	3630	21	0
2	A	6	0	0	3	0
3	A	544	0	0	7	0

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Mol	Chain	Non-H	H(model)	H(added)	Clashes	Symm-Clashes
All	All	4250	0	3630	22	0

Clashscore is defined as the number of clashes calculated for the entry per 1000 atoms (including hydrogens) of the entry. The overall clashscore for this entry is 3.

All (22) close contacts within the same asymmetric unit are listed below.

Atom-1	Atom-2	Distance(Å)	Clash(Å)
1:A:302:VAL:C	1:A:304:VAL:N	1.73	1.41
1:A:290:LEU:C	1:A:294:PRO:N	1.95	1.19
1:A:290:LEU:C	1:A:294:PRO:CA	2.65	0.64
1:A:290:LEU:CA	1:A:294:PRO:N	2.61	0.64
1:A:76:GLU:CG	3:A:2156:HOH:O	2.47	0.62
1:A:108:GLU:HG2	3:A:2211:HOH:O	2.02	0.57
1:A:389:GLY:N	3:A:2435:HOH:O	2.39	0.56
1:A:297:SER:HB2	1:A:354:TYR:OH	2.05	0.55
1:A:279:ASP:OD2	1:A:280:VAL:HG12	2.08	0.53
1:A:294:PRO:HG3	3:A:2419:HOH:O	2.08	0.52
1:A:450:LYS:NZ	2:A:600:WO5:O2	2.43	0.52
1:A:230:LYS:HE3	3:A:2322:HOH:O	2.12	0.49
1:A:302:VAL:C	1:A:304:VAL:CA	2.75	0.49
1:A:491:LEU:HD23	1:A:491:LEU:C	2.34	0.48
1:A:290:LEU:HA	1:A:294:PRO:N	2.30	0.46
1:A:286:PHE:CZ	1:A:288:PRO:HB3	2.51	0.45
1:A:286:PHE:O	1:A:287:ARG:HD3	2.18	0.44
1:A:419:LYS:NZ	1:A:514:ALA:O	2.50	0.43
1:A:500:GLN:NE2	3:A:2534:HOH:O	2.52	0.42
1:A:450:LYS:CE	2:A:600:WO5:O2	2.68	0.42
1:A:493:ASP:HB2	1:A:494:PRO:HD3	2.01	0.41
2:A:600:WO5:O3	3:A:2544:HOH:O	2.22	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	487/506 (96%)	473 (97%)	13 (3%)	1 (0%)	56 33

All (1) Ramachandran outliers are listed below:

Mol	Chain	Res	Type
1	A	296	ALA

### 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	395/406 (97%)	388 (98%)	7 (2%)	71 53

All (7) residues with a non-rotameric sidechain are listed below:

Mol	Chain	Res	Type
1	A	159	MET
1	A	172	LYS
1	A	205	LEU
1	A	289	ASP
1	A	295	THR
1	A	344	LEU
1	A	450	LYS

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

Mol	Chain	Res	Type
1	A	20	GLN
1	A	309	ASN
1	A	409	ASN

### 5.3.3 RNA ⓘ

There are no RNA chains 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 ⓘ

1 ligand is 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	$\# Z  > 2$
2	WO5	A	600	1	5,5,5	4.38	5 (100%)	0,10,10	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
2	WO5	A	600	1	-	0/0/0/0	0/0/0/0

All (5) bond length outliers are listed below:

Mol	Chain	Res	Type	Atoms	Z	Observed(Å)	Ideal(Å)
2	A	600	WO5	W-O1	5.34	2.12	1.78
2	A	600	WO5	W-O4	4.99	2.10	1.78
2	A	600	WO5	W-O2	4.71	2.08	1.78
2	A	600	WO5	W-O3	3.81	2.02	1.78
2	A	600	WO5	W-O5	2.44	1.94	1.78

There are no bond angle outliers.

There are no chirality outliers.

There are no torsion outliers.

There are no ring outliers.

## 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 ⓘ

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, 95<sup>th</sup> 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 <sup>2</sup> )	Q<0.9
1	A	493/506 (97%)	-0.05	24 (4%)	28 33	12, 20, 36, 50	14 (2%)

All (24) RSRZ outliers are listed below:

Mol	Chain	Res	Type	RSRZ
1	A	296	ALA	7.9
1	A	383	GLY	7.6
1	A	132	VAL	7.6
1	A	125	ALA	4.9
1	A	242	GLY	4.1
1	A	382	ARG	3.9
1	A	233	ILE	3.8
1	A	297	SER	3.7
1	A	389	GLY	3.6
1	A	306	LEU	3.3
1	A	279	ASP	3.1
1	A	133	ILE	2.9
1	A	304	VAL	2.9
1	A	294	PRO	2.8
1	A	514	ALA	2.8
1	A	305	GLY	2.6
1	A	243	ASN	2.6
1	A	232	ASN	2.5
1	A	244	ALA	2.5
1	A	166	PHE	2.5
1	A	151	ASN	2.3
1	A	513	ASN	2.3
1	A	295	THR	2.3
1	A	512	CYS	2.1

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

There are no non-standard protein/DNA/RNA residues in this entry.

## 6.3 Carbohydrates

There are no carbohydrates in this entry.

## 6.4 Ligands

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, 95<sup>th</sup> 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	RSR	LLDF	B-factors(Å <sup>2</sup> )	Q<0.9
2	WO5	A	600	6/6	0.06	-0.48	17,22,24,26	0

## 6.5 Other polymers

There are no such residues in this entry.