Pergamon

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Corrigendum

Corrigendum to "Novel Diphosphetene Derivatives by Reactions of Di(isopropyl)aminophosphaethyne with Chalcogens or Halogens" [Tetrahedron 56 (2000) 27]☆

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On page 32, the data for **6b** of Table 2 is incorrect, the correct data should appear as:

Table 2. Crystallographic data and parameters of the crystal structure determinations

Compound	**6a**	**6b**
Empirical formula	$C_{14}H_{28}N_2P_2S_2$	$C_{14}H_{28}N_2P_2Se_2$
Fw	350.44	444.24
Crystal size (mm)	0.23×0.20×0.26	0.20×0.22×0.25
Crystal system	Monoclinic	Monoclinic
Space group	$P2_1/n$	$P2_1/n$
a (Å)	10.635(2)	10.526(3)
b (Å)	13.081(2)	13.606(3)
c (Å)	14.919(3)	14.947(3)
β (°)	108.53	107.65
V (Å^3)	1967.9(6)	2039.9(8)
Z	4	4
ρ_{calcd} (g cm^{-3})	1.183	1.447
μ (mm^{-1})	0.427	3.776
F(000)	752	896
Temperature (K)	173(2)	153(2)
θ (degrees)	2.07–27.00	2.07–24.00
Index ranges	$0\leq h\leq 13$	$-7\leq h\leq 12$
	$0\leq k\leq 16$	$-6\leq k\leq 15$
	$-19\leq l\leq 18$	$-17\leq l\leq 16$
No. of reflns measd	4528	3386
No. of indep rflns with $I>2\sigma(I)$	3448	1829
No. of parameters	293	189
$R1$ (obs. data)	0.0319	0.0402
$wR2$ (obs. data)	0.0841	0.0728
$R1$ (all data)	0.0410	0.0848
$wR2$ (all data)	0.0866	0.0794
GooF on F^2	1.024	0.804
Resid. electron density (eÅ^{-3})	+0.353/−0.337	+0.730/−0.345

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