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## Phosphorus, Sulfur, and Silicon and the Related Elements

Publication details, including instructions for authors and subscription information:

<http://www.informaworld.com/smpp/title~content=t713618290>

## Erratum

**To cite this Article** (2006) 'Erratum', *Phosphorus, Sulfur, and Silicon and the Related Elements*, 181: 11, 2679

**To link to this Article:** DOI: 10.1080/10426500600935674

**URL:** <http://dx.doi.org/10.1080/10426500600935674>

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## I Erratum

In issue 181(6), the article ‘Synthesis and Reactions of Anhydro-Azido-thio-D-lyxofuranosides’ by D. Otzen, J. Voss, and G. Adiwidjaja on pages 1249–1270, the three compounds 26, 27, and 29 were incorrectly printed in Table 1. Following is the corrected Table 1.

**TABLE I Crystal Data and Structure Refinement for 27, 29, and 26**

Compound	27	29	26
Empirical formula	C <sub>6</sub> H <sub>9</sub> N <sub>3</sub> O <sub>3</sub> S	C <sub>6</sub> H <sub>9</sub> N <sub>3</sub> O <sub>4</sub> S	C <sub>6</sub> H <sub>9</sub> N <sub>3</sub> O <sub>4</sub> S
Formula weight	203.22	219.22	219.22
Crystal system	Orthorhombic	Monoclinic	Orthorhombic
Space group	<i>P</i> 2 <sub>1</sub> 2 <sub>1</sub> 2 <sub>1</sub>	<i>P</i> 2 <sub>1</sub>	<i>P</i> 2 <sub>1</sub> 2 <sub>1</sub> 2 <sub>1</sub>
<i>a</i> [Å]	5.494(1)	6.706(1)	8.760(1)
<i>b</i> [Å]	10.944(1)	8.522	10.209(1)
<i>c</i> [Å]	14.265(1)	8.385(1)	10.345(1)
$\beta$ [°]		110.69(1)	
<i>Z</i>	4	2	4
Crystal size [mm]	0.33 × 0.25 × 0.16	0.37 × 0.27 × 0.23	0.33 × 0.28 × 0.15
$\rho_{\text{calcd.}}$ [g cm <sup>-3</sup> ]	1.574	1.624	1.574
$\mu$ [mm <sup>-1</sup> ]	0.356	0.355	0.344
<i>F</i> (000)	424	228	456
$\theta$ limits [°]	2.35/27.48	2.60/30.02	2.80/30.02
<i>h</i> / <i>k</i> / <i>l</i> limits	0,7/0,14/0,18	0,9/0,11/–11,11	0,12/0,14/0,14
Reflections collected	9348	3565	9913
Unique reflections	1169 [ <i>R</i> (int) = 0.047]	1376 [ <i>R</i> (int) = 0.026]	1565 [ <i>R</i> (int) = 0.033]
Numb. of parameters	143	152	152
<i>R</i> indices (all data)	<i>R</i> <sub>1</sub> = 0.0318 <i>WR</i> <sub>2</sub> = 0.0770	<i>R</i> <sub>1</sub> = 0.0497 <i>wR</i> <sub>2</sub> = 0.1100	<i>R</i> <sub>1</sub> = 0.0457 <i>wR</i> <sub>2</sub> = 0.0837
Final <i>R</i> indices [ <i>I</i> > 2 $\sigma$ ( <i>I</i> )]	<i>R</i> <sub>1</sub> = 0.0268 <i>WR</i> <sub>2</sub> = 0.0739	<i>R</i> <sub>1</sub> = 0.0432 <i>wR</i> <sub>2</sub> = 0.1051	<i>R</i> <sub>1</sub> = 0.0324 <i>wR</i> <sub>2</sub> = 0.0769
Goodness of fit on <i>F</i> <sup>2</sup>	1.036	1.016	1.003
Abs. struct. parameter	0.39(10)	–0.06(13)	0.20(10)
Extinction coefficient	0.015(5)	0.43(3)	0.037(4)
Largest difference peak and hole [e Å <sup>-3</sup> ]	0.156 and –0.198	0.582 and –0.631	0.166 and –0.211