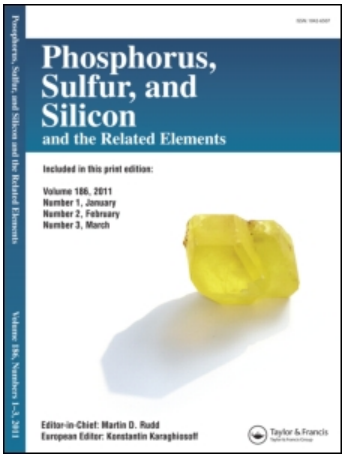

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Erratum

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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	$C_6H_9N_3O_3S$	$C_6H_9N_3O_4S$	$C_6H_9N_3O_4S$
Formula weight	203.22	219.22	219.22
Crystal system	Orthorhombic	Monoclinic	Orthorhombic
Space group	$P2_12_12_1$	$P2_1$	$P2_12_12_1$
a [Å]	5.494(1)	6.706(1)	8.760(1)
b [Å]	10.944(1)	8.522	10.209(1)
c [Å]	14.265(1)	8.385(1)	10.345(1)
β [°]		110.69(1)	
Z	4	2	4
Crystal size [mm]	0.33 × 0.25 × 0.16	0.37 × 0.27 × 0.23	0.33 × 0.28 × 0.15
$\rho_{calcd.}$ [g cm^{-1}]	1.574	1.624	1.574
μ [mm^{-1}]	0.356	0.355	0.344
$F(000)$	424	228	456
θ limits [°]	2.35/27.48	2.60/30.02	2.80/30.02
$h/k/l$ 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 [R(int) = 0.047]	1376 [R(int) = 0.026]	1565 [R(int) = 0.033]
Numb. of parameters	143	152	152
R indices (all data)	$R_1 = 0.0318$ $WR_2 = 0.0770$	$R_1 = 0.0497$ $wR_2 = 0.1100$	$R_1 = 0.0457$ $wR_2 = 0.0837$
Final R indices [$I > 2\,\sigma(I)$]	$R_1 = 0.0268$ $WR_2 = 0.0739$	$R_1 = 0.0432$ $wR_2 = 0.1051$	$R_1 = 0.0324$ $wR_2 = 0.0769$
Goodness of fit on F^2	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 A^{-3}]	0.156 and −0.198	0.582 and −0.631	0.166 and −0.211