154 ERRATA

## **Errata**

Fen-Er Chen, Xiou-Hong Ling, Yan-Ping He, Xiao-Hua Peng *Synthesis* **2001**, 1773.

In Table 2, wrong structures of substrates **1b-k** were published, the corrected Table is reprinted below. We would like to apologize for this error.

**Table 2**  $(n\text{-Bu}_4\text{N})_2\text{S}_2\text{O}_8$ -Mediated Selective Deprotection of Allyl Ethers  $(\mathbf{1a}-\mathbf{k})^a$ 

Ethers (1a-k) <sup>a</sup>				
Entry	Substrate	Time (h)	Product <sup>b</sup>	Yield of 3 (%) <sup>c</sup>
1	O O O O O O O O O O O O O O O O O O O	6	3a	84
2	AllO	8	3b	81
3	AllO OCH <sub>3</sub>	6	3c	84
4	OAII	5	3d	85
5	THPO OAll	5	3e	83
6	Allo	6	3f	82
7	OAII	6	<b>3</b> g	81
8	1g OAll	7	3h	83
9	OTr OMe	6	3I	82
10	O O O O O O O O O O O O O O O O O O O	8	<b>3</b> j	81
11	O O O O O O O O O O O O O O O O O O O	6	3k	85

<sup>&</sup>lt;sup>a</sup> All reactions were performed according to the typical procedure.

Giorgio Abbiati, Egle M. Beccalli, Alessandro Marchesini, Elisabetta Rossi *Synthesis* **2001**, 2477.

The structures for **2e** (Scheme 1) and its precursor were depicted wrongly. They are represented below in the correct way. The authors apologize for this mistake. :

The reagents for the reactions of **2a-d** with acetylenes (Scheme 2) are, Pd<sup>0</sup>, CuI, TEA, DMF, r.t.

<sup>&</sup>lt;sup>b</sup> All products were characterized by comparison of their melting points and <sup>1</sup>H NMR spectra with those of authentic samples.

<sup>&</sup>lt;sup>c</sup> Yield of isolated pure products.