

In the paper by J. Barluenga et al. published in *Chem. Eur. J.* **2001**, 7,3533, there are errors. As pointed out in the text, the absolute configuration of carbene complexes **7** was assigned to be R for C-9 on the basis of the X-ray structure of dithiolane derivative **8g**. Accordingly, Scheme 3 should be depicted as follows for clarity. Additionally, in the Experimental Section the absolute configuration should be added and/or corrected for several compounds as follows: **7a-c** ( $8S^*$ ,  $9R^*$ ), **7d** ( $8R^*$ ,  $9R^*$ ), **7j** (8R, 9R), **8a,b** ( $8S^*$ ,  $9S^*$ ), **8c** ( $8R^*$ ,  $9S^*$ ), **8d,f,g,h,j** (8S, 9S), **8e,i** (8R, 9S), **9b** ( $3S^*$ ,  $4S^*$ ), **9c** ( $3S^*$ ,  $4R^*$ ), **9e** (3S, 4R), **9f,g,j** (3S, 4S). The authors apologize for these mistakes.

Scheme 3. Structural assignments: NOESY crosspeak for compound 9e and absolute configuration for compounds 7 determined by X-ray analysis of 8g.

In the paper by J. J. Schneider et al. published in *Chem. Eur. J.* **2001**, 7, 2888, there is a mistake in Figure 11. The correct figure is given below. The authors apologize for this error.

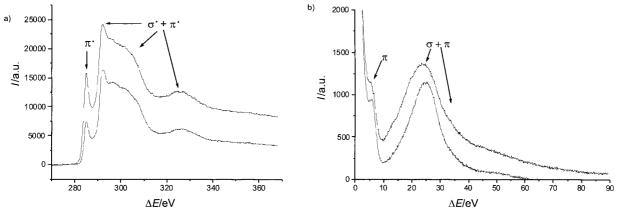


Figure 11. EEL spectra: a) core-loss region, and b)  $C_{1s}$  low-loss region of CNT fibers compared to HOPG. Upper trace: CNT fibers; lower trace: HOPG. I = intensity,  $\Delta E = \text{energy loss}$ .