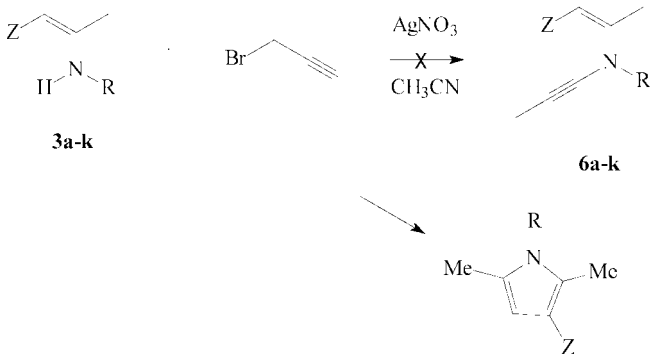
Errata

D. Gravestock, M. C. Dovey, *Synthesis*, **2003**, 523.

We have been interested in unusual chemical shifts displayed by compounds **6a**–**k**. Attempts to rationalise these observations by performing chemical transformations and *ab initio* quantum chemical calculations (Gaussian 98) met with little success. It was subsequently felt that a crystal structure might reveal valuable information as to the electronic character of the compounds in question. Following further work, a crystal suitable for X-ray diffraction studies was obtained and this revealed that the original structure proposed for **6a**–**k** was, in fact, incorrect. The correct structure is shown below (Scheme 1) and the authors apologise for this error.



Scheme 1