Tetrahedron Vol. 40, No. 2, p. 267, 1984. Pergamon Press Ltd. Printed in Great Britain.

ERRATA

PETER J. KRUEGER and ADRIAN O. FULEA: Rotation about the C–N bond in thioamides. Influence of substituents on the potential function. *Tetrahedron* **31,** 1813 (1975).

The Figure 4 is wrong because it is in contradiction to the results of the quantum mechanical treatment which leads to quantized energy and therefore should be disregarded. The conclusion that in case of compounds like **3** and **4** only an average $\overline{\Delta G^{\neq}}$ can be obtained by NMR spectroscopy is valid. We thank Dr. A. Rank for the criticism of Fig. 4.

G. PAGANI ZECCHINI, M. PAGLIALUNGA PARADISI and I. TORRINI, *In situ* selective protection of aldehydes *via* aldimines. Simple conversions of ketoaldehydes to methylene aldehydes and to methylhydroxy aldehydes. *Tetrahedron* **39,** 2709 (1983).

Page 2710, column 2, lines 1 and 7: "**2a**" should read "**1b**".