Additions and Corrections

Transition-State Modeling in Acyclic Stereoselection. A Molecular Mechanics Model for the Kinetic Formation of Lithium Enolates [*J. Am. Chem. Soc.* **1985**, *107*, 2264–2273]. DAVID W. MORELAND* and WILLIAM G. DAUBEN*

The graphical presentations of Figures 3 and 4 were transposed. They should appear as follows:

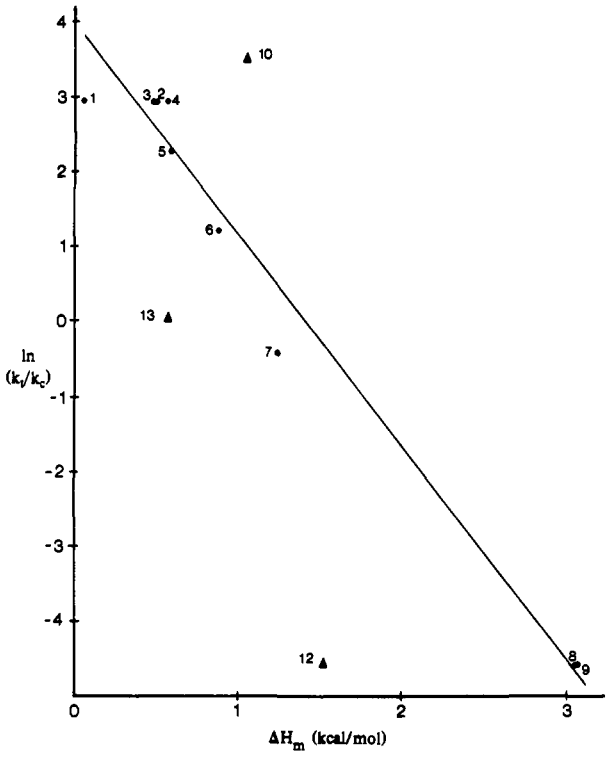

Figure 3.

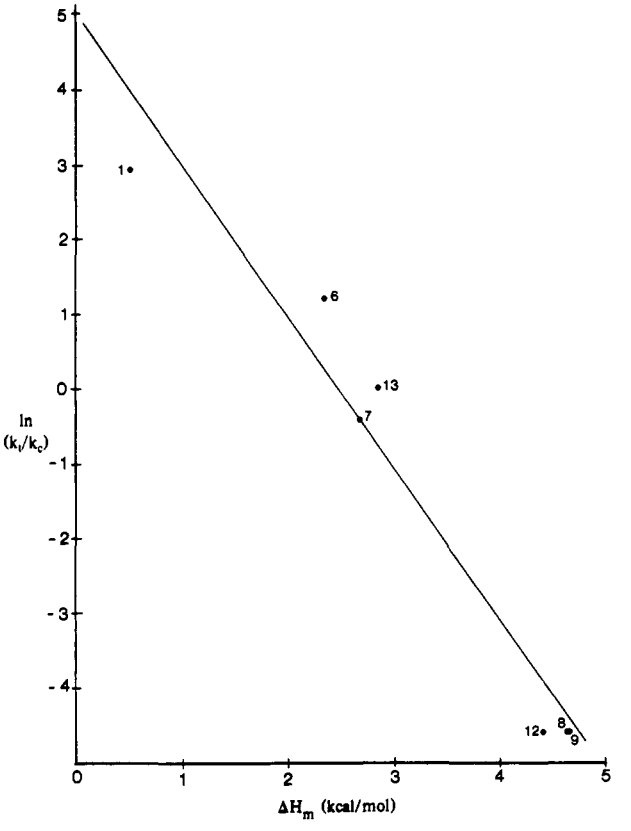

Figure 4.