## ADDITIONS AND CORRECTIONS

## 1996, Volume 100

W. L. Hase,\* H. B. Schlegel, V. Balbyshev, and M. Page: Ab Initio Study of the Transition State and Forward and Reverse Rate Constants for  $C_2H_5 \rightleftharpoons H + C_2H_4$ 

Page 5358. In the second column, 28 lines from the bottom the sentence "The resulting vibrationally adiabatic quantum barrier  $E_{0,r}$  is 3.05 kcal/mol." should read as follows: "The resulting vibrationally adiabatic quantum barrier  $E_{0,r}$  is 2.70 kcal/mol." The incorrect value resulted from an error in applying the tunneling correction while adjusting the barrier to fit experiment.

## 1997, Volume 101A

Maija Lahtela,\* Tapani A. Pakkanen, and Richard L. Rowley: Nonequilibrium Molecular Dynamics Simulations of 3-Methylhexane: The Effect of Inter- and Intramolecular Potential Models on Simulated Viscosity

Page 3449. In the paragraph simulation details the shear rate range should be  $3.5 \times 10^{10} \ s^{-1} \le \gamma \le 14.1 \times 10^{10} \ s^{-1}$ .

In Tables 3, 4, and 5 the shear rates  $\gamma$  should be G s<sup>-1</sup>.

In Figures 3, 4, and 5 the title of category (×) axis should be  $\gamma^{1/2}\ 10^{10}\ s^{-1}.$