

Time-resolved gas-phase kinetic, quantum chemical and RRKM studies of the reaction of silylene with 2,5-dihydrofuran

Article

Accepted Version

Figure 6

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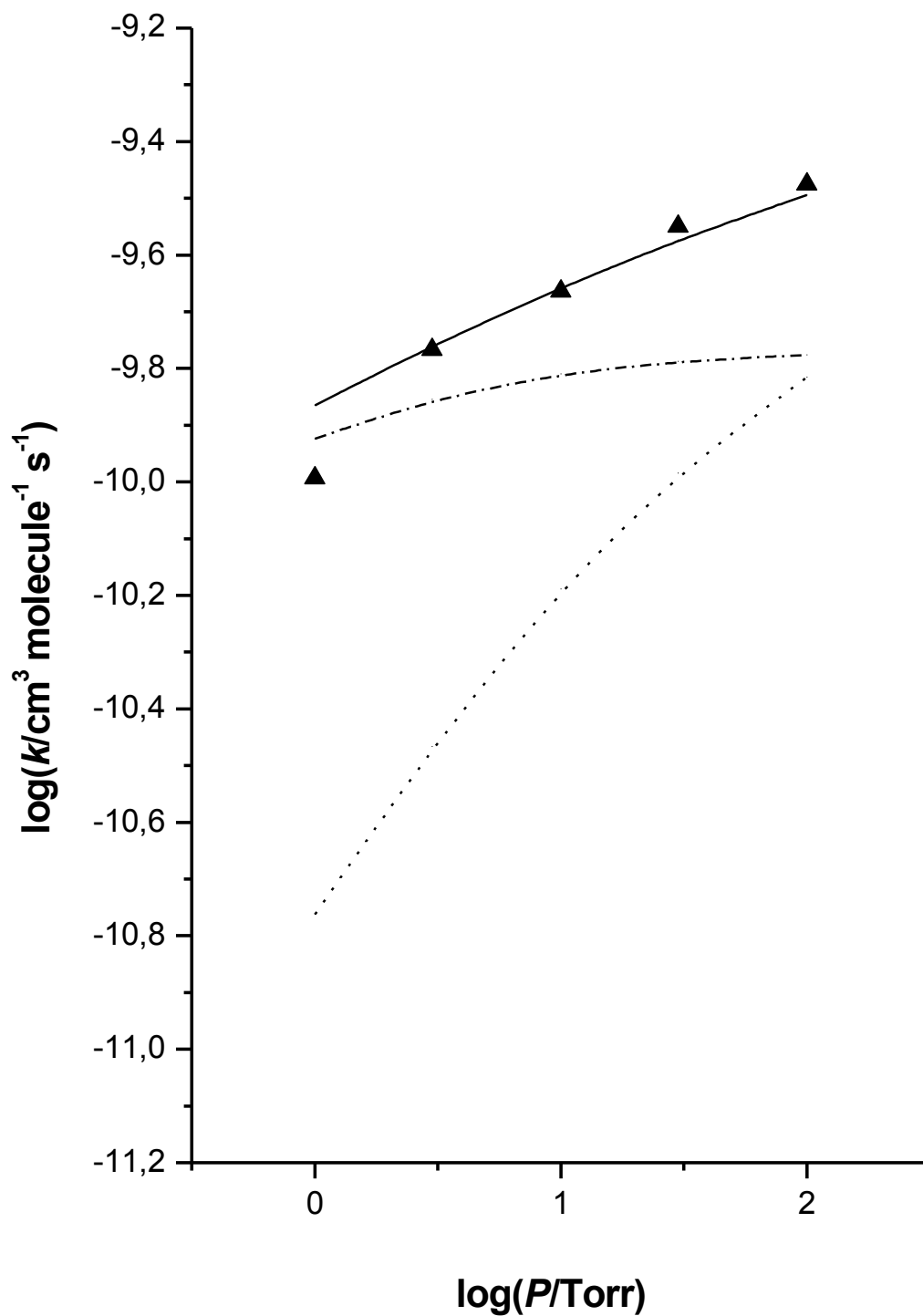


Figure 6. RRKM fit (solid line) to the data at 296 K. Individual lines show contributions from the π -pathway (·- ·) and O-donor pathway (----).

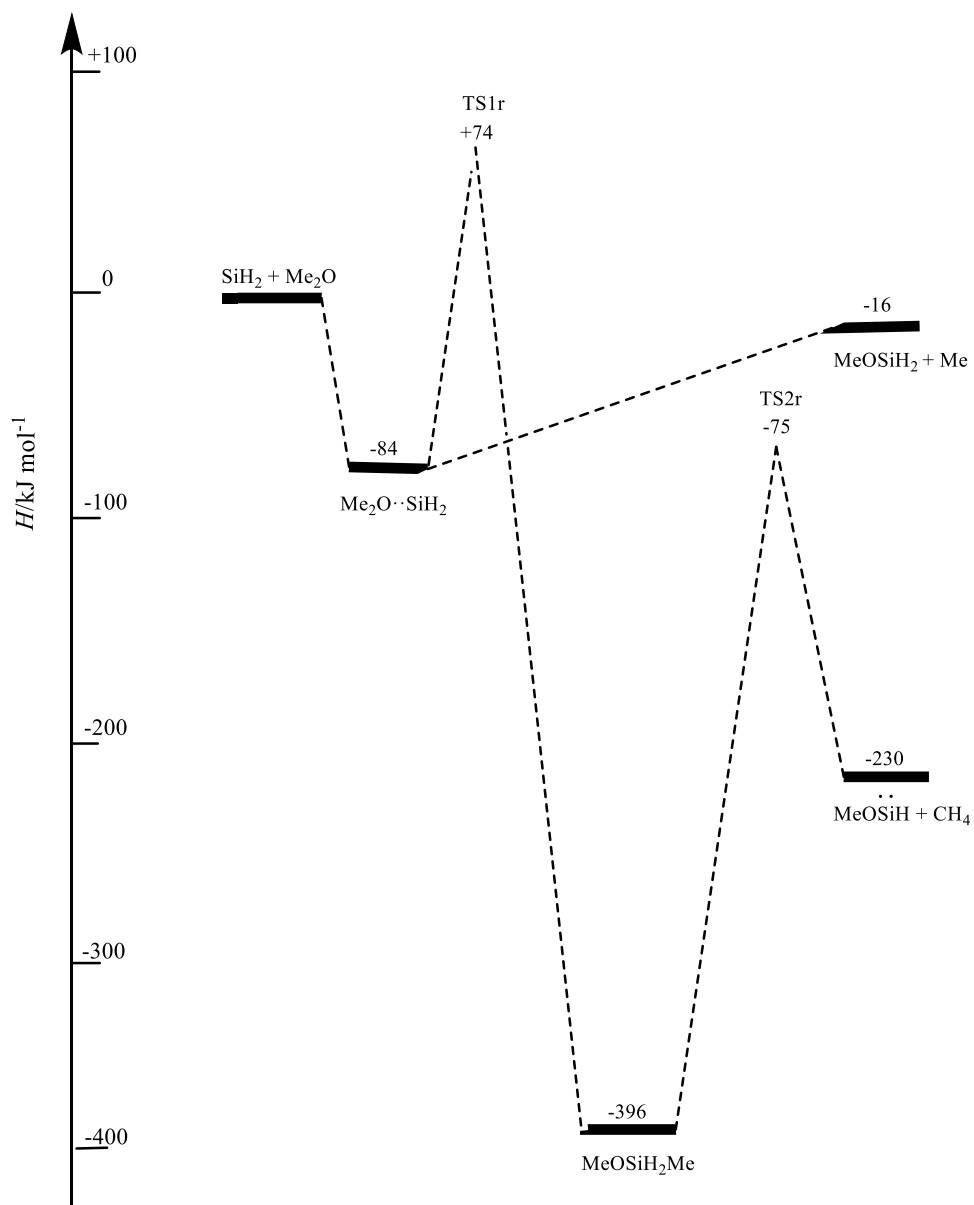


Figure 7. Potential energy (enthalpy) surface for $\text{C}_2\text{H}_8\text{SiO}$ ($\text{SiH}_2 + \text{Me}_2\text{O}$). All enthalpies are calculated at the G3 level and quoted at 298 K.

Graphic for abstract

