

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

Article

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Figures 4 & 5

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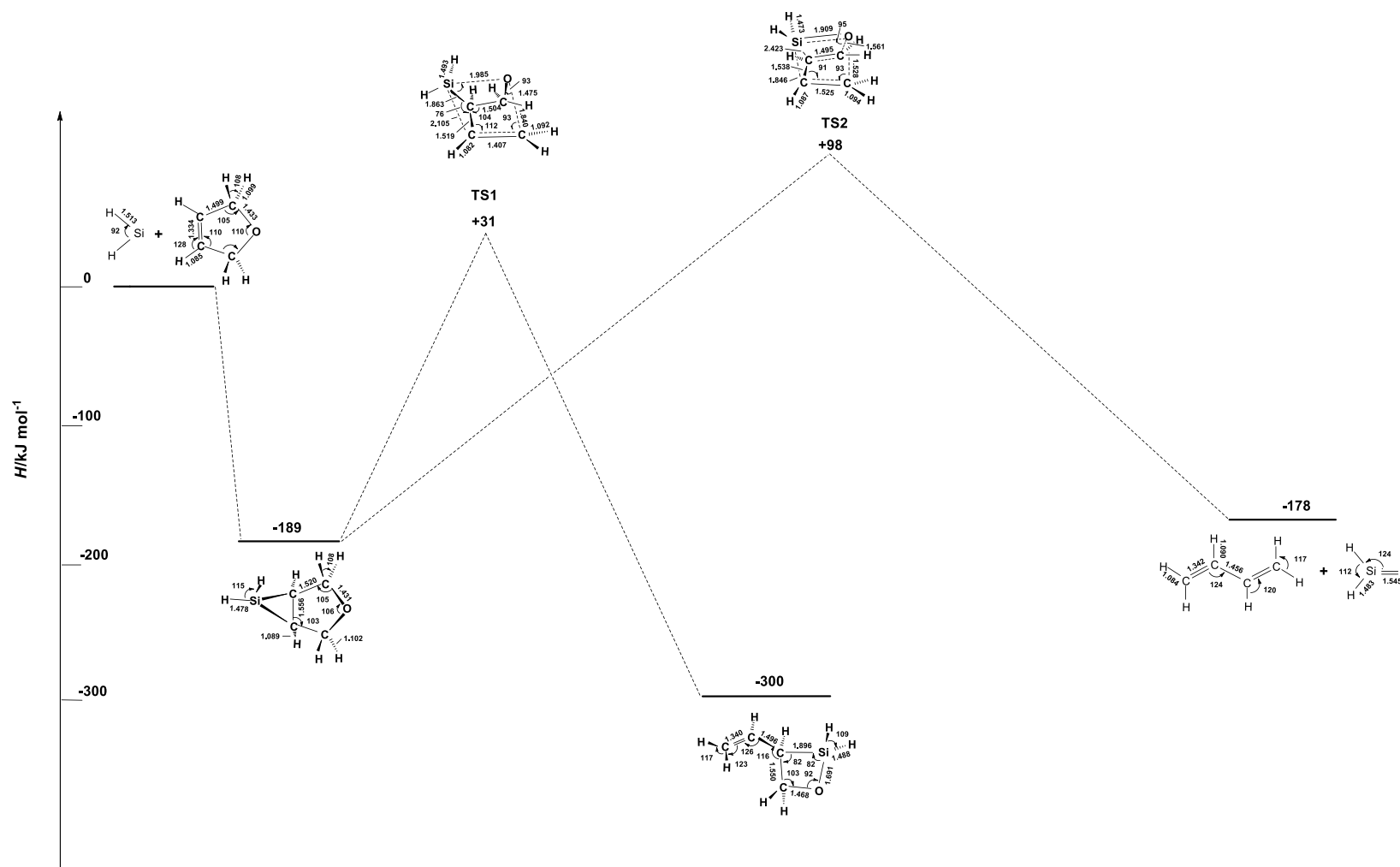
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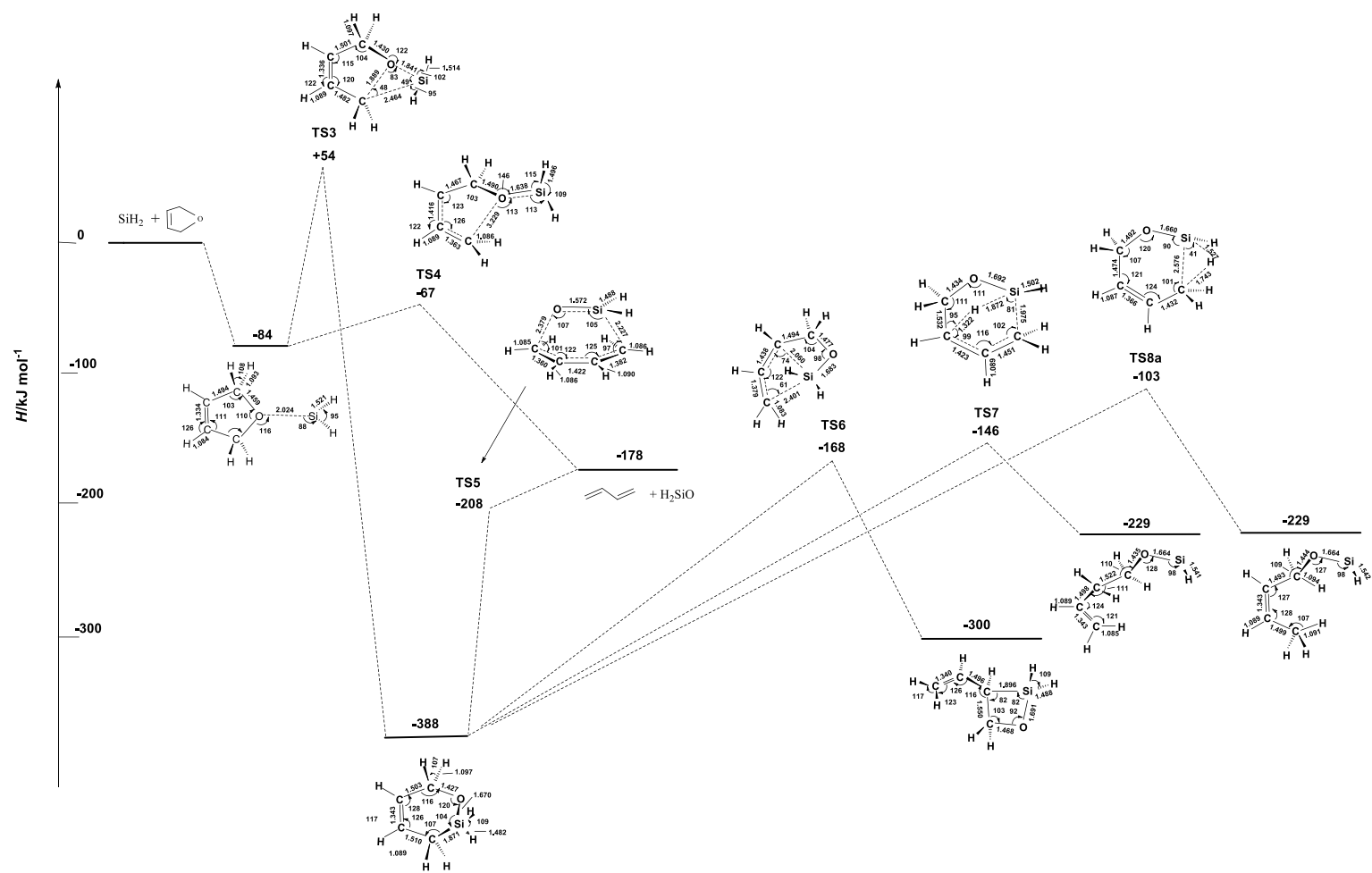
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**Figure 4.** Potential energy (enthalpy) surface for the  $\pi$ -addition pathway in the reaction of  $\text{SiH}_2$  + 2,5-DHF. All enthalpies are calculated at the G3 level and quoted at 298 K.



**Figure 5.** Potential energy (enthalpy) surface for the O-addition pathway in the reaction of  $\text{SiH}_2 + 2,5\text{-DHF}$ . All enthalpies are calculated at the G3 level and quoted at 298 K.