

Experimental Study and Simulation of Kinetics of Acetophenone Hydrosilylation with Diphenylsilane in the Presence of Rhodium Complexes in a Microreactor

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Abstract—The hydrosilylation of acetophenone with diphenylsilane in a microreactor in the presence of complexes $[\text{Rh}(\text{cod})\text{Cl}]_2$ and $[\text{Rh}(\text{CO})_2(\mu\text{-Cl})]_2$ and $(R)\text{--}(-)\text{-}cis\text{-mirtanyl-}$ and $(R)\text{--}(+)\text{-bornylamine}$ in situ was studied, the kinetics simulation of the process was performed, and the multicriteria optimization of the process was carried out. The influence of the micro-mixing effect on the reaction rate was revealed. Best results in the microreactor were obtained for the $[\text{Rh}(\text{cod})\text{Cl}]_2\text{--BornylNH}_2$ catalytic system. It was established that the formation of 1-phenylethanol and related enol silyl ethers are simultaneous competing reactions.

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