

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

V. M. Uvarov^a, E. S. Borovinskaya^a, D. A. de Vekki^a, and V. P. Reshetilovskii^b

^a *St. Petersburg State Institute of Technology, Moscovskii pr. 26, St. Petersburg, 190013 Russia
e-mail: hydrosilation@newmail.ru*

^b *Technical University of Dresden, 01062, Dresden, Germany*

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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*)-(-)-*cis*-mirtanyl- and (*R*)-(+)-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 $[\text{Rh}(\text{cod})\text{Cl}]_2$ –BornylNH₂ catalytic system. It was established that the formation of 1-phenylethanol and related enol silyl ethers are simultaneous competing reactions.

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