

Judit

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SHORT BIO:

I was born and raised in Budapest, Hungary, except for a few early years that I lived in Moscow, Russia. While obtaining a Degree in Chemistry at the Eötvös University, I worked in the group of [Prof. Tamás Turányi](#) to study correlations and uncertainties in combustion kinetic models. I spent the last year of my undergraduate studies in Leeds, U.K., as an Erasmus student working with [Prof. Mike Pilling](#), who introduced me to atmospheric chemistry. After completing my PhD in Hungary on further analysis of chemical kinetic models, I spent a year doing discharge flow experiments at the Chemical Research Centre of the Hungarian Academy of Sciences, mentored by [Dr. Sándor Dóbe](#). I started working at the Combustion Chemistry Department of Sandia in 2007 under the supervision of Dr. Craig A. Taatjes and Dr. James A. Miller (now at [Argonne](#)), studying a wide variety of combustion relevant elementary reactions, mostly using theoretical methods. I was hired as a senior staff member in the Chemistry Department in 2010.

RESEARCH INTERESTS

Theoretical reaction kinetics; Combustion chemistry; Atmospheric chemistry; Uncertainty analysis; Automatic reaction pathway search

EDUCATION

PhD in Physical Chemistry, Eötvös University, Budapest, Hungary (2006); Chemistry Diploma, Eötvös University, Budapest, Hungary (2002); Erasmus student, University of Leeds, Leeds, United Kingdom (2005-6)

AWARDS, HONORS AND MEMBERSHIPS

Distinguished Paper on Reaction Kinetics at the 32nd International Combustion Symposium (2009); Hungarian Chemical Society's Award for Outstanding Master Theses (2002); 1st in the Hungarian National High School Chemistry Contest (1997)

Combustion Institute (2007-); Reaction Kinetics and Photochemistry Working Committee of the Hungarian Academy of Sciences (2000-); Royal Society of Chemistry (2005-)

MENTEES

Brian D. Busemeyer, summer intern, 2012, now PhD student

Ewa Papajak, postdoctoral associate, 2012-

Xiaohu Li, postdoctoral associate, 2012-

Justin Kwok, summer intern, 2013

SELECTED PUBLICATIONS & PATENTS

1. **Zádor, J.**, Huang, H., Welz, O., Zetterberg, J., Osborn, D.L., Taatjes, C.A.: *Directly measuring reaction kinetics of QOOH – a crucial but elusive intermediate in hydrocarbon autoignition*. Physical Chemistry Chemical Physics, **2013** 15 10753-10760.
2. Welz, O., Klippenstein, S.J., Harding, L.B., Taatjes, C.A., **Zádor, J.**: *Unconventional peroxy chemistry in alcohol oxidation: The water elimination pathway*. Journal of Physical Chemistry Letters, **2013** 4 350-354.
3. **Zádor, J.**, Taatjes, C.A., and Fernandes, R.X.: *Kinetics of elementary reactions in low-temperature autoignition chemistry*. Progress in Energy and Combustion Science, **2011** 37 371-421.
4. **Zádor, J.**, Jasper, A.W., and Miller, J.A.: *The reaction between propene and hydroxyl*. Physical Chemistry Chemical Physics, **2009** 11 11040–11053.
5. Fernandes, R.X., **Zádor, J.**, Jusinski, L.E., Taatjes, C.A., and Miller, J.A.: *Formally direct pathways and low-temperature chain branching in hydrocarbon autoignition: The cyclohexyl + O₂ reaction at high pressure*. Physical Chemistry Chemical Physics, **2009** 11 1320-1327.
6. **Zádor, J.**, Fernandes, R.X., Georgievskii, Y., Meloni, G., Taatjes, C.A., and Miller, J.A.: *The reaction of hydroxyethyl radicals with O₂: A theoretical analysis and experimental product study*. Proceedings of the Combustion Institute, **2009** 32 271-277.
7. Kovács, G., **Zádor, J.**, Farkas, E., Nádasdi, R., Szilágyi, I., Dóbbé, S., Bérces, T., Márta, F., and Lendvai, G.: *Kinetics and mechanism of the reactions of CH₃CO and CH₃C(O)CH₂ radicals with O₂. Low-pressure discharge flow experiments and quantum chemical computations*. Physical Chemistry Chemical Physics, **2007** 9 4142-4154.
8. **Zádor, J.**, Wagner, V., Wirtz, K., and Pilling, M.J.: *Quantitative assessment of uncertainties for a model of tropospheric ethene oxidation using the European Photoreactor (EUPHORE)*. Atmospheric Environment, **2005** 39 2805-2817.
9. Zsély, I.G., **Zádor, J.**, and Turányi, T.: *Similarity of sensitivity functions of reaction kinetic models*. Journal of Physical Chemistry A, **2003** 107 2216-2238.