


<b>Title, Name and surname</b> Assist. Prof. Tomasz Prokop, PhD	
<b>Faculty and Department</b> Faculty of Energy and Fuels, Dep. Of Fundamental Research in Energy Engineering	
<b>Keywords.</b> computational material engineering, numerical simulation of heat transfer and fluid flow, numerical methods, computational fluid mechanics, mathematical modeling, fuel cells, energy, optimization	
<b>Scientific profile.</b> My research interests are related to computer modeling of mass and charge transport in materials with complex, evolving microstructures such as porous electrodes. Although I have generally wrote my own code (in C++, MATLAB, Python) for the purpose of numerical simulation and scientific visualization, I also have experience with ANSYS Fluent. I have taught courses on mathematical and numerical modeling, optimization, programming, and analysis of energy systems.	
<b>Exemplary thesis titles</b> -Numerical analysis of the microstructure of a porous material using nanotomographic data. -Analysis of reaction site density in a digital material representation of a porous electrode -Quantification of phase boundaries in a digital reconstruction of a solid oxide fuel cell microstructure	
<b>The form of conducting master's theses</b> Students participate in weekly meetings where the research progress and plans for the next week's are presented. The meeting is organized together with Ph.D. students and young staff members.	