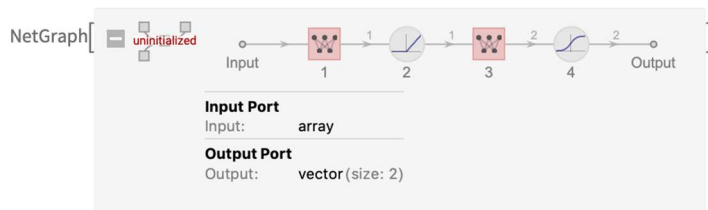


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|---|---|
| <input checked="" type="checkbox"/> Bachelor's thesis | <input checked="" type="checkbox"/> Computational |
| <input checked="" type="checkbox"/> Plant design exercise | <input type="checkbox"/> Experimental |
| <input checked="" type="checkbox"/> Paid master's thesis | <input type="checkbox"/> Plant design related |

Application of Machine Learning Methods to the Modeling of Fuel Cells and Chemical Reactors

The goal of this topic, which is open to several BSc / MSc / plant design theses, is to apply machine learning methods to model process engineering plants involving chemical reactions, including in particular fuel cells. For this purpose, approaches that have so far been used, e.g., for search engine rankings, for predictions of user



behavior or in the field of image and speech recognition [1], shall be applied to industrial operating data of process engineering plants. Examples for such

approaches are artificial neural networks or gaussian processes. For this goal, building blocks in Wolfram Mathematica from previous work can be used [2-4]. This topic requires a basic familiarity with the Wolfram Language concepts, e.g., by exploring the website <http://www.wolfram.com/wolfram-u/an-elementary-introduction-to-the-wolfram-language/>. The theses can largely be done from the home office.

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Literature: [1] <https://www.wolfram.com/language/introduction-machine-learning/>
 [2] <https://doi.org/10.1016/j.asoc.2021.107938>
 [3] <https://doi.org/10.1016/j.compchemeng.2021.107510>
 [4] <https://doi.org/10.1016/j.compchemeng.2022.107823>