

# Numerical methods for the simulation of dispersed phase systems

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Dispersed phase systems occur in chemical engineering in a large variety. They play an important role in industrial production processes, such like e.g. crystallization, granulation, polymerization or liquid-liquid extraction. A common characteristic of all those systems is that the dispersed phase consisting of e.g. particles, bubbles or drops is embedded in a continuous medium. A commonly accepted concept for the modeling of dispersed phase systems is the population balance approach [1,2,3].

The mathematical modeling of dispersed phase systems leads in general to partial integro-differential equations. For their solution different numerical methods will be discussed. The recently published 'Method of Space-Time Conservation Element and Solution Element' [4], which was originally designed for solving the Navier-Stokes and Euler Equations, was extended for the treatment of partial integro-differential equations. This rather new method will be compared to state of the art 'Method of Lines' based schemes, e.g. Finite Volume, Robust Upwind [5,6]. This comparison will be performed using a simple test model as well as a complex population balance model accounting for a continuously operated crystallizer. From the application of the discussed methods guidelines for a proper selection of numerical methods for the simulation of dispersed phase systems will be derived.

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