Smoldyn is one of the most widely used programs for stochastic simulation of individual molecules within biochemical networks. In this workshop Steve will provide an overview of Smoldyn and its capabilities followed by a hands on tutorial (no prior experience with mathematical modelling is required). Participants are encouraged to bring along specific problems that they wish to simulate.

http://www.smoldyn.org/index.html
http://www.smoldyn.org/andrews/index.html