

# **Molecular Dynamics Simulation: Elementary Methods (Monographs In Physical Chemistry Series) By J. M. Haile**

**By J. M. Haile**

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a Department of Physical Chemistry, we performed a series of molecular dynamic simulations to study the Molecular Dynamics Simulation: Elementary Methods.  
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A molecular dynamics simulation requires the definition of a potential function, or a description of the terms by which the particles in the simulation will interact.  
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