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Application of the Monte Carlo method for creation of initial models of EAP molecules for Molecular Dynamics simulation Endel Soolo, Jaanus Karo, Heiki Kasem agi, Maarja Kruusmaa and Alvo Aabloo Institute of Technology, Tartu University, Nooruse 1, 50411 Tartu, Estonia ABSTRACT A program called mcgen was written for creating initial models for Molecular Dynamics simulations with capability to

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