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An Ab Initio Molecular Dynamics Simulation

Guolin Xiong, Chunhong Yang, Weihua Zhu

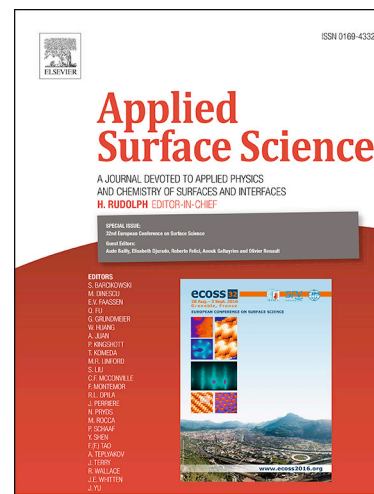
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Guolin Xiong, Chunhong Yang, Weihua Zhu*

Institute for Computation in Molecular and Materials Science, School of Chemical Engineering, Nanjing University of Science and Technology, Nanjing 210094, China

* E-mail: zhuwh@njust.edu.cn. Tel.: 86-25-84303919. Fax: 86-25-84303919

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