

Processes of Forming of Nano-Structures during Concretion of Clusters to Surface of Crystals with the Helps of Computer Modeling

A.M. Rasulov, A.A. Dzhurahalov, I.D. Yadgarov, V.G. Stelmakh

Abstract

A report is presented by computer simulation method about progress in the understanding of the properties of metallic nano-dimension particles, their interaction with surfaces subsequent to low energy slowing down and the properties of nano-structured materials formed with these particles.

Key words

Computer simulation, cluster, nanostructure, method molecular dynamics, introduction, thin films, Linked-Cell algorithm.

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Information about the Authors

A.M. Rasulov. Andizhan State University. Professor of Department of Informatics. Doctor of Physics and Mathematica Sciences. E-mail: arasulov59@mail.ru

A.A. Dzhurokhalov. Full Professor in University of Antwerp. Doctor of Physics and Mathematica Sciences. E-mail: dzhurakhalov@mail.ru

I.D. Yadgarov. Institute of Ion-Plasma and Laser Technologies Uzbek Academy of Sciences. PhD. Major Science researcher. E-mail: iyadgarov@mail.ru

V.G. Stelmakh. Institute of Ion-Plasma and Laser Technologies Uzbek Academy of Sciences. Science researcher. E-mail: stel@mail.ru

D.K. Akhmadaliev. Andizhan State University. Science researcher of Department of Informatics. E-mail: ilmiyor@mail.ru