2019 | Volume Volume - 4 - Issue Issue - 1

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Research Article

Open Access Research Article PTZAID:IJNNN-4-124

Simulation of the Behavior of Carbon Nanotori during Unfolding: A Study of **Stability and Electronic Structure**

Published On: January 29, 2018 | Pages: 004 - 008

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In this work, we present an in silico study of the dynamic behavior of a carbon nanotorus during the localized breaking of the C-C bonds in its atomic network. It is shown that the unfolding of a carbon nanotorus is accompanied by the appearance of deformation waves running along the atomic network of the carbon structure. We estimate energy stability of a carbon n ...

Abstract View Full Article View DOI: 10.17352/2455-3492.000024

Mini Review

Open Access Mini Review PTZAID:IJNNN-4-123

Sintering process and catalysis

Published On: January 08, 2018 | Pages: 001 - 003

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Sintering process has strong effect on the active surface area and catalytically active sites involved the catalyst. This process has two mechanisms depending on the thermal diffusion of the constituents of catalysts. ...

Abstract View Full Article View DOI: 10.17352/2455-3492.000023