Accepted Manuscript

Combining dual domain material point method with molecular dynamics for thermodynamic nonequilibriums

Tilak R. Dhakal, Duan Z. Zhang

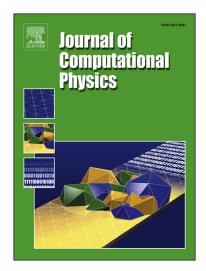
PII: S0021-9991(18)30529-1

DOI: https://doi.org/10.1016/j.jcp.2018.08.007

Reference: YJCPH 8192

To appear in: Journal of Computational Physics

Received date: 19 October 2017 Revised date: 24 May 2018 Accepted date: 2 August 2018



Please cite this article in press as: T.R. Dhakal, D.Z. Zhang, Combining dual domain material point method with molecular dynamics for thermodynamic nonequilibriums, *J. Comput. Phys.* (2018), https://doi.org/10.1016/j.jcp.2018.08.007

This is a PDF file of an unedited manuscript that has been accepted for publication. As a service to our customers we are providing this early version of the manuscript. The manuscript will undergo copyediting, typesetting, and review of the resulting proof before it is published in its final form. Please note that during the production process errors may be discovered which could affect the content, and all legal disclaimers that apply to the journal pertain.

Highlights

- Introduced a multiscale method based on material point method and molecular dynamics.
- As an example, computed a large deformation problem at thermodynamic nonequilibrium.
- \bullet The method preserves the entire material deformation history without re-initialization.
- Studied consistence conditions for inter-scale communications.

Download English Version:

https://daneshyari.com/en/article/11002780

Download Persian Version:

https://daneshyari.com/article/11002780

<u>Daneshyari.com</u>