

Accepted Manuscript

Configurational effect on ion-pair interaction energies and intermolecular potential energy functions in imidazolium-based ionic liquids: A theoretical study

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PII: S0167-7322(18)30884-5
DOI: doi:[10.1016/j.molliq.2018.04.105](https://doi.org/10.1016/j.molliq.2018.04.105)
Reference: MOLLIQ 9008
To appear in: *Journal of Molecular Liquids*
Received date: 18 February 2018
Revised date: 17 April 2018
Accepted date: 20 April 2018

Please cite this article as: Motahare Sadeghi Googheri, Mohammad Sadegh Sadeghi Googheri, Samira Hozhabr Araghi , Configurational effect on ion-pair interaction energies and intermolecular potential energy functions in imidazolium-based ionic liquids: A theoretical study. The address for the corresponding author was captured as affiliation for all authors. Please check if appropriate. Molliq(2017), doi:[10.1016/j.molliq.2018.04.105](https://doi.org/10.1016/j.molliq.2018.04.105)

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