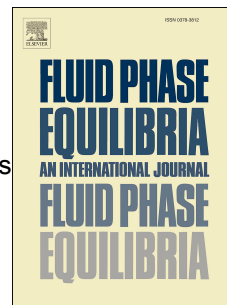


Accepted Manuscript

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PII: S0378-3812(18)30199-7

DOI: [10.1016/j.fluid.2018.05.013](https://doi.org/10.1016/j.fluid.2018.05.013)

Reference: FLUID 11835

To appear in: *Fluid Phase Equilibria*

Received Date: 20 February 2018

Revised Date: 7 May 2018

Accepted Date: 11 May 2018

Please cite this article as: R. Gillet, Angé. Fierro, L.M. Valenzuela, José.R. Pérez-Correa, Using molecular dynamics simulations to predict the effect of temperature on aqueous solubility for aromatic compounds, *Fluid Phase Equilibria* (2018), doi: 10.1016/j.fluid.2018.05.013.

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Using molecular dynamics simulations to predict the effect of temperature on aqueous solubility for aromatic compounds

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