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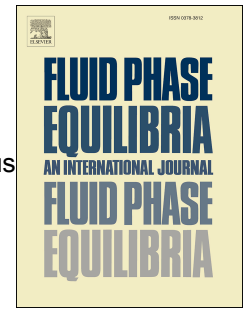

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Using molecular dynamics simulations to predict the effect of temperature on aqueous solubility for aromatic compounds

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