

# Karnesh Jain

## List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Folding Free-Energy Landscape of Î±-Synuclein (35â€“97) Via Replica Exchange Molecular Dynamics. Journal of Chemical Information and Modeling, 2021, 61, 432-443.	5.4	10
2	Application of the interface potential approach for studying wetting behavior within a molecular dynamics framework. Journal of Chemical Physics, 2019, 150, 204118.	3.0	8
3	Construction of the interface potential from a series of canonical ensemble simulations. Journal of Chemical Physics, 2019, 151, 044103.	3.0	4
4	Using isothermal-isobaric Monte Carlo simulation to study the wetting behavior of model systems. Journal of Chemical Physics, 2019, 150, 084110.	3.0	9