

# Minerva Gonzalez-Melchor

## List of Publications by Year in descending order

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24  
papers

639  
citations

759233

12  
h-index

642732

23  
g-index

24  
all docs

24  
docs citations

24  
times ranked

598  
citing authors

#	ARTICLE	IF	CITATIONS
1	Exploring electrostatic patterns of human, murine, equine and canine TLR4/MD-2 receptors. <i>Innate Immunity</i> , 2020, 26, 364-380.	2.4	2
2	Self-diffusion and structure of monovalent ions in two dimensions: A molecular dynamics study. <i>Journal of Molecular Liquids</i> , 2019, 294, 111542.	4.9	1
3	Influence of pH on the formation of a polyelectrolyte complex by dissipative particle dynamics simulation: From an extended to a compact shape. <i>Physical Review E</i> , 2019, 100, 012505.	2.1	3
4	Static dielectric constant of water within a bilayer using recent water models: a molecular dynamics study. <i>Journal of Physics Condensed Matter</i> , 2018, 30, 195001.	1.8	10
5	Dopamine and Caffeine Encapsulation within Boron Nitride (14,0) Nanotubes: Classical Molecular Dynamics and First Principles Calculations. <i>Journal of Physical Chemistry B</i> , 2018, 122, 5885-5896.	2.6	18
6	The line tension of two-dimensional ionic fluids. <i>Journal of Chemical Physics</i> , 2016, 144, 134705.	3.0	3
7	The structure and interaction mechanism of a polyelectrolyte complex: a dissipative particle dynamics study. <i>Soft Matter</i> , 2015, 11, 5889-5897.	2.7	24
8	Liquid-vapor interface varying the softness and range of the interaction potential. <i>Molecular Simulation</i> , 2013, 39, 64-71.	2.0	8
9	Influence of a neutral component on the liquid-vapor coexistence and the surface tension of an ionic fluid. <i>Journal of Molecular Liquids</i> , 2013, 185, 32-35.	4.9	0
10	First principles studies of the graphene-phenol interactions. <i>Journal of Molecular Modeling</i> , 2012, 18, 3857-3866.	1.8	36
11	Interfacial and coexistence properties of soft spheres with a short-range attractive Yukawa fluid: Molecular dynamics simulations. <i>Journal of Chemical Physics</i> , 2012, 136, 154702.	3.0	8
12	Liquid-vapor phase diagram and cluster formation of two-dimensional ionic fluids. <i>Journal of Chemical Physics</i> , 2012, 137, 054711.	3.0	13
13	Analytical static structure factors for the restricted primitive model. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2012, 391, 1759-1769.	2.6	2
14	Equilibrium structure of the multi-component screened charged hard-sphere fluid. <i>Journal of Chemical Physics</i> , 2011, 135, 014504.	3.0	8
15	Interfacial properties of charge asymmetric ionic liquids. <i>Molecular Physics</i> , 2009, 107, 357-363.	1.7	12
16	Effect of softness of the potential on the stress anisotropy in liquids. <i>Journal of Chemical Physics</i> , 2007, 126, 224511.	3.0	9
17	Finite-size effects in dissipative particle dynamics simulations. <i>Journal of Chemical Physics</i> , 2006, 124, 084104.	3.0	41
18	Electrostatic interactions in dissipative particle dynamics using the Ewald sums. <i>Journal of Chemical Physics</i> , 2006, 125, 224107.	3.0	163

#	ARTICLE	IF	CITATIONS
19	Stress anisotropy induced by periodic boundary conditions. Journal of Chemical Physics, 2005, 122, 094503.	3.0	60
20	Molecular dynamics simulations of the surface tension of ionic liquids. Journal of Chemical Physics, 2005, 122, 104710.	3.0	88
21	Influence of ion size asymmetry on the properties of ionic liquid-vapour interfaces. Journal of Physics Condensed Matter, 2005, 17, S3301-S3307.	1.8	21
22	Surface Tension of the Restrictive Primitive Model for Ionic Liquids. Physical Review Letters, 2003, 90, 135506.	7.8	43
23	Surface tension at the vapor/liquid interface in an attractive hard-core Yukawa fluid. Journal of Chemical Physics, 2001, 115, 3862-3872.	3.0	44
24	Equation of state and structure of binary mixtures of hard d-dimensional hyperspheres. Journal of Chemical Physics, 2001, 114, 4905-4911.	3.0	22