

Orlando Guzmán

List of Publications by Year in descending order

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docs citations

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483
citing authors

#	ARTICLE	IF	CITATIONS
1	Nucleation and growth of blue phase liquid crystals on chemically-patterned surfaces: a surface anchoring assisted blue phase correlation length. <i>Molecular Systems Design and Engineering</i> , 2021, 6, 534-544.	3.4	3
2	Dynamics of Nanoparticle Self-Assembly by Liquid Crystal Sorting in Two Dimensions. <i>Frontiers in Physics</i> , 2021, 9, .	2.1	1
3	Control of Monodomain Polymer-Stabilized Cuboidal Nanocrystals of Chiral Nematics by Confinement. <i>ACS Nano</i> , 2021, 15, 15972-15981.	14.6	10
4	Self-assembling and phase coexistence of SW trimers as complex amphiphile analogues. I. Simulations. <i>Molecular Physics</i> , 2020, 118, e1726519.	1.7	1
5	Specific inter-domain interactions stabilize a compact HIV-1 Gag conformation. <i>PLoS ONE</i> , 2019, 14, e0221256.	2.5	2
6	Sculpted grain boundaries in soft crystals. <i>Science Advances</i> , 2019, 5, eaax9112.	10.3	18
7	Global square-well free-energy model via singular value decomposition. <i>Molecular Physics</i> , 2018, 116, 2070-2082.	1.7	7
8	Free-energy model for nanoparticle self-assembly by liquid crystal sorting. <i>Physical Review E</i> , 2018, 97, 062704.	2.1	11
9	Liquid-Vapor Equilibria of Ionic Liquids from a SAFT Equation of State with Explicit Electrostatic Free Energy Contributions. <i>Journal of Physical Chemistry B</i> , 2015, 119, 5864-5872.	2.6	5
10	Room temperature ionic liquids: A simple model. Effect of chain length and size of intermolecular potential on critical temperature. <i>Journal of Chemical Physics</i> , 2015, 142, 154508.	3.0	3
11	Theoretically informed Monte Carlo simulation of liquid crystals by sampling of alignment-tensor fields. <i>Journal of Chemical Physics</i> , 2015, 143, 044107.	3.0	22
12	Measuring liquid crystal elastic constants with free energy perturbations. <i>Soft Matter</i> , 2014, 10, 882-893.	2.7	42
13	Self-assembly of kagome lattices, entangled webs and linear fibers with vibrating patchy particles in two dimensions. <i>Soft Matter</i> , 2014, 10, 9167-9176.	2.7	17
14	Analytical equation of state with three-body forces: Application to noble gases. <i>Journal of Chemical Physics</i> , 2013, 139, 184503.	3.0	20
15	Steric contribution of macromolecular crowding to the time and activation energy for preprotein translocation across the endoplasmic reticulum membrane. <i>Physical Review E</i> , 2013, 88, 012725.	2.1	2
16	Systematic prediction of critical point coordinates from molecular parameters of equations of state and interaction potentials. <i>Molecular Physics</i> , 2012, 110, 1261-1267.	1.7	2
17	High-precision virial coefficients of argon and carbon dioxide from integration of speed of sound data in the pressure-temperature domain. <i>Molecular Physics</i> , 2012, 110, 1349-1358.	1.7	12
18	Effective potential for three-body forces in fluids. <i>Molecular Physics</i> , 2011, 109, 955-967.	1.7	16

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19	Boundary-layer method for the analytical calculation of stable textures of bent-core liquid crystal fibers. <i>Physical Review E</i> , 2011, 84, 011701.	2.1	2
20	Third Virial Coefficients of Mixtures from a Model of Two- and Three-body Forces. , 2010, , .		0
21	Third virial coefficient of nonpolar gases from accurate binary potentials and ternary forces. <i>Journal of Physics B: Atomic, Molecular and Optical Physics</i> , 2007, 40, 3989-4003.	1.5	15
22	Effective intermolecular potentials in theoretical thermodynamics of pure substances and solutions. <i>Fluid Phase Equilibria</i> , 2007, 259, 9-22.	2.5	17
23	Anisotropic nanoparticles immersed in a nematic liquid crystal: Defect structures and potentials of mean force. <i>Physical Review E</i> , 2006, 74, 011711.	2.1	63
24	Measurement of the Azimuthal Anchoring Energy of Liquid Crystals in Contact with Oligo(ethylene) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 5 <i>Langmuir</i> , 2006, 22, 4654-4659.	3.5	31
25	Anchoring Energies of Liquid Crystals Measured on Surfaces Presenting Oligopeptides. <i>Langmuir</i> , 2006, 22, 7776-7782.	3.5	19
26	Interactions of Liquid Crystal-Forming Molecules with Phospholipid Bilayers Studied by Molecular Dynamics Simulations. <i>Biophysical Journal</i> , 2005, 89, 3141-3158.	0.5	20
27	Multiscale Simulation of Liquid Crystals. <i>NATO Science Series Series II, Mathematics, Physics and Chemistry</i> , 2005, , 221-247.	0.1	0
28	Interactions between spherical colloids mediated by a liquid crystal: A molecular simulation and mesoscale study. <i>Journal of Chemical Physics</i> , 2004, 121, 1949-1961.	3.0	49
29	An effective-colloid pair potential for Lennard-Jones colloidâ€“polymer mixtures. <i>Journal of Chemical Physics</i> , 2003, 118, 2392-2397.	3.0	17
30	Theoretical Equation of State of Dense Nonconformal Fluids from Effective Potentials. 1. Applications to Model Systems. <i>Journal of Physical Chemistry B</i> , 2001, 105, 8220-8229.	2.6	8
31	Phase-shift symmetries of the correlation and bridge functions in additive hard sphere mixtures. <i>Molecular Physics</i> , 1998, 95, 645-648.	1.7	3
32	An Integral Equation and Monte Carlo Study of Square-Well Fluid Mixtures. <i>The Journal of Physical Chemistry</i> , 1995, 99, 1587-1593.	2.9	10