

Marco Lauricella

List of Publications by Year in descending order

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63
papers

994
citations

393982

19
h-index

476904

29
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72
all docs

72
docs citations

72
times ranked

823
citing authors

#	ARTICLE	IF	CITATIONS
1	Electric field induced macroscopic cellular phase of nanoparticles. <i>Soft Matter</i> , 2022, 18, 1991-1996.	1.2	2
2	Capturing Free-Radical Polymerization by Synergetic <i>Ab Initio</i> Calculations and Topological Reactive Molecular Dynamics. <i>Macromolecules</i> , 2022, 55, 1474-1486.	2.2	3
3	Double Life of Methanol: Experimental Studies and Nonequilibrium Molecular-Dynamics Simulation of Methanol Effects on Methane-Hydrate Nucleation. <i>Journal of Physical Chemistry C</i> , 2022, 126, 6075-6081.	1.5	9
4	Stochastic Jetting and Dripping in Confined Soft Granular Flows. <i>Physical Review Letters</i> , 2022, 128, 128001.	2.9	9
5	LBcuda: A high-performance CUDA port of LBsoft for simulation of colloidal systems. <i>Computer Physics Communications</i> , 2022, 277, 108380.	3.0	6
6	Computational droplets: Where we stand and how far we can go. <i>Europhysics Letters</i> , 2022, 138, 67001.	0.7	2
7	The vortex-driven dynamics of droplets within droplets. <i>Nature Communications</i> , 2021, 12, 82.	5.8	26
8	Mesoscale modelling of droplets' self-assembly in microfluidic channels. <i>Soft Matter</i> , 2021, 17, 2374-2383.	1.2	11
9	Wet to dry self-transitions in dense emulsions: From order to disorder and back. <i>Physical Review Fluids</i> , 2021, 6, .	1.0	11
10	Optimized Modeling and Design of a PCM-Enhanced H2 Storage. <i>Energies</i> , 2021, 14, 1554.	1.6	17
11	Shear dynamics of polydisperse double emulsions. <i>Physics of Fluids</i> , 2021, 33, .	1.6	10
12	Lattice Boltzmann multicomponent model for direct-writing printing. <i>Physics of Fluids</i> , 2021, 33, .	1.6	6
13	Translocation Dynamics of High-Internal Phase Double Emulsions in Narrow Channels. <i>Langmuir</i> , 2021, 37, 9026-9033.	1.6	11
14	Microscale modelling of dielectrophoresis assembly processes. <i>Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences</i> , 2021, 379, 20200407.	1.6	2
15	Tracking droplets in soft granular flows with deep learning techniques. <i>European Physical Journal Plus</i> , 2021, 136, 864.	1.2	8
16	A fast and efficient deep learning procedure for tracking droplet motion in dense microfluidic emulsions. <i>Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences</i> , 2021, 379, 20200400.	1.6	10
17	Dynamics of polydisperse multiple emulsions in microfluidic channels. <i>Physical Review E</i> , 2021, 104, 065112.	0.8	1
18	Toward exascale design of soft mesoscale materials. <i>Journal of Computational Science</i> , 2020, 46, 101175.	1.5	6

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19	LBsoft: A parallel open-source software for simulation of colloidal systems. <i>Computer Physics Communications</i> , 2020, 256, 107455.	3.0	10
20	Shear dynamics of confined bijels. <i>AIP Advances</i> , 2020, 10, 095304.	0.6	8
21	Models of polymer solutions in electrified jets and solution blowing. <i>Reviews of Modern Physics</i> , 2020, 92, .	16.4	51
22	Multiparticle collision dynamics for fluid interfaces with near-contact interactions. <i>Journal of Chemical Physics</i> , 2020, 152, 144101.	1.2	4
23	A coupled lattice Boltzmann-Multiparticle collision method for multi-resolution hydrodynamics. <i>Journal of Computational Science</i> , 2020, 44, 101160.	1.5	1
24	A Multiresolution Mesoscale Approach for Microscale Hydrodynamics. <i>Advanced Theory and Simulations</i> , 2020, 3, 1900250.	1.3	2
25	Novel nonequilibrium steady states in multiple emulsions. <i>Physics of Fluids</i> , 2020, 32, .	1.6	20
26	Lattice Boltzmann simulations capture the multiscale physics of soft flowing crystals. <i>Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences</i> , 2020, 378, 20190406.	1.6	6
27	Concentrated phase emulsion with multicore morphology under shear: A numerical study. <i>Physical Review Fluids</i> , 2020, 5, .	1.0	10
28	Acoustic-propagation properties of methane clathrate hydrates from non-equilibrium molecular dynamics. <i>Journal of Chemical Physics</i> , 2019, 151, 144505.	1.2	8
29	Microvorticity fluctuations affect the structure of thin fluid films. <i>Physical Review E</i> , 2019, 100, 042606.	0.8	2
30	Simulating blood rheology across scales: A hybrid LB-particle approach. <i>International Journal of Modern Physics C</i> , 2019, 30, 1941003.	0.8	2
31	Mesoscale modelling of near-contact interactions for complex flowing interfaces. <i>Journal of Fluid Mechanics</i> , 2019, 872, 327-347.	1.4	48
32	Modeling realistic multiphase flows using a non-orthogonal multiple-relaxation-time lattice Boltzmann method. <i>Physics of Fluids</i> , 2019, 31, .	1.6	67
33	Mesoscale modelling of soft flowing crystals. <i>Philosophical Transactions Series A, Mathematical, Physical, and Engineering Sciences</i> , 2019, 377, 20180149.	1.6	16
34	Amplitude effects on seismic velocities: How low can we go?. <i>Journal of Chemical Physics</i> , 2019, 150, 084101.	1.2	1
35	Curvature dynamics and long-range effects on fluid-fluid interfaces with colloids. <i>Soft Matter</i> , 2019, 15, 2848-2862.	1.2	7
36	Jetting to dripping transition: Critical aspect ratio in step emulsifiers. <i>Physics of Fluids</i> , 2019, 31, .	1.6	25

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37	Ab initio accelerated molecular dynamics study of the hydride ligands in the ruthenium complex: Ru(H ₂) ₂ H ₂ (P(C ₅ H ₉)) ₃ . Physical Chemistry Chemical Physics, 2019, 21, 25247-25257.	1.3	1
38	Disordered interfaces in soft fluids with suspended colloids. International Journal of Modern Physics C, 2019, 30, 1941004.	0.8	1
39	Towards Exascale Lattice Boltzmann computing. Computers and Fluids, 2019, 181, 107-115.	1.3	40
40	Modeling pattern formation in soft flowing crystals. Physical Review Fluids, 2019, 4, .	1.0	30
41	Regularized lattice Boltzmann multicomponent models for low capillary and Reynolds microfluidics flows. Computers and Fluids, 2018, 167, 33-39.	1.3	33
42	Elastic Characterization of <i>S</i> - and <i>P</i> -Wave Velocities in Marine-like Silica: The Role of Nonequilibrium Molecular Dynamics. Journal of Physical Chemistry C, 2018, 122, 3006-3013.	1.5	2
43	Multicomponent Lattice Boltzmann Models for Biological Applications. , 2018, , 357-370.		1
44	Mechanistic modelling of drug release from multi-layer capsules. Computers in Biology and Medicine, 2018, 93, 149-157.	3.9	41
45	Entropic lattice Boltzmann model for charged leaky dielectric multiphase fluids in electrified jets. Physical Review E, 2018, 97, 033308.	0.8	19
46	Lattice propagators and Haldane-Wu fractional statistics. Europhysics Letters, 2018, 122, 10002.	0.7	0
47	Elucidating the mechanism of step emulsification. Physical Review Fluids, 2018, 3, .	1.0	27
48	Mesoscopic model for soft flowing systems with tunable viscosity ratio. Physical Review Fluids, 2018, 3, .	1.0	20
49	Mechanisms and Nucleation Rate of Methane Hydrate by Dynamical Nonequilibrium Molecular Dynamics. Journal of Physical Chemistry C, 2017, 121, 24223-24234.	1.5	30
50	Effects of orthogonal rotating electric fields on electrospinning process. Physics of Fluids, 2017, 29, .	1.6	20
51	Effects of nanoparticles on the dynamic morphology of electrified jets. Europhysics Letters, 2017, 119, 44001.	0.7	1
52	Dynamic mesh refinement for discrete models of jet electro-hydrodynamics. Journal of Computational Science, 2016, 17, 325-333.	1.5	12
53	Iontophoretic transdermal drug delivery: a multi-layered approach. Mathematical Medicine and Biology, 2016, 34, dqw017.	0.8	6
54	Three-Dimensional Model for Electrospinning Processes in Controlled Gas Counterflow. Journal of Physical Chemistry A, 2016, 120, 4884-4892.	1.1	15

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55	Structural and dynamical properties of methane clathrate hydrates from molecular dynamics: Comparison of atomistic and more coarse-grained potential models. <i>Fluid Phase Equilibria</i> , 2016, 413, 235-241.	1.4	8
56	Nonlinear Langevin model for the early-stage dynamics of electrospinning jets. <i>Molecular Physics</i> , 2015, 113, 2435-2441.	0.8	9
57	Sub-ms dynamics of the instability onset of electrospinning. <i>Soft Matter</i> , 2015, 11, 3424-3431.	1.2	29
58	Clathrate structure-type recognition: Application to hydrate nucleation and crystallisation. <i>Journal of Chemical Physics</i> , 2015, 142, 244503.	1.2	33
59	JETSPIN: A specific-purpose open-source software for simulations of nanofiber electrospinning. <i>Computer Physics Communications</i> , 2015, 197, 227-238.	3.0	19
60	Different regimes of the uniaxial elongation of electrically charged viscoelastic jets due to dissipative air drag. <i>Mechanics Research Communications</i> , 2015, 69, 97-102.	1.0	11
61	Massively parallel molecular dynamics simulation of formation of clathrate-hydrate precursors at planar water-methane interfaces: Insights into heterogeneous nucleation. <i>Journal of Chemical Physics</i> , 2014, 140, 204714.	1.2	56
62	Methane Clathrate Hydrate Nucleation Mechanism by Advanced Molecular Simulations. <i>Journal of Physical Chemistry C</i> , 2014, 118, 22847-22857.	1.5	87
63	Experimental and computational investigation of the group 11 "group 2 diatomic molecules: First determination of the AuSr and AuBa bond energies and thermodynamic stability of the copper- and silver-alkaline earth species. <i>Journal of Chemical Physics</i> , 2012, 136, 184306.	1.2	5