

# João Puibasset

## List of Publications by Year in descending order

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66  
papers

1,402  
citations

377584

21  
h-index

371746

37  
g-index

66  
all docs

66  
docs citations

66  
times ranked

2832  
citing authors

#	ARTICLE	IF	CITATIONS
1	Silica-induced electron loss of silver nanoparticles. <i>Nanoscale</i> , 2022, 14, 7280-7291.	2.8	1
2	Bulk supercooled water <i>versus</i> adsorbed films on silica surfaces: specific heat by Monte Carlo simulation. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 2275-2285.	1.3	0
3	Cavitation in heterogeneous nanopores: The chemical ink-bottle. <i>AIP Advances</i> , 2021, 11, .	0.6	3
4	Density Functional Theory Study of the Spontaneous Formation of Covalent Bonds at the Silver/Silica Interface in Silver Nanoparticles Embedded in SiO <sub>2</sub> : Implications for Ag <sup>+</sup> Release. <i>ACS Applied Nano Materials</i> , 2019, 2, 5179-5189.	2.4	13
5	Elastic Compliance and Stiffness Matrix of the FCC Lennard-Jones Thin Films: Influence of Thickness and Temperature. <i>Journal of Physical Chemistry C</i> , 2019, 123, 15027-15037.	1.5	1
6	Molecular simulation study of the heat capacity of metastable water between 100 and 300 K. <i>Molecular Simulation</i> , 2019, 45, 462-465.	0.9	2
7	Freezing and Melting of Silver Nanoparticles on Silica Substrate Using a Simple Interatomic Potential for Ag-SiO <sub>2</sub> Interaction on the Basis of <i>ab Initio</i> Calculations and Experimental Data. <i>Journal of Physical Chemistry C</i> , 2017, 121, 3615-3622.	1.5	15
8	On the Thermodynamics and Experimental Control of Twinning in Metal Nanocrystals. <i>Angewandte Chemie</i> , 2017, 129, 8773-8777.	1.6	6
9	On the Thermodynamics and Experimental Control of Twinning in Metal Nanocrystals. <i>Angewandte Chemie - International Edition</i> , 2017, 56, 8647-8651.	7.2	21
10	Adsorption-Induced Deformation of a Nanoporous Material: Influence of the Fluid-Adsorbent Interaction and Surface Freezing on the Pore-Load Modulus Measurement. <i>Journal of Physical Chemistry C</i> , 2017, 121, 18779-18788.	1.5	5
11	Adsorption-Induced Deformation in Nanopores: Unexpected Results Obtained by Molecular Simulations. , 2017, , .		0
12	Finite-size corrections in simulation of dipolar fluids. <i>Journal of Chemical Physics</i> , 2017, 147, 224110.	1.2	5
13	Structure and Permeability of Porous Silicon Investigated by Self-Diffusion NMR Measurements of Ethanol and Heptane. <i>Oil and Gas Science and Technology</i> , 2016, 71, 54.	1.4	8
14	Influence of the Silica Support on the Structure and the Morphology of Silver Nanoparticles: A Molecular Simulation Study. <i>Journal of Physical Chemistry C</i> , 2016, 120, 8323-8332.	1.5	16
15	A SIMPLE bubble chamber for dark matter searches: Testing and development. , 2015, , .		0
16	Adsorption-induced strain of a nanoscale silicon honeycomb. <i>Europhysics Letters</i> , 2015, 109, 56002.	0.7	39
17	The simple dark matter search: Present and future. <i>E3S Web of Conferences</i> , 2014, 4, 03002.	0.2	0
18	Influence of system size on the properties of a fluid adsorbed in a nanopore: Physical manifestations and methodological consequences. <i>Journal of Chemical Physics</i> , 2014, 141, 044716.	1.2	5

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19	The SIMPLE Phase II dark matter search. <i>Physical Review D</i> , 2014, 89, .	1.6	45
20	Fluid adsorption in linear pores: a molecular simulation study of the influence of heterogeneities on the hysteresis loop and the distribution of metastable states. <i>Molecular Simulation</i> , 2014, 40, 690-697.	0.9	6
21	Fabrication and response of high concentration SIMPLE superheated droplet detectors with different liquids. <i>Astroparticle Physics</i> , 2013, 49, 28-43.	1.9	13
22	Phase transitions in porous materials: influence of physico-chemical heterogeneities. <i>MATEC Web of Conferences</i> , 2013, 3, 01083.	0.1	0
23	Improving Molecular Simulation Models of Adsorption in Porous Materials: Interdependence between Domains. <i>Oil and Gas Science and Technology</i> , 2013, 68, 309-318.	1.4	3
24	Final Analysis and Results of the Phase II SIMPLE Dark Matter Search. <i>Physical Review Letters</i> , 2012, 108, 201302.	2.9	228
25	Prospects for a Phase III SIMPLE Measurement. <i>Journal of Physics: Conference Series</i> , 2012, 375, 012017.	0.3	4
26	Bridge function for the dipolar fluid from simulation. <i>Journal of Chemical Physics</i> , 2012, 136, 154503.	1.2	19
27	Stability intervals of metastable states in hysteretic systems. <i>Physical Review E</i> , 2011, 84, 061126.	0.8	2
28	Numerical characterization of the density of metastable states within the hysteresis loop in disordered systems. <i>Journal of Physics Condensed Matter</i> , 2011, 23, 035106.	0.7	4
29	First Results of the Phase II SIMPLE Dark Matter Search. <i>Physical Review Letters</i> , 2010, 105, 211301.	2.9	62
30	Counting metastable states within the adsorption/desorption hysteresis loop: A molecular simulation study of confinement in heterogeneous pores. <i>Journal of Chemical Physics</i> , 2010, 133, 104701.	1.2	14
31	Vapor-Liquid Equilibrium. , 2010, , 213-240.		1
32	Effect of the reservoir size on gas adsorption in inhomogeneous porous media. <i>Journal of Physics Condensed Matter</i> , 2009, 21, 155102.	0.7	14
33	Influence of reservoir size on the adsorption path in an ideal pore. <i>Journal of Chemical Physics</i> , 2009, 131, 124123.	1.2	21
34	Monte-Carlo Multiscale Simulation Study of Argon Adsorption/Desorption Hysteresis in Mesoporous Heterogeneous Tubular Pores like MCM-41 or Oxidized Porous Silicon. <i>Langmuir</i> , 2009, 25, 903-911.	1.6	46
35	Molecular simulations of water in hydrophobic microporous solids. <i>Adsorption</i> , 2008, 14, 733-742.	1.4	24
36	Grand Canonical Monte Carlo Simulation Study of Water Adsorption in Silicalite at 300 K. <i>Journal of Physical Chemistry B</i> , 2008, 112, 6390-6397.	1.2	83

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37	Pseudocritical or hysteresis temperature versus pore size for simple fluids confined in cylindrical nanopores. <i>Journal of Chemical Physics</i> , 2008, 129, 024705.	1.2	6
38	Surface excess free energy of simple fluids confined in cylindrical pores by isothermal-isobaric Monte Carlo: Influence of pore size. <i>Journal of Chemical Physics</i> , 2007, 126, 184701.	1.2	7
39	Thermodynamic pressure of simple fluids confined in cylindrical nanopores by isothermal-isobaric Monte Carlo: Influence of fluid/substrate interactions. <i>Journal of Chemical Physics</i> , 2007, 127, 074702.	1.2	1
40	Influence of temperature on water adsorption / desorption hysteresis loop in disordered mesoporous silica glass by Grand Canonical Monte Carlo simulation method. <i>Studies in Surface Science and Catalysis</i> , 2007, , 535-541.	1.5	3
41	Effect of pore morphology and topology on capillary condensation in nanopores: a theoretical and molecular simulation study. <i>Studies in Surface Science and Catalysis</i> , 2007, 160, 1-8.	1.5	11
42	Adsorption-desorption hysteresis of simple fluids confined in realistic heterogeneous silica mesopores of micrometric length: A new analysis exploiting a multiscale Monte Carlo approach. <i>Journal of Chemical Physics</i> , 2007, 127, 154701.	1.2	40
43	Water confined in mesoporous silica glasses: Influence of temperature on adsorption/desorption hysteresis loop and fluid structure. <i>European Physical Journal: Special Topics</i> , 2007, 141, 41-44.	1.2	12
44	Generalized isobaric-isothermal ensemble: application to capillary condensation and cavitation in heterogeneous nanopores. <i>Molecular Physics</i> , 2006, 104, 3021-3032.	0.8	5
45	SIMPLE limits on spin-dependent WIMP interactions. <i>Journal of Physics: Conference Series</i> , 2006, 39, 114-116.	0.3	2
46	The simple SDD. <i>Radiation Protection Dosimetry</i> , 2006, 120, 503-508.	0.4	2
47	Influence of surface chemical heterogeneities on adsorption/desorption hysteresis and coexistence diagram of metastable states within cylindrical pores. <i>Journal of Chemical Physics</i> , 2006, 125, 074707.	1.2	23
48	RECENT RESULTS FROM THE SIMPLE DARK MATTER SEARCH. , 2005, , .		3
49	SIMPLE dark matter search results. <i>Physics Letters, Section B: Nuclear, Elementary Particle and High-Energy Physics</i> , 2005, 621, 233-238.	1.5	55
50	Water adsorption in disordered mesoporous silica (Vycor) at 300K and 650K: A Grand Canonical Monte Carlo simulation study of hysteresis. <i>Journal of Chemical Physics</i> , 2005, 122, 094704.	1.2	82
51	Phase coexistence in heterogeneous porous media: A new extension to Gibbs ensemble Monte Carlo simulation method. <i>Journal of Chemical Physics</i> , 2005, 122, 134710.	1.2	26
52	Capillary Condensation in a Geometrically and a Chemically Heterogeneous Pore: A Molecular Simulation Study. <i>Journal of Physical Chemistry B</i> , 2005, 109, 4700-4706.	1.2	24
53	Grand Potential, Helmholtz Free Energy, and Entropy Calculation in Heterogeneous Cylindrical Pores by the Grand Canonical Monte Carlo Simulation Method. <i>Journal of Physical Chemistry B</i> , 2005, 109, 480-487.	1.2	39
54	Thermodynamic Characterization of Fluids Confined in Heterogeneous Pores by Monte Carlo Simulations in the Grand Canonical and the Isobaric-Isenthalpic Ensembles. <i>Journal of Physical Chemistry B</i> , 2005, 109, 8185-8194.	1.2	20

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55	New extension of the Gibbs ensemble Monte Carlo simulation method: analysis of phase coexistence of simple fluids adsorbed in heterogeneous pores. <i>European Journal of Control</i> , 2005, 30, 401-410.	1.6	0
56	A grand canonical Monte Carlo simulation study of water adsorption on Vycor-like hydrophilic mesoporous silica at different temperatures. <i>Journal of Physics Condensed Matter</i> , 2004, 16, S5329-S5343.	0.7	22
57	A comparison of water adsorption on ordered and disordered silica substrates. <i>Physical Chemistry Chemical Physics</i> , 2004, 6, 1933-1937.	1.3	33
58	Confinement effect on thermodynamic and structural properties of water in hydrophilic mesoporous silica. <i>European Physical Journal E</i> , 2003, 12, 67-70.	0.7	14
59	Grand canonical Monte Carlo simulation study of water structure on hydrophilic mesoporous and plane silica substrates. <i>Journal of Chemical Physics</i> , 2003, 119, 9226-9232.	1.2	65
60	Water adsorption on hydrophilic mesoporous and plane silica substrates: A grand canonical Monte Carlo simulation study. <i>Journal of Chemical Physics</i> , 2003, 118, 5613-5622.	1.2	97
61	A grand canonical Monte Carlo simulation study of water adsorption in a Vycor-like disordered mesoporous material at 300K.. <i>Studies in Surface Science and Catalysis</i> , 2002, , 371-378.	1.5	0
62	The Status of SIMPLE in 2002. , 2002, , 524-528.		0
63	The Status of SIMPLE-2000. , 2001, , 598-603.		1
64	WIMP SEARCHES WITH SUPERHEATED DROPLET DETECTORS: STATUS AND PROSPECTS. , 2001, ,		0
65	Prospects for SIMPLE 2000: a large-mass, low-background superheated droplet detector for WIMP searches. <i>New Journal of Physics</i> , 2000, 2, 14-14.	1.2	31
66	First Dark Matter Limits from a Large-Mass, Low-Background, Superheated Droplet Detector. <i>Physical Review Letters</i> , 2000, 85, 3083-3086.	2.9	49