

Oscar Pinto

List of Publications by Year in descending order

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28
papers

251
citations

1040056

9
h-index

996975

15
g-index

28
all docs

28
docs citations

28
times ranked

184
citing authors

#	ARTICLE	IF	CITATIONS
1	Theoretical approach to energy levels applied to modified surfaces. <i>Physical Chemistry Chemical Physics</i> , 2022, , .	2.8	0
2	Mean field approach applied to surface deposition on a modified electrode. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 20247-20254.	2.8	1
3	Addressing the surface coverage of Au nano-agglomerates and the electrochemical properties of modified carbon paste electrodes: Experimental and theoretical studies on ascorbic acid oxidation. <i>Colloids and Surfaces B: Biointerfaces</i> , 2021, 200, 111585.	5.0	2
4	Effect of Temperature on The Kinetics of Electrochemical Insertion of Li-Ions into a Graphite Electrode Studied by Kinetic Monte Carlo. <i>Journal of the Electrochemical Society</i> , 2020, 167, 013533.	2.9	25
5	Kinetic Monte Carlo applied to the electrochemical study of the Li-ion graphite system. <i>Electrochimica Acta</i> , 2020, 331, 135439.	5.2	23
6	Structural surface and thermodynamics analysis of nanoparticles with defects. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 23148-23157.	2.8	0
7	Fractional and integer stages of lithium ion-graphite systems: the role of electrostatic and elastic contributions. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 16174-16183.	2.8	5
8	Numerical simulations of cyclic voltammetry for lithium-ion intercalation in nanosized systems: finiteness of diffusion versus electrode kinetics. <i>Journal of Solid State Electrochemistry</i> , 2020, 24, 3279-3287.	2.5	16
9	Electrosorption of a repulsive binary mixture on modified electrodes. <i>Surface Science</i> , 2020, 695, 121587.	1.9	4
10	Interaction of semiochemicals with model lipid membranes: A biophysical approach. <i>Colloids and Surfaces B: Biointerfaces</i> , 2018, 161, 413-419.	5.0	3
11	Electrosorption of a modified electrode in the vicinity of phase transition: A Monte Carlo study. <i>Applied Surface Science</i> , 2018, 433, 705-712.	6.1	7
12	The kinetic origin of the Daumas-HÅ©rld model for the Li-ion/graphite intercalation system. <i>Electrochemistry Communications</i> , 2018, 93, 133-137.	4.7	39
13	Grand Canonical Monte Carlo Study of Li Intercalation into Graphite. <i>Journal of the Electrochemical Society</i> , 2018, 165, A2019-A2025.	2.9	16
14	Electrochemical behavior of a typical redox mediator on a modified electrode surface: Experiment and computer simulations. <i>Surface Science</i> , 2017, 658, 15-21.	1.9	6
15	Simulation of selective thermodynamic deposition in nanoholes. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 1601-1609.	2.8	2
16	Criticality of the phase transition on stage two in a lattice-gas model of a graphite anode in a lithium-ion battery. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 23138-23145.	2.8	13
17	Quasi-chemical approach for adsorption of mixtures with non-additive lateral interactions. <i>Applied Surface Science</i> , 2017, 392, 1088-1096.	6.1	6
18	Monomolecular adsorption on nanoparticles with repulsive interactions: a Monte Carlo study. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 14610-14618.	2.8	8

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19	Phase diagrams for the adsorption of monomers with non-additive interactions. <i>Surface Science</i> , 2016, 651, 62-69.	1.9	8
20	The adsorption of a mixture of particles with non-additive interactions: a Monte Carlo study. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 3050-3058.	2.8	8
21	Microthermodynamic Interpretation of Fluid States from FTIR Measurements in Lipid Membranes: A Monte Carlo Study. <i>Journal of Physical Chemistry B</i> , 2014, 118, 10436-10443.	2.6	11
22	Computer simulation and detailed mean-field approximation applied to adsorption on nanoparticles. <i>Physical Review E</i> , 2013, 88, 062407.	2.1	4
23	Configurational entropy of systems with non-additive lateral interactions. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2012, 391, 6390-6398.	2.6	1
24	Computer simulation of adsorption on nanoparticles: The case of attractive interactions. <i>Physical Review E</i> , 2012, 86, 061602.	2.1	8
25	Lattice-gas model of nonadditive interacting particles on nanotube bundles. <i>Journal of Chemical Physics</i> , 2011, 134, 064702.	3.0	8
26	Statistical thermodynamics of straight rigid rods with nonadditive lateral interactions: Theory and Monte Carlo simulations. <i>Physical Review E</i> , 2011, 84, 061142.	2.1	4
27	Adsorption thermodynamics of a lattice-gas model with non-additive lateral interactions on triangular and honeycomb lattices. <i>Physica A: Statistical Mechanics and Its Applications</i> , 2010, 389, 3456-3464.	2.6	9
28	Adsorption thermodynamics of a lattice-gas model with non-additive lateral interactions. <i>Surface Science</i> , 2008, 602, 1763-1769.	1.9	14