

Gennady Gor

List of Publications by Year in descending order

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37
papers

817
citations

567281

15
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501196

28
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38
all docs

38
docs citations

38
times ranked

701
citing authors

#	ARTICLE	IF	CITATIONS
1	Adsorption-induced deformation of nanoporous materials—A review. <i>Applied Physics Reviews</i> , 2017, 4, .	11.3	189
2	Elastic response of mesoporous silicon to capillary pressures in the pores. <i>Applied Physics Letters</i> , 2015, 106, .	3.3	64
3	Revisiting Bangham's law of adsorption-induced deformation: changes of surface energy and surface stress. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 9788-9798.	2.8	50
4	Deformation of Microporous Carbons during N ₂ , Ar, and CO ₂ Adsorption: Insight from the Density Functional Theory. <i>Langmuir</i> , 2016, 32, 8265-8274.	3.5	49
5	Adsorption-Induced Deformation of Hierarchically Structured Mesoporous Silica—Effect of Pore-Level Anisotropy. <i>Langmuir</i> , 2017, 33, 5592-5602.	3.5	47
6	Relation between pore size and the compressibility of a confined fluid. <i>Journal of Chemical Physics</i> , 2015, 143, 194506.	3.0	38
7	Adsorption Stress Changes the Elasticity of Liquid Argon Confined in a Nanopore. <i>Langmuir</i> , 2014, 30, 13564-13569.	3.5	35
8	Sorption-Induced Deformation and Elastic Weakening of Bentheim Sandstone. <i>Journal of Geophysical Research: Solid Earth</i> , 2018, 123, 8589-8601.	3.4	29
9	Gassmann Theory Applies to Nanoporous Media. <i>Geophysical Research Letters</i> , 2018, 45, 146-155.	4.0	27
10	Modulus—pressure equation for confined fluids. <i>Journal of Chemical Physics</i> , 2016, 145, 164505.	3.0	26
11	Effect of pore geometry on the compressibility of a confined simple fluid. <i>Journal of Chemical Physics</i> , 2018, 148, 054503.	3.0	23
12	Single Parameter for Predicting the Morphology of Atmospheric Black Carbon. <i>Environmental Science & Technology</i> , 2018, 52, 14169-14179.	10.0	19
13	Adsorption-Induced Surface Stresses of the Water/Quartz Interface: Ab Initio Molecular Dynamics Study. <i>Langmuir</i> , 2016, 32, 5259-5266.	3.5	18
14	Force Fields for Molecular Modeling of Sarin and its Simulants: DMMP and DIMP. <i>Journal of Physical Chemistry B</i> , 2021, 125, 4086-4098.	2.6	18
15	Elastocapillarity in nanopores: Sorption strain from the actions of surface tension and surface stress. <i>Physical Review Materials</i> , 2018, 2, .	2.4	18
16	Porous Structure of Silica Colloidal Crystals. <i>Langmuir</i> , 2019, 35, 2230-2235.	3.5	15
17	Elastic properties of confined fluids from molecular modeling to ultrasonic experiments on porous solids. <i>Applied Physics Reviews</i> , 2021, 8, .	11.3	14
18	Pore-Size Distribution of Silica Colloidal Crystals from Nitrogen Adsorption Isotherms. <i>Langmuir</i> , 2019, 35, 14975-14982.	3.5	13

#	ARTICLE	IF	CITATIONS
19	Molecular Simulations Shed Light on Potential Uses of Ultrasound in Nitrogen Adsorption Experiments. <i>Langmuir</i> , 2018, 34, 15650-15657.	3.5	12
20	Mechanical Characterization of Hierarchical Structured Porous Silica by in Situ Dilatometry Measurements during Gas Adsorption. <i>Langmuir</i> , 2019, 35, 2948-2956.	3.5	12
21	In Situ Small-Angle Neutron Scattering Investigation of Adsorption-Induced Deformation in Silica with Hierarchical Porosity. <i>Langmuir</i> , 2019, 35, 11590-11600.	3.5	11
22	Non-covalent interactions of nitrous oxide with aromatic compounds: Spectroscopic and computational evidence for the formation of 1:1 complexes. <i>Journal of Chemical Physics</i> , 2014, 140, 144304.	3.0	10
23	Density Functional Theory Model for Adsorption-Induced Deformation of Mesoporous Materials with Nonconvex Pore Geometry. <i>Journal of Physical Chemistry C</i> , 2020, 124, 20046-20054.	3.1	10
24	Compressibility of Supercritical Methane in Nanopores: A Molecular Simulation Study. <i>Energy & Fuels</i> , 2020, 34, 1506-1513.	5.1	10
25	Adsorption-induced deformation of mesoporous materials with corrugated cylindrical pores. <i>Journal of Chemical Physics</i> , 2020, 153, 194703.	3.0	10
26	Models of adsorption-induced deformation: ordered materials and beyond. <i>Journal of Physics Condensed Matter</i> , 2022, 34, 063002.	1.8	9
27	Solvation pressure in spherical mesopores: Macroscopic theory and molecular simulations. <i>AIChE Journal</i> , 2021, 67, e16542.	3.6	7
28	Modeling elastic properties of Vycor glass saturated with liquid and solid adsorbates. <i>Adsorption</i> , 2019, 25, 973-982.	3.0	5
29	Adsorption-induced deformation of hierarchical organised carbon materials with ordered, non-convex mesoporosity. <i>Molecular Physics</i> , 2021, 119, .	1.7	5
30	Bulk Modulus of Not-So-Bulk Fluid. , 2017, , .		4
31	Capillary bridge formation between hexagonally ordered carbon nanorods. <i>Adsorption</i> , 2020, 26, 563-578.	3.0	4
32	Kinetic model for competitive condensation of vapor between concave and convex surfaces in a soot aggregate. <i>Aerosol Science and Technology</i> , 2021, 55, 302-315.	3.1	4
33	Compressibility of a Simple Fluid in Cylindrical Confinement: Molecular Simulation and Equation of State Modeling. <i>Industrial & Engineering Chemistry Research</i> , 2020, 59, 8393-8402.	3.7	3
34	The effect of interconnections on gas adsorption in materials with spherical mesopores: A Monte Carlo simulation study. <i>Journal of Chemical Physics</i> , 2021, 154, 114706.	3.0	3
35	Molecular Simulations of Vapor-Liquid Equilibrium of Isocyanates. <i>Journal of Physical Chemistry B</i> , 2021, 125, 12528-12538.	2.6	2
36	Adsorption from binary liquid solutions into mesoporous silica: a capacitance isotherm on 5CB nematogen/cyclohexane mixtures. <i>Molecular Physics</i> , 2021, 119, .	1.7	1

#	ARTICLE	IF	CITATIONS
37	Molecular Simulation of Benzene Adsorption in Graphitic and Amorphous Carbon Slit Pores. Journal of Chemical & Engineering Data, 2022, 67, 1765-1778.	1.9	1