

Elena Butyrskaya

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9142505/publications.pdf>

Version: 2024-02-01

20
papers

81
citations

1684188

5
h-index

1588992

8
g-index

20
all docs

20
docs citations

20
times ranked

46
citing authors

#	ARTICLE	IF	CITATIONS
1	Mechanistic aspects of functional layer formation in hybrid one-step designed GOx/Nafion/Pd-NPs nanobiosensors. <i>Analyst</i> , The, 2021, 146, 2172-2185.	3.5	5
2	Cluster model of the step-shaped adsorption isotherm in metal-organic frameworks. <i>Microporous and Mesoporous Materials</i> , 2021, 322, 111146.	4.4	4
3	Sorption Interaction between Carbon Nanotubes and Histidine Enantiomers in Aqueous Solutions. <i>Russian Journal of Physical Chemistry A</i> , 2021, 95, 2280-2286.	0.6	1
4	Mechanism of Sorption Interaction between L-Alanine and Carbon Nanotubes. <i>Russian Journal of Physical Chemistry A</i> , 2019, 93, 710-716.	0.6	1
5	Separation of ethylene glycol and alkali metal salts on carbon nanotubes and mosaic membranes. <i>Zavodskaya Laboratoriya Diagnostika Materialov</i> , 2018, 84, 11-17.	0.5	0
6	Sorption interactions between ethylene glycol and carbon nanotubes. <i>Russian Journal of Physical Chemistry A</i> , 2017, 91, 567-571.	0.6	3
7	Computer simulation of amino acid sorption on carbon nanotubes. <i>Journal of Structural Chemistry</i> , 2017, 58, 217-225.	1.0	6
8	Sorption Interactions between L-Alanine and Carbon Nanotubes in Aqueous Solutions. <i>Journal of Physical Chemistry C</i> , 2017, 121, 20524-20531.	3.1	9
9	Effect of the calculation method and the basis set on the structure and electrical properties of (4,4) carbon nanotubes with different lengths and open ends. <i>Journal of Structural Chemistry</i> , 2016, 57, 649-657.	1.0	1
10	Computer simulation of size effects and adsorption properties of one-wall carbon nanotubes (6,6). <i>Russian Journal of General Chemistry</i> , 2016, 86, 1684-1691.	0.8	4
11	Determining role of hydrogen bonding in electrically driven membrane transport: Quantum-chemical and molecular dynamics study. <i>Petroleum Chemistry</i> , 2015, 55, 918-926.	1.4	1
12	Extraction of ethylene glycol from aqueous salt solutions. <i>Russian Journal of Physical Chemistry A</i> , 2012, 86, 1721-1725.	0.6	3
13	Standardless structural-group analysis of supramolecular systems. <i>Journal of Analytical Chemistry</i> , 2009, 64, 1000-1006.	0.9	4
14	Computer simulation of the infrared spectra of endohedral metallofullerenes Li ₂ C ₆₀ and Na ₂ C ₆₀ . <i>Physics of the Solid State</i> , 2009, 51, 649-656.	0.6	4
15	Transfer of glycols through an MA-41 ion-exchange membrane from aqueous and salt-aqueous solutions. <i>Russian Journal of Physical Chemistry A</i> , 2007, 81, 405-409.	0.6	1
16	Separation of mineral salts and nonelectrolytes (ethylene glycol) by dialysis through ion-exchange membranes. <i>Journal of Analytical Chemistry</i> , 2007, 62, 710-715.	0.9	8
17	Interpretation of hypsochromic and bathochromic shifts of vibrational frequencies of a cation exchanger. <i>Journal of Analytical Chemistry</i> , 2007, 62, 930-934.	0.9	5
18	Computer Simulation of Cation-Exchange Membrane Structure: An Elementary Act of Hydrated Ion Transport. <i>Russian Journal of Electrochemistry</i> , 2004, 40, 767-770.	0.9	10

#	ARTICLE	IF	CITATIONS
19	Ab Initio Calculation of the Structure and Functions of Sulfo Cation Exchangers. Journal of Structural Chemistry, 2003, 44, 1062-1066.	1.0	3
20	Interpretation of infrared spectra for ion-exchange systems. Optics and Spectroscopy (English) Tj ETQq0 0 0 rgBT /Overlock 10 Tf 50 70	0.6	8