CITATION REPORT List of articles citing



DOI: 10.1021/acs.jchemed.7b00385 Journal of Chemical Education, 2018, 95, 888-894.

Source: https://exaly.com/paper-pdf/68895007/citation-report.pdf

Version: 2024-04-10

This report has been generated based on the citations recorded by exaly.com for the above article. For the latest version of this publication list, visit the link given above.

The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

#	Paper	IF	Citations
14	Realistic Implementation of the Particle Model for the Visualization of Nanoparticle Precipitation and Growth. <i>Journal of Chemical Education</i> , 2019 , 96, 1654-1662	2.4	4
13	Dynamical aspects of supercooled TIP3P-water in the grooves of DNA. <i>Journal of Chemical Physics</i> , 2019 , 150, 235101	3.9	4
12	Process Oriented Guided Inquiry Learning Computational Chemistry Experiments: Revisions and Extensions Based on Lessons Learned from Implementation. <i>ACS Symposium Series</i> , 2019 , 65-77	0.4	1
11	Discovery-Based Computational Activities in the Undergraduate Chemistry Curriculum. <i>ACS Symposium Series</i> , 2019 , 227-243	0.4	2
10	FatsILoveHate Relationships: A Molecular Dynamics Simulation and Hands-On Experiment Outreach Activity to Introduce the Amphiphilic Nature and Biological Functions of Lipids to Young Students and the General Public. <i>Journal of Chemical Education</i> , 2020 , 97, 1360-1367	2.4	2
9	Molecular dynamics simulations and analysis for bioinformatics undergraduate students. <i>Biochemistry and Molecular Biology Education</i> , 2021 , 49, 570-582	1.3	4
8	Transforming traditional teaching laboratories for effective remote delivery Areview. <i>Education for Chemical Engineers</i> , 2021 , 35, 96-104	2.4	13
7	Systematic Comparison of the Structural and Dynamic Properties of Commonly Used Water Models for Molecular Dynamics Simulations. <i>Journal of Chemical Information and Modeling</i> , 2021 , 61, 4521-4536	6.1	14
6	Calculation of diffusion coefficients of pesticides by employing molecular dynamics simulations. <i>Journal of Molecular Liquids</i> , 2021 , 340, 117106	6	1
5	Development of an electro-pneumatic system for the practical training of pneumatic processes in the university environment. <i>Journal of Physics: Conference Series</i> , 2021 , 2073, 012016	0.3	1
4	Describing the adsorption of sodium tripolyphosphate on kaolinite surfaces in a saline medium by molecular dynamics. <i>Minerals Engineering</i> , 2022 , 175, 107280	4.9	O
3	Mechanism of the distinct toxicity level of imidacloprid and thiacloprid against honey bees: An in silico study based on cytochrome P450 9Q3. <i>Journal of Molecular Graphics and Modelling</i> , 2022 , 116, 108	3 2 57	
2	The Performance of Different Water Models on the Structure and Function of Cytochrome P450 Enzymes.		2
1	Experimental and Simulation Studies on Hematite Interaction with Na-Metasilicate Pentahydrate. 2023 , 28, 3629		0