Xinwen Ou

List of Publications by Year in descending order

Source: https://exaly.com/author-pdf/810938/publications.pdf

Version: 2024-02-01

		1478505	1125743	
13	176	6	13	
papers	citations	h-index	g-index	
13	13	13	181	
all docs	docs citations	times ranked	citing authors	

#	Article	IF	CITATIONS
1	Spontaneous desorption of protein from self-assembled monolayer (SAM)-coated gold nanoparticles induced by high temperature. Physical Chemistry Chemical Physics, 2022, 24, 2363-2370.	2.8	2
2	Distinct lipid membrane interaction and uptake of differentially charged nanoplastics in bacteria. Journal of Nanobiotechnology, 2022, 20, 191.	9.1	30
3	Mechanistic Insight on General Protein-Binding Ability of ATP and the Impacts of Arginine Residues. Journal of Physical Chemistry B, 2022, 126, 4647-4658.	2.6	5
4	Membrane Insertion of MoS2 Nanosheets: Fresh vs. Aged. Frontiers in Chemistry, 2021, 9, 706917.	3.6	6
5	ATP Can Efficiently Stabilize Protein through a Unique Mechanism. Jacs Au, 2021, 1, 1766-1777.	7.9	24
6	Structure and Mechanical Stabilities of the Three-Way Junction Motifs in Prohead RNA. Journal of Physical Chemistry B, 2021, 125, 12125-12134.	2.6	2
7	Spontaneous Translocation of Single-Stranded DNA in Graphene–MoS ₂ Heterostructure Nanopores: Shape Effect. Journal of Physical Chemistry B, 2020, 124, 9490-9496.	2.6	12
8	Dynamic behaviors of interfacial water on the self-assembly monolayer (SAM) heterogeneous surface. Journal of Chemical Physics, 2020, 153, 124705.	3.0	3
9	Structure and sequence features of mussel adhesive protein lead to its salt-tolerant adhesion ability. Science Advances, 2020, 6, .	10.3	47
10	Effects of the surface polarity of nanomaterials on their interaction with complement protein gC1q. RSC Advances, 2020, 10, 41993-42000.	3.6	2
11	Surface microstructure engenders unusual hydrophobicity in phyllosilicates. Chemical Communications, 2018, 54, 5418-5421.	4.1	17
12	Unintentionally doped hydrogen removal mechanism in Li doped ZnO. AIP Advances, 2018, 8, .	1.3	6
13	Heterogeneous Condensation of Water on the Mica (001) Surface: A Molecular Dynamics Simulation Work. Journal of Physical Chemistry C, 2017, 121, 6813-6819.	3.1	20