

Andre Muniz

List of Publications by Citations

Source: <https://exaly.com/author-pdf/6598927/andre-muniz-publications-by-citations.pdf>

Version: 2024-04-24

This document has been generated based on the publications and citations recorded by exaly.com. 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.

47
papers

750
citations

17
h-index

26
g-index

52
ext. papers

924
ext. citations

5
avg, IF

4.86
L-index

#	Paper	IF	Citations
47	Ab initio studies of thermodynamic and electronic properties of phosphorene nanoribbons. <i>Physical Review B</i> , 2014 , 90,	3.3	112
46	Opening and tuning of band gap by the formation of diamond superlattices in twisted bilayer graphene. <i>Physical Review B</i> , 2012 , 86,	3.3	46
45	Mechanical behavior of interlayer-bonded nanostructures obtained from bilayer graphene. <i>Carbon</i> , 2015 , 81, 663-677	10.4	45
44	Functionalized diamond nanothreads from benzene derivatives. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 7132-7137	3.6	42
43	First-principles calculation of the mechanical properties of diamond nanothreads. <i>Carbon</i> , 2017 , 113, 260-265	10.4	40
42	Superlattices of Fluorinated Interlayer-Bonded Domains in Twisted Bilayer Graphene. <i>Journal of Physical Chemistry C</i> , 2013 , 117, 7315-7325	3.8	34
41	Elastic properties of graphene nanomeshes. <i>Applied Physics Letters</i> , 2014 , 104, 141911	3.4	33
40	Tunable mechanical properties of diamond superlattices generated by interlayer bonding in twisted bilayer graphene. <i>Applied Physics Letters</i> , 2013 , 103, 013113	3.4	30
39	An open and extensible sigma-profile database for COSMO-based models. <i>AIChE Journal</i> , 2018 , 64, 3443-3455	3.6	28
38	Formation of fullerene superlattices by interlayer bonding in twisted bilayer graphene. <i>Journal of Applied Physics</i> , 2012 , 111, 043513	2.5	28
37	Experimental and computational analysis of carbon molecular sieve membrane formation upon polyetherimide pyrolysis. <i>Carbon</i> , 2017 , 119, 21-29	10.4	24
36	Elucidation of deactivation phenomena in cobalt catalyst for Fischer-Tropsch synthesis using SSITKA. <i>Journal of Catalysis</i> , 2016 , 344, 669-679	7.3	23
35	Mechanical behavior and fracture of graphene nanomeshes. <i>Journal of Applied Physics</i> , 2015 , 117, 024302	2.5	23
34	Formation and Mechanical Behavior of Nanocomposite Superstructures from Interlayer Bonding in Twisted Bilayer Graphene. <i>ACS Applied Materials & Interfaces</i> , 2018 , 10, 28898-28908	9.5	20
33	Effects of hydrogen chemisorption on the structure and deformation of single-walled carbon nanotubes. <i>Applied Physics Letters</i> , 2009 , 94, 103108	3.4	20
32	Electronic and Mechanical Properties of Partially Saturated Carbon and Carbon Nitride Nanothreads. <i>Journal of Physical Chemistry C</i> , 2019 , 123, 3886-3891	3.8	18
31	Carbon nanothreads from polycyclic aromatic hydrocarbon molecules. <i>Carbon</i> , 2018 , 140, 644-652	10.4	18

30	On the hydrogen storage capacity of carbon nanotube bundles. <i>Applied Physics Letters</i> , 2009 , 95, 163111	3.4	17
29	Analysis of diamond nanocrystal formation from multiwalled carbon nanotubes. <i>Physical Review B</i> , 2009 , 80,	3.3	17
28	Carbon-Based Nanostructures Derived from Bilayer Graphene with Zero Thermal Expansion Behavior. <i>Journal of Physical Chemistry C</i> , 2015 , 119, 17458-17465	3.8	13
27	Diamond nanothread-based 2D and 3D materials: Diamond nanomeshes and nanofoams. <i>Carbon</i> , 2018 , 139, 789-800	10.4	13
26	Hydrogenation effects on the structure and morphology of graphene and single-walled carbon nanotubes. <i>Journal of Applied Physics</i> , 2010 , 108, 113532	2.5	11
25	Mechanical properties of hydrogenated electron-irradiated graphene. <i>Journal of Applied Physics</i> , 2016 , 120, 124301	2.5	11
24	Fluorescence Based Platform to Discriminate Protein Using Carbon Quantum Dots. <i>ChemistrySelect</i> , 2019 , 4, 5619-5627	1.8	8
23	EVALUATION OF BIOMASS AND COAL CO-GASIFICATION OF BRAZILIAN FEEDSTOCK USING A CHEMICAL EQUILIBRIUM MODEL. <i>Brazilian Journal of Chemical Engineering</i> , 2016 , 33, 401-414	1.7	8
22	Nanoporous carbon superstructures based on covalent bonding of porous fullerenes. <i>Carbon</i> , 2018 , 130, 424-432	10.4	7
21	First-principles study of carbon nanothreads derived from five-membered heterocyclic rings: thiophene, furan and pyrrole. <i>Physical Chemistry Chemical Physics</i> , 2021 , 23, 2055-2062	3.6	7
20	Structure-properties relations in graphene derivatives and metamaterials obtained by atomic-scale modeling. <i>Molecular Simulation</i> , 2019 , 45, 1173-1202	2	5
19	Thermal annealing of graphite oxide under high pressure: An experimental and computational study. <i>Carbon</i> , 2018 , 139, 1035-1047	10.4	5
18	High-order finite volume method for solving viscoelastic fluid flows. <i>Brazilian Journal of Chemical Engineering</i> , 2008 , 25, 153-166	1.7	5
17	Prediction of β -P and β P: Two New Strain-Interconvertible Phosphorene Allotropes. <i>Journal of Physical Chemistry C</i> , 2020 , 124, 21207-21214	3.8	5
16	Tuning the band structure of graphene nanoribbons through defect-interaction-driven edge patterning. <i>Physical Review B</i> , 2017 , 96,	3.3	4
15	Molecular dynamics study of the effects of static and oscillating electric fields in ovalbumin. <i>Innovative Food Science and Emerging Technologies</i> , 2022 , 75, 102911	6.8	4
14	Strain-Tunable Carbon Nanothread-Derived Membranes for Water Desalination. <i>Journal of Physical Chemistry B</i> , 2021 , 125, 7311-7319	3.4	4
13	Chain- and chainmail-like nanostructures from carbon nanotube rings. <i>Computational Materials Science</i> , 2019 , 161, 76-82	3.2	3

12	Molecular-Dynamics Analysis of Nanoindentation of Graphene Nanomeshes: Implications for 2D Mechanical Metamaterials. <i>ACS Applied Nano Materials</i> , 2020 , 3, 3613-3624	5.6	3
11	Flexible carbon nanothread-based membranes with strain-dependent gas transport properties. <i>Journal of Membrane Science</i> , 2019 , 585, 184-190	9.6	2
10	Unveiling the Origin of the Giant Barocaloric Effect in Natural Rubber. <i>Macromolecules</i> , 2020 , 53, 2606-2615	9.5	2
9	Reduction-Driven 3D to 2D Transformation of Cu Nanoparticles.. <i>Small</i> , 2022 , e2106583	11	2
8	Análise teórico-experimental da dispersão de poluentes líquidos em solos. <i>Engenharia Sanitaria E Ambiental</i> , 2007 , 12, 410-416	0.4	2
7	Peering into the Formation of Template-Free Hierarchical Flowerlike Nanostructures of SrTiO. <i>ACS Omega</i> , 2020 , 5, 33007-33016	3.9	2
6	Molecular-Dynamics Simulations on Nanoindentation of Graphene-Diamond Composite Superstructures in Interlayer-Bonded Twisted Bilayer Graphene: Implications for Mechanical Metamaterials. <i>ACS Applied Nano Materials</i> , 2021 , 4, 8611-8625	5.6	2
5	Nature of the interactions between Fe and Zr for the methane dehydroaromatization reaction in ZSM-5. <i>Journal of Molecular Structure</i> , 2020 , 1220, 128720	3.4	1
4	Uma nova metodologia para a simulação de escoamentos de fluidos viscoelásticos. <i>Polimeros</i> , 2005 , 15, 53-58	1.6	1
3	Experimental and DFT analysis of the acid and reduction properties of Fe-Cu/ZSM-5. <i>Microporous and Mesoporous Materials</i> , 2021 , 314, 110860	5.3	1
2	The effect of subgrid-scale modeling on LES of turbulent coaxial jets. <i>Journal of the Brazilian Society of Mechanical Sciences and Engineering</i> , 2021 , 43, 1	2	0
1	Reduction-Driven 3D to 2D Transformation of Cu Nanoparticles (Small 7/2022). <i>Small</i> , 2022 , 18, 2270032	11	1