

Hakim Amara

List of Publications by Year in descending order

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62
papers

2,458
citations

218381

26
h-index

197535

49
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63
all docs

63
docs citations

63
times ranked

2868
citing authors

#	ARTICLE	IF	CITATIONS
1	Effect of size on the surface energy of noble metal nanoparticles from analytical and numerical approaches. <i>Physical Review B</i> , 2022, 105, .	1.1	10
2	Simulation of thermodynamic properties of magnetic transition metals from an efficient tight-binding model: The case of cobalt and beyond. <i>Physical Review B</i> , 2022, 105, .	1.1	1
3	Colloidal synthesis of nanoparticles: from bimetallic to high entropy alloys. <i>Nanoscale</i> , 2022, 14, 9832-9841.	2.8	13
4	Size-dependent hydrogen trapping in palladium nanoparticles. <i>Journal of Materials Chemistry A</i> , 2021, 9, 10354-10363.	5.2	15
5	Quantitative In Situ Visualization of Thermal Effects on the Formation of Gold Nanocrystals in Solution. <i>Advanced Materials</i> , 2021, 33, e2102514.	11.1	15
6	Quantitative Study of Temperature Effects on The Nucleation and Growth of Gold Nanocrystals in Water. <i>Microscopy and Microanalysis</i> , 2021, 27, 29-30.	0.2	0
7	Morphology control of metallic nanoparticles supported on carbon substrates in catalytic conditions. <i>Carbon</i> , 2020, 159, 504-511.	5.4	9
8	Strain and electronic properties at the van der Waals interface of phosphorus/boron nitride heterobilayers. <i>Physical Review B</i> , 2020, 102, .	1.1	1
9	Selective shortening of gold nanorods: when surface functionalization dictates the reactivity of nanostructures. <i>Nanoscale</i> , 2020, 12, 22658-22667.	2.8	13
10	A deep learning approach for determining the chiral indices of carbon nanotubes from high-resolution transmission electron microscopy images. <i>Carbon</i> , 2020, 169, 465-474.	5.4	27
11	Revealing the Dynamics of Functional Nanomaterials in Their Formation and Application Media with Liquid and Gas-phase TEM. <i>Microscopy and Microanalysis</i> , 2020, 26, 196-198.	0.2	1
12	High density synthesis of topological point defects in graphene on 6H α -SiC(0001 \bar{A}). <i>Carbon</i> , 2020, 170, 174-181.	5.4	9
13	Quantitative insights into the growth mechanisms of nanopores in hexagonal boron nitride. <i>Physical Review Materials</i> , 2020, 4, .	0.9	8
14	Tuning bimetallic catalysts for a selective growth of SWCNTs. <i>Nanoscale</i> , 2019, 11, 4091-4100.	2.8	16
15	Revealing the Surface Energetics and Reactivity of Bimetallic Copper-Gold Catalyst Nanoparticles by In Situ Environmental TEM. <i>Microscopy and Microanalysis</i> , 2019, 25, 33-34.	0.2	1
16	Exciton interference in hexagonal boron nitride. <i>Physical Review B</i> , 2018, 97, .	1.1	23
17	Direct Measurement of the Surface Energy of Bimetallic Nanoparticles: Evidence of Vegard \hat{A} 's Rule-like Dependence. <i>Physical Review Letters</i> , 2018, 120, 025901.	2.9	19
18	Growth modes and chiral selectivity of single-walled carbon nanotubes. <i>Nanoscale</i> , 2018, 10, 6744-6750.	2.8	67

#	ARTICLE	IF	CITATIONS
19	Direct and indirect excitons in boron nitride polymorphs: A story of atomic configuration and electronic correlation. <i>Physical Review B</i> , 2018, 98, .	1.1	63
20	Entropy-driven stability of chiral single-walled carbon nanotubes. <i>Science</i> , 2018, 362, 212-215.	6.0	75
21	Two-photon absorption in two-dimensional materials: The case of hexagonal boron nitride. <i>Physical Review B</i> , 2018, 98, .	1.1	22
22	Ni ₂ C surface carbide to catalyze low-temperature graphene growth. <i>Physical Review B</i> , 2018, 97, .	1.1	2
23	Excitons in few-layer hexagonal boron nitride: Davydov splitting and surface localization. <i>2D Materials</i> , 2018, 5, 045017.	2.0	63
24	Probing the role of carbon solubility in transition metal catalyzing single-walled carbon nanotubes growth. <i>Carbon</i> , 2017, 120, 226-232.	5.4	37
25	Modeling the Growth of Single-Wall Carbon Nanotubes. <i>Topics in Current Chemistry</i> , 2017, 375, 55.	3.0	26
26	Structural Properties of Double-Walled Carbon Nanotubes Driven by Mechanical Interlayer Coupling. <i>ACS Nano</i> , 2017, 11, 4840-4847.	7.3	21
27	Linking growth mode to lengths of single-walled carbon nanotubes. <i>Carbon</i> , 2017, 113, 231-236.	5.4	75
28	Magnetism as indirect tool for carbon content assessment in nickel nanoparticles. <i>Journal of Applied Physics</i> , 2017, 122, 213902.	1.1	2
29	Excitons in boron nitride single layer. <i>Physical Review B</i> , 2016, 94, .	1.1	68
30	Size Dependent Phase Diagrams of Nickel-Carbon Nanoparticles. <i>Physical Review Letters</i> , 2015, 115, 205502.	2.9	54
31	Charge transfer and electronic doping in nitrogen-doped graphene. <i>Scientific Reports</i> , 2015, 5, 14564.	1.6	79
32	Electronic Interaction between Nitrogen Atoms in Doped Graphene. <i>ACS Nano</i> , 2015, 9, 670-678.	7.3	69
33	Key roles of carbon solubility in single-walled carbon nanotube nucleation and growth. <i>Nanoscale</i> , 2015, 7, 20284-20289.	2.8	27
34	Interdependency of Subsurface Carbon Distribution and Grapheneâ€Catalyst Interaction. <i>Journal of the American Chemical Society</i> , 2014, 136, 13698-13708.	6.6	95
35	Magnetism: the driving force of order in CoPt, a first-principles study. <i>Journal of Physics Condensed Matter</i> , 2013, 25, 056005.	0.7	26
36	Random vs realistic amorphous carbon models for high resolution microscopy and electron diffraction. <i>Journal of Applied Physics</i> , 2013, 114, .	1.1	18

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37	Atomistic modelling of CVD synthesis of carbon nanotubes and graphene. <i>Nanoscale</i> , 2013, 5, 6662.	2.8	88
38	Computational studies of catalyst-free single walled carbon nanotube growth. <i>Journal of Chemical Physics</i> , 2013, 139, 054308.	1.2	1
39	Role of defect healing on the chirality of single-wall carbon nanotubes. <i>Physical Review B</i> , 2012, 85, .	1.1	18
40	Carbon solubility in nickel nanoparticles: A grand canonical Monte Carlo study. <i>Physica Status Solidi (B): Basic Research</i> , 2012, 249, 2629-2634.	0.7	26
41	Computational studies of graphene growth mechanisms. <i>Physical Review B</i> , 2012, 85, .	1.1	20
42	Long-range interactions between substitutional nitrogen dopants in graphene: Electronic properties calculations. <i>Physical Review B</i> , 2012, 86, .	1.1	91
43	Importance of Carbon Solubility and Wetting Properties of Nickel Nanoparticles for Single Wall Nanotube Growth. <i>Physical Review Letters</i> , 2012, 109, 185501.	2.9	85
44	Evidence of Correlation between Catalyst Particles and the Single-Wall Carbon Nanotube Diameter: A First Step towards Chirality Control. <i>Physical Review Letters</i> , 2012, 108, 195503.	2.9	119
45	Imaging the symmetry breaking of molecular orbitals in single-wall carbon nanotubes. <i>Physical Review B</i> , 2010, 81, .	1.1	8
46	Nickel-Assisted Healing of Defective Graphene. <i>ACS Nano</i> , 2010, 4, 6114-6120.	7.3	79
47	Aluminum and vacancies in α -iron: Dissolution, diffusion, and clustering. <i>Physical Review B</i> , 2010, 81, .	1.1	53
48	Density functional theory and tight binding-based dynamical studies of carbon metal systems of relevance to carbon nanotube growth. <i>Nano Research</i> , 2009, 2, 774-782.	5.8	7
49	Spin-Unrestricted Calculations of Bare-Edged Nanographenes Using DFT and Many-Body Perturbation Theory. <i>Journal of Chemical Theory and Computation</i> , 2009, 5, 1719-1722.	2.3	12
50	Computational Studies of Metal-Carbon Nanotube Interfaces for Regrowth and Electronic Transport. <i>Nano Letters</i> , 2009, 9, 1117-1120.	4.5	31
51	Early Stages in the Nucleation Process of Carbon Nanotubes. <i>ACS Nano</i> , 2009, 3, 511-516.	7.3	75
52	Tight-binding potential for atomistic simulations of carbon interacting with transition metals: Application to the Ni-C system. <i>Physical Review B</i> , 2009, 79, .	1.1	109
53	Interaction of carbon clusters with Ni(100): Application to the nucleation of carbon nanotubes. <i>Surface Science</i> , 2008, 602, 77-83.	0.8	21
54	Understanding the Nucleation Mechanisms of Carbon Nanotubes in Catalytic Chemical Vapor Deposition. <i>Physical Review Letters</i> , 2008, 100, 056105.	2.9	141

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55	A Tight-Binding Grand Canonical Monte Carlo Study of the Catalytic Growth of Carbon Nanotubes. Journal of Nanoscience and Nanotechnology, 2008, 8, 6099-6104.	0.9	8
56	INVESTIGATION OF THE POLARIZABILITY AND OF THE LOCAL ELECTRIC FIELD OF FULLERENES WITH VARIOUS SHAPES AND DEFECTS USING A MONOPOLE-DIPOLE INTERACTION MODEL. , 2007, , .		0
57	Catalytically Assisted Tip Growth Mechanism for Single-Wall Carbon Nanotubes. ACS Nano, 2007, 1, 202-207.	7.3	44
58	Scanning tunneling microscopy fingerprints of point defects in graphene: A theoretical prediction. Physical Review B, 2007, 76, .	1.1	164
59	Study of the polarizability of fullerenes with a monopoleâ€“dipole interaction model. Diamond and Related Materials, 2007, 16, 2145-2149.	1.8	19
60	Characterization of single-walled carbon nanotubes containing defects from their local vibrational densities of states. Carbon, 2007, 45, 349-356.	5.4	24
61	Formation of carbon nanostructures on nickel surfaces: A tight-binding grand canonical Monte Carlo study. Physical Review B, 2006, 73, .	1.1	73
62	Nucleation and Growth of Single-Walled Nanotubes: The Role of Metallic Catalysts. Journal of Nanoscience and Nanotechnology, 2004, 4, 346-359.	0.9	71