José Eduardo Padilha

List of Publications by Year in descending order

Source: https://exaly.com/author-pdf/4454209/publications.pdf

Version: 2024-02-01

32 papers

1,388 citations

471509 17 h-index 30 g-index

32 all docs 32 docs citations

times ranked

32

2174 citing authors

#	Article	IF	CITATIONS
1	Oxidation of two-dimensional electrides: Structural transition and the formation of half-metallic channels protected by oxide layers. Physical Review B, 2022, 105, .	3.2	1
2	Green synthesis of templated carbon porous materials from simple raw materials. Materials Advances, 2021, 2, 403-412.	5.4	7
3	Structural Transition in Oxidized Ca2N Electrenes: CaO/CaN 2D Heterostructures. Journal of Physical Chemistry C, 2020, 124, 14706-14712.	3.1	4
4	Interatomic potential for atomistic simulation of self-catalyzed GaAs nanowires growth. Computational Materials Science, 2020, 183, 109805.	3.0	5
5	Metal Chalcogenides Janus Monolayers for Efficient Hydrogen Generation by Photocatalytic Water Splitting. ACS Applied Nano Materials, 2019, 2, 890-897.	5.0	93
6	Structural evolution and the role of native defects in subnanometer MoS nanowires. Physical Review B, 2019, 100 , .	3.2	7
7	Electronic and optical properties of hydrogenated group-IV multilayer materials. Physical Chemistry Chemical Physics, 2018, 20, 8112-8118.	2.8	12
8	Tunable magnetism and spin-polarized electronic transport in graphene mediated by molecular functionalization of extended defects. Physical Review B, $2018, 97, .$	3.2	9
9	Microscopic Description of the Ferroism in Lead-Free AlFeO3. Scientific Reports, 2018, 8, 6420.	3.3	17
10	Layer-dependent band alignment of few layers of blue phosphorus and their van der Waals heterostructures with graphene. Physical Review B, 2018, 97, .	3.2	45
11	Stacking-dependent transport properties in few-layers graphene. Solid State Communications, 2017, 250, 70-74.	1.9	10
12	Nanodots of transition metal dichalcogenides embedded in MoS2 and MoSe2: first-principles calculations. Physical Chemistry Chemical Physics, 2017, 19, 26240-26247.	2.8	0
13	Two-dimensional van der Waals <i>p-n</i> junction of InSe/phosphorene. Physical Review B, 2017, 95, .	3.2	68
14	Directional dependence of the electronic and transport properties of 2D borophene and borophane. Physical Chemistry Chemical Physics, 2016, 18, 25491-25496.	2.8	92
15	Fully and partially iodinated germanane as a platform for the observation of the quantum spin Hall effect. Physical Review B, 2016, 93, .	3.2	18
16	Electron density distribution and electronic structure as tools to study the origin of ferroic states in ferroelectric and magnetic materials. Ferroelectrics, 2016, 500, 26-36.	0.6	2
17	Substrate-supported large-band-gap quantum spin Hall insulator based on III-V bismuth layers. Physical Review B, 2016, 94, .	3.2	4
18	A new class of large band gap quantum spin hall insulators: 2D fluorinated group-IV binary compounds. Scientific Reports, 2016, 6, 26123.	3.3	17

#	Article	IF	Citations
19	Electronic and transport properties of structural defects in monolayer germanene: An ab initio investigation. Solid State Communications, 2016, 225, 38-43.	1.9	50
20	Topological phase transitions of (BixSb1 \hat{a}° x)2Se3 alloys by density functional theory. Journal of Physics Condensed Matter, 2015, 27, 255501.	1.8	10
21	van der Waals Heterostructure of Phosphorene and Graphene: Tuning the Schottky Barrier and Doping by Electrostatic Gating. Physical Review Letters, 2015, 114, 066803.	7.8	445
22	Free-Standing Bilayer Silicene: The Effect of Stacking Order on the Structural, Electronic, and Transport Properties. Journal of Physical Chemistry C, 2015, 119, 3818-3825.	3.1	73
23	Nature and evolution of the band-edge states in <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML"><mml:msub><mml:mi>MoS</mml:mi><mml:mn>2<td>:m8.2<td>าl:m<u>s</u>ab></td></td></mml:mn></mml:msub></mml:math>	:m 8.2 <td>าl:m<u>s</u>ab></td>	า l:m<u>s</u>a b>
24	Directional Control of the Electronic and Transport Properties of Graphynes. Journal of Physical Chemistry C, 2014, 118, 18793-18798.	3.1	18
25	Quantum spin Hall effect on germanene nanorod embedded in completely hydrogenated germanene. Physical Review B, 2014, 89, .	3.2	57
26	Graphene nanoribbon intercalated with hexagonal boron nitride: Electronic transport properties from ab initio calculations. Solid State Communications, 2013, 173, 24-29.	1.9	10
27	xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"> <mml:msub><mml:mrow /><mml:mi>x</mml:mi></mml:mrow </mml:msub> Ge <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline"><mml:msub><mml:mrow /><mml:mrow><mml:mn>1</mml:mn><mml:mo>a^3</mml:mo><mml:mi>x</mml:mi></mml:mrow><td>3.2 > <td>24 ath valloy</td></td></mml:mrow </mml:msub></mml:math 	3.2 > <td>24 ath valloy</td>	24 ath valloy
28	Physical Review B, 2013, 88. Bilayer graphene on h-BN substrate: investigating the breakdown voltage and tuning the bandgap by electric field. Journal of Physics Condensed Matter, 2012, 24, 075301.	1.8	22
29	Bilayer graphene dual-gate nanodevice: An <i>ab initio</i> simulation. Physical Review B, 2011, 84, .	3.2	36
30	IxV curves of boron and nitrogen doping zigzag graphene nanoribbons. International Journal of Quantum Chemistry, 2011, 111, 1379-1386.	2.0	17
31	Energetics and stability of vacancies in carbon nanotubes. Solid State Communications, 2011, 151, 482-486.	1.9	42
32	Transport properties of single vacancies in nanotubes. Physical Review B, 2008, 77, .	3.2	35