

Eduardo Costa Girão

List of Publications by Year in descending order

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49
papers

843
citations

567144

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501076

28
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50
all docs

50
docs citations

50
times ranked

1091
citing authors

#	ARTICLE	IF	CITATIONS
1	Electronic properties of boron-rich graphene nanowiggles. Computational Materials Science, 2022, 201, 110907.	1.4	1
2	Structural and electronic properties of nonconventional \hat{I}^{\pm} -graphyne nanocarbons. Physical Review Materials, 2022, 6, .	0.9	2
3	Electronic properties of carbon sheets and nanoribbons based on acepentallene-like building blocks. Computational Materials Science, 2022, 211, 111520.	1.4	0
4	Electronic and magnetic properties of tripentaphene nanoribbons. Physical Review Materials, 2022, 6, .	0.9	2
5	Phenyl- and naphthyl-type heteroatom substitution blocks in naphthylene- $\langle \text{mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" display="inline" id="d1e883" altimg="si39.svg" \rangle \langle \text{mml:mi} \rangle \hat{I}^3 \langle \text{mml:mi} \rangle \langle \text{mml:math} \rangle$: A DFT study. Computational Materials Science, 2022, 213, 111578.	1.4	0
6	Computational study of elastic, structural stability and dynamics properties of penta-graphene membrane. Chemical Physics, 2021, 542, 111052.	0.9	16
7	Electronic properties of N-rich graphene nano-chevron. Physical Chemistry Chemical Physics, 2021, 23, 13204-13215.	1.3	6
8	Electronic and Transport Properties of Graphene Nanoribbons Based on Super-Heptazethrene Molecular Blocks. Journal of Physical Chemistry C, 2021, 125, 11235-11248.	1.5	7
9	Mechanical properties of single-walled penta-graphene-based nanotubes: A DFT and Classical molecular dynamics study. Chemical Physics, 2021, 547, 111187.	0.9	6
10	Structural and electronic properties of double-walled \hat{I}^{\pm} -graphyne nanotubes. Computational Materials Science, 2021, 200, 110768.	1.4	1
11	Electronic properties of 2D and 1D carbon allotropes based on a triphenylene structural unit. Physical Chemistry Chemical Physics, 2021, 23, 25114-25125.	1.3	7
12	Tripentaphenes: two-dimensional acepentallene-based nanocarbon allotropes. Physical Chemistry Chemical Physics, 2020, 22, 23195-23206.	1.3	10
13	Electronic and structural properties of tetragraphenes. Carbon, 2020, 167, 403-413.	5.4	11
14	Naphthylene- $\langle \text{mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" \rangle \langle \text{mml:mi} \rangle \hat{I}^3 \langle \text{mml:mi} \rangle \langle \text{mml:math} \rangle$: 1D and 2D carbon allotropes based on the fusion of phenyl- and naphthyl-like groups. Physical Review Materials, 2020, 4, .	0.9	4
15	Elastic properties of graphyne-based nanotubes. Computational Materials Science, 2019, 170, 109153.	1.4	25
16	Naphthylenes: 1D and 2D carbon allotropes based on naphthyl units. Carbon, 2019, 153, 792-803.	5.4	23
17	Modeling the Kondo effect of a magnetic atom adsorbed on graphene. 2D Materials, 2019, 6, 035038.	2.0	3
18	Structural and electronic properties of nanotubes constructed from fragmented fullerenes. Carbon, 2019, 147, 616-627.	5.4	10

#	ARTICLE	IF	CITATIONS
19	Electronic properties of tetragraphene nanoribbons. <i>Physical Review Materials</i> , 2019, 3, .	0.9	14
20	Mechanical Properties of Pentagraphene-based Nanotubes: A Molecular Dynamics Study. <i>MRS Advances</i> , 2018, 3, 97-102.	0.5	10
21	Mechanical Properties of Phagraphene Membranes: A Fully Atomistic Molecular Dynamics Investigation. <i>MRS Advances</i> , 2018, 3, 67-72.	0.5	6
22	High efficiency spin-valve and spin-filter in a doped rhombic graphene quantum dot device. <i>Journal of Magnetism and Magnetic Materials</i> , 2018, 451, 532-539.	1.0	8
23	Spin-Negative Differential Resistance in Zigzag Graphene Nanoribbons with Side-Attached Porphine Molecule. <i>Journal of Physical Chemistry C</i> , 2018, 122, 15911-15921.	1.5	14
24	Quantum Dots in Graphene Nanoribbons. <i>Nano Letters</i> , 2017, 17, 4277-4283.	4.5	99
25	Tuning the electronic and quantum transport properties of nitrogenated holey graphene nanoribbons. <i>Journal of Materials Chemistry C</i> , 2017, 5, 11856-11866.	2.7	13
26	One- and two-dimensional carbon nanostructures based on unfolded buckyballs: An <i>ab initio</i> investigation of their electronic properties. <i>Physical Review B</i> , 2017, 95, .	1.1	13
27	Electronic, transport, and magnetic properties of punctured carbon nanotubes. <i>Physical Review B</i> , 2016, 94, .	1.1	3
28	Improved All-Carbon Spintronic Device Design. <i>Scientific Reports</i> , 2015, 5, 7634.	1.6	52
29	Heterospin Junctions in Zigzag-Edged Graphene Nanoribbons. <i>Applied Sciences (Switzerland)</i> , 2014, 4, 351-365.	1.3	1
30	Electronic transport in three-terminal triangular carbon nanopatches. <i>Nanotechnology</i> , 2014, 25, 045706.	1.3	3
31	Quantifying energetics of topological frustration in carbon nanostructures. <i>Physical Review B</i> , 2014, 89, .	1.1	9
32	Electronic properties of three-terminal graphitic nanowiggles. <i>Physical Review B</i> , 2014, 90, .	1.1	4
33	Electronic and magnetic structures of coronene-based graphitic nanoribbons. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 3603.	1.3	10
34	Emergent magnetism in irradiated graphene nanostructures. <i>Carbon</i> , 2014, 78, 196-203.	5.4	9
35	Quasiparticle band gaps of graphene nanowiggles and their magnetism on Au(111). <i>Physical Review B</i> , 2013, 88, .	1.1	15
36	Patchwork algorithm for the parallel computation of the Green's function in open systems. <i>Journal of Computational Electronics</i> , 2013, 12, 123-133.	1.3	17

#	ARTICLE	IF	CITATIONS
37	Electronic and transport properties of graphene nanoribbon barbell-shaped heterojunctions. <i>Physica Status Solidi (B): Basic Research</i> , 2013, 250, 2417-2423.	0.7	3
38	Modeling and Simulation of Electron Transport at the Nanoscale: Illustrations in Low-Dimensional Carbon Nanostructures. <i>Advances in Atom and Single Molecule Machines</i> , 2013, , 123-133.	0.0	0
39	Enhanced thermoelectric figure of merit in assembled graphene nanoribbons. <i>Physical Review B</i> , 2012, 86, .	1.1	81
40	Electronic Transport Properties of Assembled Carbon Nanoribbons. <i>ACS Nano</i> , 2012, 6, 6483-6491.	7.3	29
41	Topographic and Spectroscopic Characterization of Electronic Edge States in CVD Grown Graphene Nanoribbons. <i>Nano Letters</i> , 2012, 12, 1928-1933.	4.5	104
42	Structural and electronic properties of graphitic nanowiggles. <i>Physical Review B</i> , 2012, 85, .	1.1	24
43	Electronic transport properties of carbon nanotoroids. <i>Nanotechnology</i> , 2011, 22, 075701.	1.3	6
44	Emergence of Atypical Properties in Assembled Graphene Nanoribbons. <i>Physical Review Letters</i> , 2011, 107, 135501.	2.9	69
45	Electronic transmission selectivity in multiterminal graphitic nanorings. <i>Applied Physics Letters</i> , 2011, 98, 112111.	1.5	5
46	Nicotine adsorption on single wall carbon nanotubes. <i>Journal of Hazardous Materials</i> , 2010, 184, 678-683.	6.5	19
47	Functionalization of single-wall carbon nanotubes through chloroform adsorption: theory and experiment. <i>Physical Chemistry Chemical Physics</i> , 2010, 12, 1518.	1.3	27
48	A single molecule rectifier with strong push-pull coupling. <i>Journal of Chemical Physics</i> , 2008, 129, 204701.	1.2	17
49	First principles study of 1,2-dichlorobenzene adsorption on metallic carbon nanotubes. <i>International Journal of Quantum Chemistry</i> , 2006, 106, 2558-2563.	1.0	23