

Neeraj K Jaiswal

List of Publications by Year in descending order

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81
papers

840
citations

471509

17
h-index

552781

26
g-index

81
all docs

81
docs citations

81
times ranked

427
citing authors

#	ARTICLE	IF	CITATIONS
1	Label-free biosensor using nanogap embedded dielectric modulated schottky tunneling source impact ionization MOS. <i>Microelectronic Engineering</i> , 2016, 149, 129-134.	2.4	54
2	Fe-Doped Armchair Graphene Nanoribbons for Spintronic/Interconnect Applications. <i>IEEE Nanotechnology Magazine</i> , 2013, 12, 685-691.	2.0	51
3	Adsorption of CO gas molecules on zigzag BN/AlN nanoribbons for nano sensor applications. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2019, 383, 532-538.	2.1	43
4	Adsorption of COCl ₂ gas molecule on armchair boron nitride nanoribbons for nano sensor applications. <i>Microelectronic Engineering</i> , 2015, 146, 62-67.	2.4	39
5	First principles calculations of cobalt doped zigzag graphene nanoribbons. <i>Solid State Communications</i> , 2012, 152, 1489-1492.	1.9	37
6	First-principles investigation of armchair boron nitride nanoribbons for sensing PH ₃ gas molecules. <i>Superlattices and Microstructures</i> , 2014, 73, 350-358.	3.1	31
7	Inducing half-metallicity with enhanced stability in zigzag graphene nanoribbons via fluorine passivation. <i>Applied Surface Science</i> , 2017, 396, 471-479.	6.1	31
8	Structural stability and electronic properties of Ni-doped armchair graphene nanoribbons. <i>Solid State Communications</i> , 2011, 151, 1490-1495.	1.9	29
9	Realizing Negative Differential Resistance/Switching Phenomena in Zigzag GaN Nanoribbons by Edge Fluorination: A DFT Investigation. <i>Advanced Materials Interfaces</i> , 2017, 4, 1700400.	3.7	29
10	First principles calculations of armchair graphene nanoribbons interacting with Cu atoms. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2011, 44, 75-79.	2.7	28
11	First-Principles Investigation of Pd-Doped Armchair Graphene Nanoribbons as a Potential Rectifier. <i>Journal of Electronic Materials</i> , 2021, 50, 1196-1206.	2.2	27
12	Reconstructed graphene nanoribbon as a sensor for nitrogen based molecules. <i>Applied Surface Science</i> , 2015, 357, 55-59.	6.1	24
13	Edge-Oxidized Germanene Nanoribbons for Nanoscale Metal Interconnect Applications. <i>IEEE Transactions on Electron Devices</i> , 2018, 65, 3893-3900.	3.0	23
14	Chlorine sensing properties of zigzag boron nitride nanoribbons. <i>Solid State Communications</i> , 2014, 185, 41-46.	1.9	21
15	Enhanced metallicity and spin polarization in zigzag graphene nanoribbons with Fe impurities. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2013, 54, 103-108.	2.7	20
16	Possibility of spin-polarized transport in edge fluorinated armchair boron nitride nanoribbons. <i>RSC Advances</i> , 2016, 6, 11014-11022.	3.6	17
17	Tailoring the electronic properties of zigzag graphene nanoribbons via sp ² /sp ³ edge functionalization with H/F. <i>Organic Electronics</i> , 2017, 51, 25-37.	2.6	17
18	Structural and electronic properties of armchair graphene nanoribbons functionalized with fluorine. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2019, 383, 125949.	2.1	17

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19	First-principle study of NO ₂ adsorption and detection on the edges of zigzag nitride nanoribbons. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2019, 114, 113575.	2.7	17
20	Tailoring the electronic properties of a Z-shaped graphene field effect transistor via B/N doping. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2015, 379, 710-718.	2.1	16
21	Half-metallicity in armchair boron nitride nanoribbons: A first-principles study. <i>Solid State Communications</i> , 2015, 212, 19-24.	1.9	16
22	First-principles study of sensing SO ₂ adsorption on III-V nitride nanoribbons. <i>Materials Chemistry and Physics</i> , 2020, 242, 122437.	4.0	16
23	First-principle investigations of cove edged GaN nanoribbon for nanoscale resonant tunneling applications. <i>Solid State Communications</i> , 2021, 340, 114486.	1.9	16
24	Prospects of asymmetrically H-terminated zigzag germanene nanoribbons for spintronic application. <i>Applied Surface Science</i> , 2017, 396, 1352-1359.	6.1	15
25	Structural & electronic properties of zigzag silicene nanoribbons with symmetric/asymmetric edge passivations via fluorine and hydrogen. <i>Physica B: Condensed Matter</i> , 2021, 615, 413072.	2.7	15
26	Electronic and Transport Properties of Zigzag Boron Nitride Nanoribbons. <i>Journal of Computational and Theoretical Nanoscience</i> , 2013, 10, 368-375.	0.4	13
27	First-principles investigation of CO ₂ and NH ₃ adsorption on antimonene nanoribbons. <i>Materials Today: Proceedings</i> , 2020, 28, 65-69.	1.8	10
28	First-principle investigations of negative differential resistance in zigzag boron nitride nanoribbons. <i>Physica E: Low-Dimensional Systems and Nanostructures</i> , 2021, 134, 114861.	2.7	10
29	Selective Edge-Hydrogenated Zigzag Boron Nitride Nanoribbons for Giant Magnetoresistance and Rectifying Behavior. <i>IEEE Transactions on Electron Devices</i> , 2021, 68, 5894-5900.	3.0	9
30	Analytical model for 2DEG charge density in $\text{In}_{1-x}\text{Ga}_x\text{As}$ nanowires. <i>IEEE Transactions on Electron Devices</i> , 2022, 37, 025002.	2.0	9
31	Structural, electronic, and magnetic properties of Mn-Doped InP nanowire. <i>Superlattices and Microstructures</i> , 2016, 92, 134-142.	3.1	8
32	First-principle insights of CO and NO detection via antimonene nanoribbons. <i>Applied Physics A: Materials Science and Processing</i> , 2020, 126, 1.	2.3	8
33	First-Principles Investigations of N-Vacancy Induced Zigzag Boron Nitride Nanoribbons for Nanoscale Resonant Tunneling Applications. <i>Journal of Electronic Materials</i> , 2021, 50, 5664-5681.	2.2	8
34	Giant Magnetoresistance and Rectification Behavior in Fluorinated Zigzag Boron Nitride Nanoribbon for Spintronic Nanodevices. <i>IEEE Nanotechnology Magazine</i> , 2022, 21, 244-250.	2.0	8
35	Density functional insights of iodine interaction with graphene and its nanoribbon with zigzag edges. <i>Physica B: Condensed Matter</i> , 2018, 545, 268-274.	2.7	7
36	Sarin chemisorbent based on cobalt-doped graphene. <i>Applied Surface Science</i> , 2019, 480, 759-764.	6.1	7

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37	First-principle investigations for electronic transport in nitrogen-doped disconnected zigzag graphene nanoribbons. <i>Microelectronic Engineering</i> , 2018, 199, 96-100.	2.4	6
38	Zn-Passivated Zigzag Boron Nitride Nanoribbons for Perfect Spin-Filtering and Negative Differential Resistance Based Devices. <i>IEEE Nanotechnology Magazine</i> , 2022, 21, 299-306.	2.0	6
39	Role of Device Parameters in Optimizing 2DEG Charge Density in \hat{I}^2 -(Al _x) _{1-x} GaN Heterostructure. Analytical Approach. <i>IEEE Transactions on Electron Devices</i> , 2022, 69, 3876-3883.	3.0	6
40	Potential spin-polarized transport in gold-doped armchair graphene nanoribbons. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2015, 379, 835-842.	2.1	5
41	First-principles study of electronic properties of GaN nanowires: Effect of surface orientation, passivation, and Mn doping. <i>Physica Status Solidi (B): Basic Research</i> , 2016, 253, 2185-2196.	1.5	5
42	Intrinsic half metallicity in lithium terminated zigzag graphene nanoribbons. <i>Solid State Communications</i> , 2017, 250, 112-118.	1.9	5
43	First-principles design of nano-porous graphene membranes for efficient separation of halogen gases. <i>Diamond and Related Materials</i> , 2020, 108, 107911.	3.9	5
44	First-principle investigations of zigzag III-V nitride nanoribbons as CS ₂ scavengers. <i>Applied Surface Science</i> , 2021, 545, 148969.	6.1	5
45	Ab-Initio Study of Transition Metal (Ni) Interaction with Zigzag Graphene Nanoribbons. <i>Journal of Computational and Theoretical Nanoscience</i> , 2012, 9, 555-559.	0.4	4
46	Tailoring the Electronic Structure of Zigzag Graphene Nanoribbons via Cu Impurities. <i>Journal of Computational and Theoretical Nanoscience</i> , 2013, 10, 1441-1445.	0.4	4
47	Enhancing the performance of BN nanosheets as promising anode material for Li-ion batteries with carbon-doping. <i>Journal of Molecular Graphics and Modelling</i> , 2022, 115, 108213.	2.4	4
48	First-principles study of CO adsorption on zigzag ZnO nanoribbons towards nanosensor application. <i>Journal of Molecular Graphics and Modelling</i> , 2022, 116, 108232.	2.4	4
49	First-principles predictions of tunable half metallicity in zigzag GaN nanoribbons with possible applications in CO detection and spintronics. <i>Journal Physics D: Applied Physics</i> , 2022, 55, 405002.	2.8	4
50	COCl ₂ Gas Adsorption on Zigzag Boron Nitride Nanoribbons: First Principles Study. <i>Journal of Computational and Theoretical Nanoscience</i> , 2015, 12, 2472-2476.	0.4	3
51	Electronic properties and migration pathways of Cl functionalized zigzag graphene nanoribbons. <i>AIP Conference Proceedings</i> , 2017, , .	0.4	3
52	First principle insights of NO ₂ detection via III-V nitride nanoribbons with armchair edges. <i>Nano Express</i> , 2020, 1, 010059.	2.4	3
53	First-principles Investigation of Antimonene Nanoribbons for Sensing Toxic NO ₂ Gas. <i>Physica Status Solidi (B): Basic Research</i> , 2020, 257, 2000034.	1.5	3
54	Tuning the Electronic Structure of Zigzag Boron Nitride Nanoribbons via sp ² /sp ³ Edge Functionalization. <i>Journal of Electronic Materials</i> , 2022, 51, 3299-3307.	2.2	3

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55	On the evolution and electronic properties of self-assembled gold nanowires. Computational Materials Science, 2017, 130, 222-231.	3.0	2
56	Structural, magnetic and electronic properties of armchair graphene nanoribbons interacting with Co: DFT investigations. Ferroelectrics, 2017, 519, 178-186.	0.6	2
57	Effect of Positional Cu Doping on the Electronic and Transport Properties of Graphene Nanoribbons. Graphene, 2014, 2, 40-51.	0.2	2
58	Comparative Analysis of T-Gate and L-Gate Dielectric Modulated Schottky Tunneling Source Impact Ionization MOS for Label-Free Detection of Toxic Gases. Journal of Nanoelectronics and Optoelectronics, 2018, 13, 501-508.	0.5	2
59	Enhancing metallicity in zigzag graphene nanoribbons with adsorbed/substitutionally doped copper atoms. , 2012, , .		1
60	Transport properties of Ni-interacted armchair graphene nanoribbons. , 2013, , .		1
61	Ab initio study of gold-doped zigzag graphene nanoribbons. Applied Physics A: Materials Science and Processing, 2014, 117, 1997-2008.	2.3	1
62	Electronic properties of armchair graphene nanoribbons doped with Cobalt atoms. IOP Conference Series: Materials Science and Engineering, 2015, 73, 012140.	0.6	1
63	First principles study of Li terminated armchair graphene nanoribbons. Materials Today: Proceedings, 2016, 3, 2241-2246.	1.8	1
64	First-principles study for I-V characteristics of halogen functionalized zigzag graphene nanoribbons. AIP Conference Proceedings, 2019, , .	0.4	1
65	First-principle investigations of Cl decorated armchair GaN nanoribbons. AIP Conference Proceedings, 2020, , .	0.4	1
66	First-principles study of structural and electronic properties of cove-edged zigzag ZnO nanoribbons. Materials Today: Proceedings, 2022, 67, 616-619.	1.8	1
67	First principles study of adsorbed and substitutionally doped Fe atoms in zigzag graphene nanoribbons. , 2011, , .		0
68	First-Principles Study of Cu Doped Zigzag Graphene Nanoribbons. , 2011, , .		0
69	First Principle Calculations of Absorption of Cu-Doped Boron Nanotube. , 2011, , .		0
70	Ab-initio study of co-doped zigzag graphene nanoribbons. , 2012, , .		0
71	Transport properties of cu-doped armchair graphene nanoribbons in two probe geometry. , 2013, , .		0
72	Stability analysis of zigzag boron nitride nanoribbons. AIP Conference Proceedings, 2015, , .	0.4	0

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73	Spintronic and transport properties of linear atomic strings of transition metals (Fe, Co, Ni). AIP Conference Proceedings, 2016, , .	0.4	0
74	Shape dependent electronic properties of wurzite GaN nanowire. AIP Conference Proceedings, 2016, , .	0.4	0
75	Functionalization and migration of bromine adatoms on zigzag graphene nanoribbons: A first-principles study. AIP Conference Proceedings, 2018, , .	0.4	0
76	First-principle study of zigzag GaN nanoribbons with Cl functionalization. AIP Conference Proceedings, 2019, , .	0.4	0
77	Density functional investigations of Li adatom migration on perfect and vacancy engineered BN nanosheets. AIP Conference Proceedings, 2020, , .	0.4	0
78	First-Principle Study of Cl Functionalized Zigzag AlN Nanoribbons. Integrated Ferroelectrics, 2020, 205, 114-121.	0.7	0
79	First-Principle Study of Cl Functionalized Armchair AlN Nanoribbons. Springer Proceedings in Mathematics and Statistics, 2021, , 151-157.	0.2	0
80	Half Metallic Transition In Silver-adsorbed Zigzag Graphene Nanoribbons. Advanced Materials Letters, 2016, 7, 708-712.	0.6	0
81	Effect Of Phases On The Energetics Of Pristine InP Nanowires: An Ab-initio Approach. Advanced Materials Letters, 2016, 7, 831-835.	0.6	0