Niraj K Nepal

List of Publications by Year
in descending order

Source: https://exaly.com/author-pdf/5880105/publications.pdf

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14	244	933447 10	996975 15
papers	citations	h-index	g-index

15	15	15	220
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	First-principles wave-vector- and frequency-dependent exchange-correlation kernel for jellium at all densities. Physical Review B, 2022, 105, .	3.2	7
2	Bending as a control knob for the electronic and optical properties of phosphorene nanoribbons. Physical Review Materials, 2022, 6, .	2.4	4
3	Tunable band gaps and optical absorption properties of bent MoS2 nanoribbons. Scientific Reports, 2022, 12, 3008.	3.3	14
4	Interpretations of ground-state symmetry breaking and strong correlation in wavefunction and density functional theories. Proceedings of the National Academy of Sciences of the United States of America, 2021, 118, .	7.1	53
5	Describing adsorption of benzene, thiophene, and xenon on coinage metals by using the Zarembaâ€“Kohn theory-based model. Journal of Chemical Physics, 2021, 154, 124705.	3.0	4
6	Opening band gaps of low-dimensional materials at the meta-GGA level of density functional approximations. Physical Review Materials, 2021, 5, .	2.4	18
7	The Fermiâ€“LÃ¶wdin self-interaction correction for ionization energies of organic molecules. Journal of Chemical Physics, 2020, 153, 184303.	3.0	12
8	Formation energy puzzle in intermetallic alloys: Random phase approximation fails to predict accurate formation energies. Physical Review B, 2020, 102, .	3.2	12
9	Understanding plasmon dispersion in nearly free electron metals: Relevance of exact constraints for exchange-correlation kernels within time-dependent density functional theory. Physical Review B, 2020, 101, .	3.2	3
10	Constraint-based wave vector and frequency dependent exchange-correlation kernel of the uniform electron gas. Physical Review B, 2020, 101, .	3.2	17
11	Stretched or noded orbital densities and self-interaction correction in density functional theory. Journal of Chemical Physics, 2019, 150, 174102.	3.0	46
12	Treating different bonding situations: Revisiting Au-Cu alloys using the random phase approximation. Physical Review B, 2019, 100, .	3.2	10
13	First-principles study of mechanical and electronic properties of bent monolayer transition metal dichalcogenides. Physical Review Materials, 2019, 3, .	2.4	28
14	Rocksalt or cesium chloride: Investigating the relative stability of the cesium halide structures with random phase approximation based methods. Physical Review B, 2018, 97, .	3.2	13