

# Guodong Yu

## List of Publications by Year in descending order

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papers

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623734  
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docs citations

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times ranked

1092  
citing authors

#	ARTICLE	IF	CITATIONS
1	Interlayer hybridization in graphene quasicrystal and other bilayer graphene systems. Physical Review B, 2022, 105, .	3.2	4
2	Polarization-Dependent Selection Rules and Optical Spectrum Atlas of Twisted Bilayer Graphene Quantum Dots. Physical Review X, 2022, 12, .	8.9	8
3	Discovery of multivalley Fermi surface responsible for the high thermoelectric performance in Yb $\text{MnSb}_{11}$ and Yb $\text{MgSb}_{11}$ . Science Advances, 2021, 7, .	10.3	34
4	Structureâ€“Compositionâ€“Property Relationships in Antiperovskite Nitrides: Guiding a Rational Alloy Design. ACS Applied Materials & Interfaces, 2021, 13, 48516-48524.	8.0	14
5	Electronic properties and quasiparticle model of monolayer $\text{MoSi}_4$ . Physical Review B, 2021, 104, .	3.2	17
6	Electronic structure of $\text{N}_{30}$ twisted double bilayer graphene. Physical Review B, 2020, 102, .		
7	Tunability of multiple ultraflat bands and effect of spin-orbit coupling in twisted bilayer transition metal dichalcogenides. Physical Review B, 2020, 102, .	3.2	31
8	Pressure and electric field dependence of quasicrystalline electronic states in $\text{N}_{30}$ twisted bilayer graphene. Physical Review B, 2020, 102, .		
9	Interplay between in-plane and flexural phonons in electronic transport of two-dimensional semiconductors. Physical Review B, 2019, 100, .	3.2	11
10	Origins of ultralow thermal conductivity in 1-2-1-4 quaternary selenides. Journal of Materials Chemistry A, 2019, 7, 2589-2596.	10.3	28
11	Dodecagonal bilayer graphene quasicrystal and its approximants. Npj Computational Materials, 2019, 5, .	8.7	53
12	Computationally driven high-throughput identification of CaTe and $\text{Li}_3\text{P}$ as promising candidates for high-mobility transparent conducting materials. Physical Review Materials, 2019, 3, .	2.4	16
13	Resonant Bonding, Multiband Thermoelectric Transport, and Native Defects in n-Type $\text{BaBiTe}_3\text{xSex}$ ( $x = \frac{1}{7}$ )	6.7	14
14	PyCDT: A Python toolkit for modeling point defects in semiconductors and insulators. Computer Physics Communications, 2018, 226, 165-179.	7.5	142
15	Electronic and mechanical properties of few-layer borophene. Physical Review B, 2018, 98, .	3.2	83
16	Tunable half-metallicity and edge magnetism of H-saturated InSe nanoribbons. Physical Review Materials, 2018, 2, .	2.4	11
17	A computational assessment of the electronic, thermoelectric, and defect properties of bournonite ( $\text{CuPbSb}_3$ ) and related substitutions. Physical Chemistry Chemical Physics, 2017, 19, 6743-6756.	2.8	47
18	Two-dimensional Kagome phosphorus and its edge magnetism: a density functional theory study. Journal of Physics Condensed Matter, 2015, 27, 255006.	1.8	16

#	ARTICLE		IF	CITATIONS
19	Surface magnetism of the carbon foam: An ab-initio theoretical study. <i>Applied Physics Letters</i> , 2014, 105, 061601.		3.3	7
20	First-principles study of 3d transition metal atom adsorption onto graphene: the role of the extended line defect. <i>Journal of Materials Chemistry C</i> , 2014, 2, 9767-9774.		5.5	18
21	Structural, electronic and magnetic properties of transition-metal embedded zigzag-edged graphene nanoribbons. <i>Journal Physics D: Applied Physics</i> , 2013, 46, 375303.		2.8	9
22	Electronic properties of four typical zigzag-edged graphyne nanoribbons. <i>Journal of Physics Condensed Matter</i> , 2013, 25, 285502.		1.8	23
23	A valley-filtering switch based on the Stone-Wales defect array in carbon nanotube. <i>Europhysics Letters</i> , 2013, 103, 47008.		2.0	1
24	Line-defect-induced Fano interference in an armchair graphene nanoribbon. <i>Europhysics Letters</i> , 2013, 103, 18003.		2.0	7
25	RKKY interaction in AB-stacked multilayer graphene. <i>Journal of Physics Condensed Matter</i> , 2012, 24, 206003.		1.8	17
26	A simple tight-binding model for typical graphyne structures. <i>New Journal of Physics</i> , 2012, 14, 113007.		2.9	41