

# Xianqing Lin

## List of Publications by Year in descending order

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29  
papers

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566801

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docs citations

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times ranked

1478  
citing authors

#	ARTICLE	IF	CITATIONS
1	High pressure induced secondary and tertiary gaps in relaxed graphene on hexagonal boron nitride. Physical Review B, 2022, 105, .	1.1	2
2	Misalignment instability in magic-angle twisted bilayer graphene on hexagonal boron nitride. 2D Materials, 2021, 8, 025025.	2.0	11
3	Monte Carlo simulation study of the influence of defects on two-dimensional ferromagnetic order. AIP Advances, 2021, 11, 085016.	0.6	0
4	Emergence of intrinsically isolated flat bands and their topology in fully relaxed twisted multilayer graphene. Physical Review B, 2021, 104, .	1.1	4
5	Symmetry breaking in the double moiré superlattices of relaxed twisted bilayer graphene on hexagonal boron nitride. Physical Review B, 2020, 102, .	1.1	17
6	Band Engineering of Large-Twist-Angle Graphene/hBN Moiré Superlattices with Pressure. Physical Review Letters, 2020, 125, 226403.	1.7	17
7	Periodically Gated Bilayer Graphene as an Electronic Metamaterial. Physical Review Applied, 2020, 13, .	1.5	4
8	Pressure-induced gap modulation and topological transitions in twisted bilayer and twisted double bilayer graphene. Physical Review B, 2020, 101, .	1.1	19
9	Moiré effects in graphene-hBN heterostructures. Physical Review Research, 2020, 2, .	1.3	9
10	Layer-dependent intrinsic anomalous Hall effect in Fe <sub>3</sub> C <sub>2</sub> alloy films. Physical Review B, 2019, 100, .	1.3	10
11	Magnetostriction, Soft Magnetism, and Microwave Properties in Fe <sup>2+</sup> /Co Alloy Films. Physical Review Applied, 2019, 12, .	1.6	16
12	Effective lattice model of graphene moiré superlattices on hexagonal boron nitride. Physical Review B, 2019, 100, .	1.1	17
13	Shear instability in twisted bilayer graphene. Physical Review B, 2018, 98, .	1.1	31
14	Microscopic Mechanism of the Helix-to-Layer Transformation in Elemental Group VI Solids. Nano Letters, 2018, 18, 4908-4913.	4.5	19
15	Minimum model for the electronic structure of twisted bilayer graphene and related structures. Physical Review B, 2018, 98, .	1.1	34
16	Topologically insulating states in ternary transition metal dichalcogenides. Journal of Applied Physics, 2017, 121, 024303.	1.1	5
17	Switch effect of the nonquantized intrinsic spin Hall conductivity in monolayered monoclinic transition metal dichalcogenides. Journal of Physics Condensed Matter, 2017, 29, 295302.	0.7	1
18	Topological phase transition due to strain-controlled evolution of the inverted bands in 1T <sub>2</sub> C <sub>2</sub> . Physical Review B, 2017, 95, .	1.6	16

#	ARTICLE	IF	CITATIONS
19	Magnetism and electronic phase transitions in monoclinic transition metal dichalcogenides with transition metal atoms embedded. <i>Journal of Applied Physics</i> , 2016, 120, 064305.	1.1	10
20	Insulator-metal transition in 1T-MoS <sub>2</sub> under uniaxial strain. <i>Physics Letters, Section A: General, Atomic and Solid State Physics</i> , 2015, 379, 2883-2889.	0.9	7
21	Dirac points and van Hove singularities of silicene under uniaxial strain. <i>Journal of Applied Physics</i> , 2015, 117, 164305.	1.1	14
22	Tailoring of the structural, energetic and electronic properties of silicene-based nanostructures. <i>Journal of Physics: Conference Series</i> , 2014, 491, 012005.	0.3	4
23	The extraordinary magnetoelectric response in silicene doped with Fe and Cr atoms. <i>Applied Physics Letters</i> , 2014, 105, .	1.5	18
24	Charge and magnetic states of Mn-, Fe-, and Co-doped monolayer MoS <sub>2</sub> . <i>Journal of Applied Physics</i> , 2014, 116, .	1.1	92
25	Adsorption capacity of H <sub>2</sub> O, NH <sub>3</sub> , CO, and NO <sub>2</sub> on the pristine graphene. <i>Journal of Applied Physics</i> , 2013, 113, .	1.1	80
26	Flat bands near Fermi level of topological line defects on graphite. <i>Applied Physics Letters</i> , 2012, 101, .	1.5	30
27	Electronic and magnetic properties of substitutionally Fe-, Co-, and Ni-doped BC <sub>3</sub> honeycomb structure. <i>Journal of Applied Physics</i> , 2012, 111, .	1.1	12
28	Much stronger binding of metal adatoms to silicene than to graphene: A first-principles study. <i>Physical Review B</i> , 2012, 86, .	1.1	208
29	Half-metallicity in graphene nanoribbons with topological line defects. <i>Physical Review B</i> , 2011, 84, .	1.1	108