

# Geun Ho Gu

## List of Publications by Year in descending order

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Version: 2024-02-01

25  
papers

1,241  
citations

471509

17  
h-index

552781

26  
g-index

26  
all docs

26  
docs citations

26  
times ranked

1578  
citing authors

#	ARTICLE	IF	CITATIONS
1	Bimetallic Gold-Silver Nanostructures Drive Low Overpotentials for Electrochemical Carbon Dioxide Reduction. <i>ACS Applied Materials &amp; Interfaces</i> , 2022, 14, 6604-6614.	8.0	14
2	Perovskite synthesizability using graph neural networks. <i>Npj Computational Materials</i> , 2022, 8, .	8.7	16
3	Automated exploitation of the big configuration space of large adsorbates on transition metals reveals chemistry feasibility. <i>Nature Communications</i> , 2022, 13, 2087.	12.8	8
4	Predicting potentially hazardous chemical reactions using an explainable neural network. <i>Chemical Science</i> , 2021, 12, 11028-11037.	7.4	3
5	Autobifunctional Mechanism of Jagged Pt Nanowires for Hydrogen Evolution Kinetics via End-to-End Simulation. <i>Journal of the American Chemical Society</i> , 2021, 143, 5355-5363.	13.7	33
6	Understanding potential-dependent competition between electrocatalytic dinitrogen and proton reduction reactions. <i>Nature Communications</i> , 2021, 12, 4353.	12.8	78
7	Uncertainty Quantification and Error Propagation in the Enthalpy and Entropy of Surface Reactions Arising from a Single DFT Functional. <i>Journal of Physical Chemistry C</i> , 2021, 125, 18187-18196.	3.1	8
8	Reduced graphene oxides with engineered defects enable efficient electrochemical reduction of dinitrogen to ammonia in wide pH range. <i>Nano Energy</i> , 2020, 68, 104323.	16.0	64
9	Generative Adversarial Networks for Crystal Structure Prediction. <i>ACS Central Science</i> , 2020, 6, 1412-1420.	11.3	102
10	Structure-Based Synthesizability Prediction of Crystals Using Partially Supervised Learning. <i>Journal of the American Chemical Society</i> , 2020, 142, 18836-18843.	13.7	59
11	Highly stable two-dimensional bismuth metal-organic frameworks for efficient electrochemical reduction of CO <sub>2</sub> . <i>Applied Catalysis B: Environmental</i> , 2020, 277, 119241.	20.2	109
12	Progress in Computational and Machine Learning Methods for Heterogeneous Small-Molecule Activation. <i>Advanced Materials</i> , 2020, 32, e1907865.	21.0	46
13	Uncertainty-Quantified Hybrid Machine Learning/Density Functional Theory High Throughput Screening Method for Crystals. <i>Journal of Chemical Information and Modeling</i> , 2020, 60, 1996-2003.	5.4	31
14	Practical Deep-Learning Representation for Fast Heterogeneous Catalyst Screening. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 3185-3191.	4.6	63
15	Machine-enabled inverse design of inorganic solid materials: promises and challenges. <i>Chemical Science</i> , 2020, 11, 4871-4881.	7.4	88
16	Lattice Convolutional Neural Network Modeling of Adsorbate Coverage Effects. <i>Journal of Physical Chemistry C</i> , 2019, 123, 18951-18959.	3.1	21
17	Machine learning for renewable energy materials. <i>Journal of Materials Chemistry A</i> , 2019, 7, 17096-17117.	10.3	207
18	Unveiling new stable manganese based photoanode materials via theoretical high-throughput screening and experiments. <i>Chemical Communications</i> , 2019, 55, 13418-13421.	4.1	18

#	ARTICLE	IF	CITATIONS
19	Microkinetic modeling of aqueous phase biomass conversion: Application to ethylene glycol reforming. <i>Chemical Engineering Science</i> , 2019, 197, 415-418.	3.8	12
20	Thermochemistry of gas-phase and surface species via LASSO-assisted subgraph selection. <i>Reaction Chemistry and Engineering</i> , 2018, 3, 454-466.	3.7	29
21	Group Additivity for Aqueous Phase Thermochemical Properties of Alcohols on Pt(111). <i>Journal of Physical Chemistry C</i> , 2017, 121, 21510-21519.	3.1	27
22	Mechanism of Dehydration of Phenols on Noble Metals via First-Principles Microkinetic Modeling. <i>ACS Catalysis</i> , 2016, 6, 3047-3055.	11.2	69
23	Group Additivity for Thermochemical Property Estimation of Lignin Monomers on Pt(111). <i>Journal of Physical Chemistry C</i> , 2016, 120, 19234-19241.	3.1	18
24	Guaiacol Hydrodeoxygenation Mechanism on Pt(111): Insights from Density Functional Theory and Linear Free Energy Relations. <i>ChemSusChem</i> , 2015, 8, 315-322.	6.8	109
25	Tailoring electrode hydrophobicity to improve anode performance in alkaline media. <i>Journal of Power Sources</i> , 2013, 242, 581-588.	7.8	7