

Jinggang Lan

List of Publications by Year in descending order

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23
papers

854
citations

623734

14
h-index

677142

22
g-index

34
all docs

34
docs citations

34
times ranked

1065
citing authors

#	ARTICLE	IF	CITATIONS
1	Origin of Efficient Catalytic Combustion of Methane over $\text{Co}_3\text{O}_4(110)$: Active Low-Coordination Lattice Oxygen and Cooperation of Multiple Active Sites. <i>ACS Catalysis</i> , 2016, 6, 5508-5519.	11.2	116
2	Bifunctional Single Atom Electrocatalysts: Coordination-Performance Correlations and Reaction Pathways. <i>ACS Nano</i> , 2020, 14, 13279-13293.	14.6	107
3	Dynamics and control of active sites in hierarchically nanostructured cobalt phosphide/chalcogenide-based electrocatalysts for water splitting. <i>Energy and Environmental Science</i> , 2022, 15, 727-739.	30.8	96
4	Pure Siliceous Zeolite-Supported Ru Single-Atom Active Sites for Ammonia Synthesis. <i>Chemistry of Materials</i> , 2019, 31, 9413-9421.	6.7	83
5	Ultralow-temperature CO oxidation on an In_2O_3 - Co_3O_4 catalyst: a strategy to tune CO adsorption strength and oxygen activation simultaneously. <i>Chemical Communications</i> , 2014, 50, 6835-6838.	4.1	73
6	Modeling Electrified $\text{Pt}(111)\text{-H}_2\text{O}$ Interfaces from Ab Initio Molecular Dynamics. <i>Jacs Au</i> , 2021, 1, 569-577.	7.9	56
7	Ionization of Water as an Effect of Quantum Delocalization at Aqueous Electrode Interfaces. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 3724-3730.	4.6	39
8	Characterization of the Platinum-Hydrogen Bond by Surface-Sensitive Time-Resolved Infrared Spectroscopy. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 1254-1259.	4.6	38
9	Simulating the ghost: quantum dynamics of the solvated electron. <i>Nature Communications</i> , 2021, 12, 766.	12.8	36
10	First-Principles Simulations of an Aqueous $\text{CO}/\text{Pt}(111)$ Interface. <i>Journal of Physical Chemistry C</i> , 2018, 122, 24068-24076.	3.1	35
11	Solution Phase and Surface Photoisomerization of a Hydrazone Switch with a Long Thermal Half-Life. <i>Journal of the American Chemical Society</i> , 2019, 141, 17637-17645.	13.7	30
12	In situ spectroelectrochemical probing of CO redox landscape on copper single-crystal surfaces. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2022, 119, .	7.1	27
13	Inexpensive modeling of quantum dynamics using path integral generalized Langevin equation thermostats. <i>Journal of Chemical Physics</i> , 2020, 152, 124104.	3.0	26
14	In-situ nanospectroscopic imaging of plasmon-induced two-dimensional [4+4]-cycloaddition polymerization on $\text{Au}(111)$. <i>Nature Communications</i> , 2021, 12, 4557.	12.8	24
15	Monitoring surface transformations of metal carbodiimide water oxidation catalysts by operando XAS and Raman spectroscopy. <i>Dalton Transactions</i> , 2018, 47, 10759-10766.	3.3	11
16	CO_2 adsorption on the pristine and reduced $\text{CeO}_2(111)$ surface: Geometries and vibrational spectra by first principles simulations. <i>Journal of Chemical Physics</i> , 2021, 154, 094702.	3.0	11
17	Toward High-level Machine Learning Potential for Water Based on Quantum Fragmentation and Neural Networks. <i>Journal of Physical Chemistry A</i> , 2022, 126, 3926-3936.	2.5	11
18	Efficient Quantum Vibrational Spectroscopy of Water with High-Order Path Integrals: From Bulk to Interfaces. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 9108-9114.	4.6	9

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19	Shallow and deep trap states of solvated electrons in methanol and their formation, electronic excitation, and relaxation dynamics. <i>Chemical Science</i> , 2022, 13, 3837-3844.	7.4	9
20	Nanoscale Chemical Imaging of Coadsorbed Thiolate Self-Assembled Monolayers on Au(111) by Tip-Enhanced Raman Spectroscopy. <i>Analytical Chemistry</i> , 2022, 94, 1645-1653.	6.5	5
21	Effect of the alkyl linker length on the photoisomerization of hydrazone switches on metal surfaces. <i>Materials Today Chemistry</i> , 2022, 24, 100797.	3.5	4
22	Nuclear quantum effects at aqueous metal interfaces captured by molecular dynamics simulations. <i>Current Opinion in Electrochemistry</i> , 2022, 33, 100934.	4.8	2
23	Theoretical Study of NO Dissociation on an Open Flat Ru(101̄...1) Surface. <i>Journal of Physical Chemistry C</i> , 2019, 123, 5488-5494.	3.1	0