Benjamin Wei Jie Chen

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Computational Methods in Heterogeneous Catalysis. Chemical Reviews, 2021, 121, 1007-1048.	47.7	198
2	Zeoliteâ€Encaged Pd–Mn Nanocatalysts for CO ₂ Hydrogenation and Formic Acid Dehydrogenation. Angewandte Chemie - International Edition, 2020, 59, 20183-20191.	13.8	175
3	Palladiumâ€Based Nanocatalyst for Oneâ€Pot Synthesis of Polysubstituted Quinolines. ChemCatChem, 2013, 5, 277-283.	3.7	55
4	Atomic and Molecular Adsorption on Ag(111). Journal of Physical Chemistry C, 2019, 123, 7551-7566.	3.1	39
5	Formic Acid: A Hydrogen-Bonding Cocatalyst for Formate Decomposition. ACS Catalysis, 2020, 10, 10812-10825.	11.2	36
6	Density functional theory study of thermodynamic and kinetic isotope effects of H2/D2 dissociative adsorption on transition metals. Catalysis Science and Technology, 2018, 8, 3321-3335.	4.1	26
7	How coverage influences thermodynamic and kinetic isotope effects for H ₂ /D ₂ dissociative adsorption on transition metals. Catalysis Science and Technology, 2020, 10, 671-689.	4.1	26
8	Zeoliteâ€Encaged Pd–Mn Nanocatalysts for CO ₂ Hydrogenation and Formic Acid Dehydrogenation. Angewandte Chemie, 2020, 132, 20358-20366.	2.0	22
9	Kinetic Isolation between Turnovers on Au ₁₈ Nanoclusters: Formic Acid Decomposition One Molecule at a Time. ACS Catalysis, 2019, 9, 9446-9457.	11.2	20
10	Role of Hydrogen-bonded Bimolecular Formic Acid–Formate Complexes for Formic Acid Decomposition on Copper: A Combined First-Principles and Microkinetic Modeling Study. ACS Catalysis, 2021, 11, 4349-4361.	11.2	19
11	Effects of composition and morphology on the hydrogen storage properties of transition metal hydrides: Insights from PtPd nanoclusters. Nano Energy, 2019, 63, 103858.	16.0	15
12	Unraveling the Synergistic Effect of Re and Cs Promoters on Ethylene Epoxidation over Silver Catalysts with Machine Learning-Accelerated First-Principles Simulations. ACS Catalysis, 2022, 12, 2540-2551.	11.2	13
13	An automated cluster surface scanning method for exploring reaction paths on metal-cluster surfaces. Computational Materials Science, 2021, 186, 110010.	3.0	10
14	Exploring driving forces for length growth in graphene nanoribbons during chemical vapor deposition of hydrocarbons on Ge(0Â0Â1) via kinetic Monte Carlo simulations. Applied Surface Science, 2020, 527, 146784.	6.1	8
15	Identification of stable adsorption sites and diffusion paths on nanocluster surfaces: an automated scanning algorithm. Npj Computational Materials, 2019, 5, .	8.7	6
16	Carboxylic acid formation by hydroxyl insertion into acyl moieties on late transition metals. Catalysis Science and Technology, 2017, 7, 5365-5375.	4.1	2