

Bart Zijlstra

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/4330389/publications.pdf>

Version: 2024-02-01

22
papers

1,112
citations

430874

18
h-index

677142

22
g-index

23
all docs

23
docs citations

23
times ranked

1652
citing authors

#	ARTICLE	IF	CITATIONS
1	Optimum Cu nanoparticle catalysts for CO ₂ hydrogenation towards methanol. <i>Nano Energy</i> , 2018, 43, 200-209.	16.0	133
2	Efficient Base-Metal NiMn/TiO ₂ Catalyst for CO ₂ Methanation. <i>ACS Catalysis</i> , 2019, 9, 7823-7839.	11.2	124
3	Understanding carbon dioxide activation and carbon-carbon coupling over nickel. <i>Nature Communications</i> , 2019, 10, 5330.	12.8	124
4	Theoretical Approach To Predict the Stability of Supported Single-Atom Catalysts. <i>ACS Catalysis</i> , 2019, 9, 3289-3297.	11.2	101
5	Influence of Carbon Deposits on the Cobalt-Catalyzed Fischer-Tropsch Reaction: Evidence of a Two-Site Reaction Model. <i>ACS Catalysis</i> , 2018, 8, 1580-1590.	11.2	61
6	An Active Alkali-Exchanged Faujasite Catalyst for <i>p</i> -Xylene Production via the One-Pot Diels-Alder Cycloaddition/Dehydration Reaction of 2,5-Dimethylfuran with Ethylene. <i>ACS Catalysis</i> , 2018, 8, 760-769.	11.2	54
7	Mechanism of Cobalt-Catalyzed CO Hydrogenation: 1. Methanation. <i>ACS Catalysis</i> , 2017, 7, 8050-8060.	11.2	53
8	Optimum Particle Size for Gold-Catalyzed CO Oxidation. <i>Journal of Physical Chemistry C</i> , 2018, 122, 8327-8340.	3.1	45
9	Improved Pd/CeO ₂ Catalysts for Low-Temperature NO Reduction: Activation of CeO ₂ Lattice Oxygen by Fe Doping. <i>ACS Catalysis</i> , 2021, 11, 5614-5627.	11.2	44
10	Mechanism of Carbon Monoxide Dissociation on a Cobalt Fischer-Tropsch Catalyst. <i>ChemCatChem</i> , 2018, 10, 136-140.	3.7	39
11	Coverage Effects in CO Dissociation on Metallic Cobalt Nanoparticles. <i>ACS Catalysis</i> , 2019, 9, 7365-7372.	11.2	37
12	The Vital Role of Step-Edge Sites for Both CO Activation and Chain Growth on Cobalt Fischer-Tropsch Catalysts Revealed through First-Principles-Based Microkinetic Modeling Including Lateral Interactions. <i>ACS Catalysis</i> , 2020, 10, 9376-9400.	11.2	37
13	Quantum-Chemical DFT Study of Direct and H- and C-Assisted CO Dissociation on the γ -Fe ₅ C ₂ H ₂ Agg Carbide. <i>Journal of Physical Chemistry C</i> , 2018, 122, 9929-9938.	3.1	34
14	First-principles microkinetics simulations of electrochemical reduction of CO ₂ over Cu catalysts. <i>Electrochimica Acta</i> , 2020, 335, 135665.	5.2	32
15	First-principles based microkinetic modeling of transient kinetics of CO hydrogenation on cobalt catalysts. <i>Catalysis Today</i> , 2020, 342, 131-141.	4.4	29
16	Enumerating Active Sites on Metal Nanoparticles: Understanding the Size Dependence of Cobalt Particles for CO Dissociation. <i>ACS Catalysis</i> , 2021, 11, 8484-8492.	11.2	26
17	A theoretical study of the reverse water-gas shift reaction on Ni(111) and Ni(311) surfaces. <i>Canadian Journal of Chemical Engineering</i> , 2020, 98, 740-748.	1.7	25
18	Kinetic aspects of chain growth in Fischer-Tropsch synthesis. <i>Faraday Discussions</i> , 2017, 197, 153-164.	3.2	18

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19	A quantum-chemical study of the CO dissociation mechanism on low-index Miller planes of γ -Fe ₃ C. <i>Catalysis Today</i> , 2020, 342, 152-160.	4.4	15
20	A quantum-chemical DFT study of CO dissociation on Fe-promoted stepped Rh surfaces. <i>Catalysis Today</i> , 2016, 275, 111-118.	4.4	12
21	Quantum-chemical-based microkinetics simulations of syngas conversion over MoS ₂ (110) surface. <i>Chemical Engineering Science</i> , 2019, 198, 166-183.	3.8	5
22	Mechanism of Carbon Monoxide Dissociation on a Cobalt Fischer-Tropsch Catalyst. <i>ChemCatChem</i> , 2018, 10, 5-5.	3.7	1