

Ivana Matanovic

List of Publications by Year in descending order

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70
papers

3,409
citations

136740

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docs citations

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times ranked

4133
citing authors

#	ARTICLE	IF	CITATIONS
1	Protonated phosphonic acid electrodes for high power heavy-duty vehicle fuel cells. <i>Nature Energy</i> , 2022, 7, 248-259.	19.8	65
2	Highly Durable and Selective Fe- and Mo-Based Atomically Dispersed Electrocatalysts for Nitrate Reduction to Ammonia via Distinct and Synergized NO ₂ Pathways. <i>ACS Catalysis</i> , 2022, 12, 6651-6662.	5.5	58
3	Robust palladium hydride catalyst for electrocatalytic formate formation with high CO tolerance. <i>Applied Catalysis B: Environmental</i> , 2022, 316, 121659.	10.8	11
4	Synergistically integrated phosphonated poly(pentafluorostyrene) for fuel cells. <i>Nature Materials</i> , 2021, 20, 370-377.	13.3	112
5	Identification of durable and non-durable Fe _x sites in Fe-N-C materials for proton exchange membrane fuel cells. <i>Nature Catalysis</i> , 2021, 4, 10-19.	16.1	368
6	First principles inelastic mean free paths coupled with Monte Carlo simulation of secondary electron yield of Cu-Ni, Cu-Zn, and Mo-Li. <i>Journal of Applied Physics</i> , 2021, 129, .	1.1	8
7	Protocol for rapid ammonia detection via surface-enhanced Raman spectroscopy. <i>STAR Protocols</i> , 2021, 2, 100599.	0.5	0
8	Self-Anchored Platinum-Decorated Antimony-Doped-Tin Oxide as a Durable Oxygen Reduction Electrocatalyst. <i>ACS Catalysis</i> , 2021, 11, 7006-7017.	5.5	17
9	Calculation of Secondary Electron Yield of Alloys: Single Pole Approximation. , 2021, , .		0
10	Energetics of Base-Acid Pairs for the Design of High-Temperature Fuel Cell Polymer Electrolytes. <i>Journal of Physical Chemistry B</i> , 2020, 124, 7725-7734.	1.2	23
11	Towards defect engineering in hexagonal MoS ₂ nanosheets for tuning hydrogen evolution and nitrogen reduction reactions. <i>Applied Materials Today</i> , 2020, 21, 100812.	2.3	16
12	Facile All-Optical Method for In Situ Detection of Low Amounts of Ammonia. <i>IScience</i> , 2020, 23, 101757.	1.9	12
13	Assessing Stability of Transition Metal Nitrides in Aqueous Environments: The Case of Molybdenum, Iron, Vanadium and Nickel Nitride. <i>Journal of the Electrochemical Society</i> , 2020, 167, 046518.	1.3	7
14	Spectro-Electrochemical Microfluidic Platform for Monitoring Multi-Step Cascade Reactions. <i>ChemElectroChem</i> , 2019, 6, 246-251.	1.7	10
15	Understanding the Oxygen Reduction Reaction Activity and Oxidative Stability of Pt Supported on Nb-Doped TiO ₂ . <i>ChemSusChem</i> , 2019, 12, 3409-3409.	3.6	0
16	Investigating the Nature of the Active Sites for the CO ₂ Reduction Reaction on Carbon-Based Electrocatalysts. <i>ACS Catalysis</i> , 2019, 9, 7668-7678.	5.5	58
17	Understanding Active Sites in Pyrolyzed Fe-N-C Catalysts for Fuel Cell Cathodes by Bridging Density Functional Theory Calculations and ⁵⁷ Fe Mössbauer Spectroscopy. <i>ACS Catalysis</i> , 2019, 9, 9359-9371.	5.5	167
18	Adsorption of Polyaromatic Backbone Impacts the Performance of Anion Exchange Membrane Fuel Cells. <i>Chemistry of Materials</i> , 2019, 31, 4195-4204.	3.2	91

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19	Modular Microfluidic Paper-Based Devices for Multi-Modal Cascade Catalysis. <i>ChemElectroChem</i> , 2019, 6, 2448-2455.	1.7	8
20	The energetics of phosphoric acid interactions reveals a new acid loss mechanism. <i>Journal of Materials Chemistry A</i> , 2019, 7, 9867-9876.	5.2	83
21	Understanding the Oxygen Reduction Reaction Activity and Oxidative Stability of Pt Supported on Nb-Doped TiO ₂ . <i>ChemSusChem</i> , 2019, 12, 3468-3480.	3.6	39
22	Thermochemical analysis of Mo-C-H system for synthesis of molybdenum carbides. <i>Thermochimica Acta</i> , 2019, 676, 27-32.	1.2	4
23	Selectivity control for electroreduction of ketones. <i>Nature Catalysis</i> , 2019, 2, 186-187.	16.1	5
24	Phenyl Oxidation Impacts the Durability of Alkaline Membrane Water Electrolyzer. <i>ACS Applied Materials & Interfaces</i> , 2019, 11, 9696-9701.	4.0	79
25	Design of Pd-Pb Catalysts for Glycerol and Ethylene Glycol Electrooxidation in Alkaline Medium. <i>Electrocatalysis</i> , 2018, 9, 480-485.	1.5	20
26	Effect of pH on the Activity of Platinum Group Metal-Free Catalysts in Oxygen Reduction Reaction. <i>ACS Catalysis</i> , 2018, 8, 3041-3053.	5.5	158
27	Nitrogen electroreduction and hydrogen evolution on cubic molybdenum carbide: a density functional study. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 14679-14687.	1.3	55
28	Inhibition of Surface Chemical Moieties by Tris(hydroxymethyl)aminomethane: A Key to Understanding Oxygen Reduction on Iron-Nitrogen-Carbon Catalysts. <i>ACS Applied Energy Materials</i> , 2018, 1, 1942-1949.	2.5	18
29	Role of Surface Chemistry on Catalyst/Ionomer Interactions for Transition Metal-Nitrogen-Carbon Electrocatalysts. <i>ACS Applied Energy Materials</i> , 2018, 1, 68-77.	2.5	44
30	Understanding PGM-free catalysts by linking density functional theory calculations and structural analysis: Perspectives and challenges. <i>Current Opinion in Electrochemistry</i> , 2018, 9, 137-144.	2.5	85
31	Impact of ionomer adsorption on alkaline hydrogen oxidation activity and fuel cell performance. <i>Current Opinion in Electrochemistry</i> , 2018, 12, 189-195.	2.5	55
32	Mechanism of Oxygen Reduction Reaction on Transition Metal-Nitrogen-Carbon Catalysts: Establishing the Role of Nitrogen-containing Active Sites. <i>ACS Applied Energy Materials</i> , 2018, 1, 5948-5953.	2.5	54
33	Rational design of polyaromatic ionomers for alkaline membrane fuel cells with >1 W cm ⁻² power density. <i>Energy and Environmental Science</i> , 2018, 11, 3283-3291.	15.6	209
34	Surface Adsorption Affects the Performance of Alkaline Anion-Exchange Membrane Fuel Cells. <i>ACS Catalysis</i> , 2018, 8, 9429-9439.	5.5	55
35	Cascade Kinetics of an Artificial Metabolon by Molecular Dynamics and Kinetic Monte Carlo. <i>ACS Catalysis</i> , 2018, 8, 7719-7726.	5.5	13
36	Electro-Reduction of Nitrogen on Molybdenum Carbides: A Density Functional Theory Study. <i>ECS Meeting Abstracts</i> , 2018, . .	0.0	0

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37	Oxygen Binding to Active Sites of Fe-N-C ORR Electrocatalysts Observed by Ambient-Pressure XPS. <i>Journal of Physical Chemistry C</i> , 2017, 121, 2836-2843.	1.5	135
38	Air Breathing Cathodes for Microbial Fuel Cell using Mn-, Fe-, Co- and Ni-containing Platinum Group Metal-free Catalysts. <i>Electrochimica Acta</i> , 2017, 231, 115-124.	2.6	131
39	Novel Hybrid Catalyst for the Oxidation of Organic Acids: Pd Nanoparticles Supported on Mn-Graphene Nanosheets. <i>ChemElectroChem</i> , 2017, 4, 2336-2344.	1.7	5
40	Outer membrane cytochromes/flavin interactions in <i>Shewanella</i> spp. A molecular perspective. <i>Biointerphases</i> , 2017, 12, 021004.	0.6	24
41	Hybrid molecular/enzymatic catalytic cascade for complete electro-oxidation of glycerol using a promiscuous NAD-dependent formate dehydrogenase from <i>Candida boidinii</i> . <i>Chemical Communications</i> , 2017, 53, 5368-5371.	2.2	23
42	Benzene Adsorption: A Significant Inhibitor for the Hydrogen Oxidation Reaction in Alkaline Conditions. <i>Journal of Physical Chemistry Letters</i> , 2017, 8, 4918-4924.	2.1	55
43	Evaluation of Pt Alloys as Electrocatalysts for Oxalic Acid Oxidation: A Combined Experimental and Computational Study. <i>Journal of the Electrochemical Society</i> , 2016, 163, H787-H795.	1.3	4
44	Effect of Organic Cations on Hydrogen Oxidation Reaction of Carbon Supported Platinum. <i>Journal of the Electrochemical Society</i> , 2016, 163, F1503-F1509.	1.3	29
45	Core Level Shifts of Hydrogenated Pyridinic and Pyrrolic Nitrogen in the Nitrogen-Containing Graphene-Based Electrocatalysts: In-Plane vs Edge Defects. <i>Journal of Physical Chemistry C</i> , 2016, 120, 29225-29232.	1.5	123
46	Highly active and selective nickel molybdenum catalysts for direct hydrazine fuel cell. <i>Electrochimica Acta</i> , 2016, 215, 420-426.	2.6	59
47	Cation-Hydroxide-Water Coadsorption Inhibits the Alkaline Hydrogen Oxidation Reaction. <i>Journal of Physical Chemistry Letters</i> , 2016, 7, 4464-4469.	2.1	57
48	Protein-Support Interactions for Rationally Designed Bilirubin Oxidase Based Cathode: A Computational Study. <i>Journal of Physical Chemistry B</i> , 2016, 120, 3634-3641.	1.2	24
49	Tolerance of non-platinum group metals cathodes proton exchange membrane fuel cells to air contaminants. <i>Journal of Power Sources</i> , 2016, 324, 556-571.	4.0	34
50	Functional interfaces for biomimetic energy harvesting: CNTs-DNA matrix for enzyme assembly. <i>Biochimica Et Biophysica Acta - Bioenergetics</i> , 2016, 1857, 612-620.	0.5	5
51	Highly-active Pd-Cu electrocatalysts for oxidation of ubiquitous oxygenated fuels. <i>Applied Catalysis B: Environmental</i> , 2016, 191, 76-85.	10.8	61
52	Role of Quinones in Electron Transfer of PQQ-Glucose Dehydrogenase Anodes-Mediation or Orientation Effect. <i>Journal of the American Chemical Society</i> , 2015, 137, 7754-7762.	6.6	34
53	Predicting Electrocatalytic Properties: Modeling Structure-Activity Relationships of Nitroxyl Radicals. <i>Journal of the American Chemical Society</i> , 2015, 137, 16179-16186.	6.6	91
54	CuCo ₂ O ₄ ORR/OER Bi-Functional Catalyst: Influence of Synthetic Approach on Performance. <i>Journal of the Electrochemical Society</i> , 2015, 162, F449-F454.	1.3	104

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55	Bio-inspired design of electrocatalysts for oxalate oxidation: a combined experimental and computational study of Mn ^{II} -N ^{II} -C catalysts. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 13235-13244.	1.3	26
56	The study of secondary effects in vibrational and hydrogen bonding properties of 2- and 3-ethynylpyridine and ethynylbenzene by IR spectroscopy. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2014, 132, 215-224.	2.0	10
57	Quinone-Modified Surfaces for Enhanced Enzyme-Electrode Interactions in Pyrroloquinoline-Quinone-Dependent Glucose Dehydrogenase Anodes. <i>ChemElectroChem</i> , 2014, 1, 2017-2028.	1.7	14
58	Electro-reduction of nitrogen on molybdenum nitride: structure, energetics, and vibrational spectra from DFT. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 3014.	1.3	55
59	Applicability of density functional theory in reproducing accurate vibrational spectra of surface bound species. <i>Journal of Computational Chemistry</i> , 2014, 35, 1921-1929.	1.5	2
60	Effect of enzymatic orientation through the use of syringaldazine molecules on multiple multi-copper oxidase enzymes. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 13367-13375.	1.3	39
61	Density Functional Theory Study of Oxygen Reduction Activity on Ultrathin Platinum Nanotubes. <i>Journal of Physical Chemistry C</i> , 2012, 116, 16499-16510.	1.5	18
62	Theoretical Study of Electrochemical Processes on Pt-Ni Alloys. <i>Journal of Physical Chemistry C</i> , 2011, 115, 10640-10650.	1.5	79
63	Fluxionality of Hydrogen Ligands in Fe(H) ₂ (H ₂) ₂ (PEtPh) ₂ ₃ . <i>Inorganic Chemistry</i> , 2011, 50, 10740-10747.	1.9	15
64	Sulfur Ylide Promoted Synthesis of N-Protected Aziridines: A Combined Experimental and Computational Approach. <i>Chemistry - A European Journal</i> , 2010, 16, 11744-11752.	1.7	23
65	Bergman Cyclization of Acyclic Amino Acid Derived Eneidyne Leads to the Formation of 2,3-Dihydrobenzo[f]isoindoles. <i>Journal of Organic Chemistry</i> , 2010, 75, 6219-6228.	1.7	14
66	Methane molecule confined in the small and large cages of structure I clathrate hydrate: Quantum six-dimensional calculations of the coupled translation-rotation eigenstates. <i>Journal of Chemical Physics</i> , 2009, 131, 224308.	1.2	20
67	Generalized approximation to the reaction path: The formic acid dimer case. <i>Journal of Chemical Physics</i> , 2008, 128, 084103.	1.2	26
68	Theoretical modeling of the formic acid dimer infrared spectrum: Shaping the O-H stretch band. <i>Chemical Physics</i> , 2007, 338, 121-126.	0.9	22
69	Supramolecular amide and thioamide synthons in hydrogen bonding patterns of N-aryl-furamides and N-aryl-thiofuramides. <i>Structural Chemistry</i> , 2006, 17, 275-285.	1.0	9
70	Exploring the potential energy surface for proton transfer in acetylacetone. <i>Chemical Physics</i> , 2004, 306, 201-207.	0.9	31