

Joseph H Montoya

List of Publications by Year in descending order

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Version: 2024-02-01

37
papers

6,556
citations

218677

26
h-index

330143

37
g-index

39
all docs

39
docs citations

39
times ranked

9036
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|------|-----------|
| 1 | Materials for solar fuels and chemicals. <i>Nature Materials</i> , 2017, 16, 70-81. | 27.5 | 1,163 |
| 2 | The Challenge of Electrochemical Ammonia Synthesis: A New Perspective on the Role of Nitrogen Scaling Relations. <i>ChemSusChem</i> , 2015, 8, 2180-2186. | 6.8 | 1,018 |
| 3 | Theoretical Insights into a CO Dimerization Mechanism in CO ₂ Electroreduction. <i>Journal of Physical Chemistry Letters</i> , 2015, 6, 2032-2037. | 4.6 | 606 |
| 4 | Accelerating the discovery of materials for clean energy in the era of smart automation. <i>Nature Reviews Materials</i> , 2018, 3, 5-20. | 48.7 | 489 |
| 5 | Matminer: An open source toolkit for materials data mining. <i>Computational Materials Science</i> , 2018, 152, 60-69. | 3.0 | 446 |
| 6 | Understanding trends in C-H bond activation in heterogeneous catalysis. <i>Nature Materials</i> , 2017, 16, 225-229. | 27.5 | 387 |
| 7 | Insights into C-C Coupling in CO ₂ Electroreduction on Copper Electrodes. <i>ChemCatChem</i> , 2013, 5, 737-742. | 3.7 | 339 |
| 8 | A Review on Challenges and Successes in Atomic-Scale Design of Catalysts for Electrochemical Synthesis of Hydrogen Peroxide. <i>ACS Catalysis</i> , 2020, 10, 7495-7511. | 11.2 | 254 |
| 9 | Atomate: A high-level interface to generate, execute, and analyze computational materials science workflows. <i>Computational Materials Science</i> , 2017, 139, 140-152. | 3.0 | 223 |
| 10 | Electrochemical Stability of Metastable Materials. <i>Chemistry of Materials</i> , 2017, 29, 10159-10167. | 6.7 | 168 |
| 11 | Acetaldehyde as an Intermediate in the Electroreduction of Carbon Monoxide to Ethanol on Oxide-Derived Copper. <i>Angewandte Chemie - International Edition</i> , 2016, 55, 1450-1454. | 13.8 | 166 |
| 12 | Autonomous experimentation systems for materials development: A community perspective. <i>Matter</i> , 2021, 4, 2702-2726. | 10.0 | 143 |
| 13 | Robust and synthesizable photocatalysts for CO ₂ reduction: a data-driven materials discovery. <i>Nature Communications</i> , 2019, 10, 443. | 12.8 | 125 |
| 14 | Rutile Alloys in the Mn-Sb-O System Stabilize Mn ³⁺ To Enable Oxygen Evolution in Strong Acid. <i>ACS Catalysis</i> , 2018, 8, 10938-10948. | 11.2 | 97 |
| 15 | Trends in adsorption of electrocatalytic water splitting intermediates on cubic ABO ₃ oxides. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 3813-3818. | 2.8 | 94 |
| 16 | Random forest machine learning models for interpretable X-ray absorption near-edge structure spectrum-property relationships. <i>Npj Computational Materials</i> , 2020, 6, . | 8.7 | 94 |
| 17 | Computational Design of Active Site Structures with Improved Transition-State Scaling for Ammonia Synthesis. <i>ACS Catalysis</i> , 2018, 8, 4017-4024. | 11.2 | 80 |
| 18 | A high-throughput framework for determining adsorption energies on solid surfaces. <i>Npj Computational Materials</i> , 2017, 3, . | 8.7 | 70 |

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|----|--|------|-----------|
| 19 | Efficient Pourbaix diagrams of many-element compounds. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 25323-25327. | 2.8 | 69 |
| 20 | Predicting aqueous stability of solid with computed Pourbaix diagram using SCAN functional. <i>Npj Computational Materials</i> , 2020, 6, . | 8.7 | 69 |
| 21 | Theoretical evaluation of the surface electrochemistry of perovskites with promising photon absorption properties for solar water splitting. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 2634-2640. | 2.8 | 58 |
| 22 | Machine learningâ€“accelerated design and synthesis of polyelemental heterostructures. <i>Science Advances</i> , 2021, 7, eabj5505. | 10.3 | 53 |
| 23 | Autonomous intelligent agents for accelerated materials discovery. <i>Chemical Science</i> , 2020, 11, 8517-8532. | 7.4 | 49 |
| 24 | Rational Solid-State Synthesis Routes for Inorganic Materials. <i>Journal of the American Chemical Society</i> , 2021, 143, 9244-9259. | 13.7 | 48 |
| 25 | Origins of the Instability of Nonprecious Hydrogen Evolution Reaction Catalysts at Open-Circuit Potential. <i>ACS Energy Letters</i> , 2021, 6, 2268-2274. | 17.4 | 44 |
| 26 | Acetaldehyde as an Intermediate in the Electroreduction of Carbon Monoxide to Ethanol on Oxideâ€“Derived Copper. <i>Angewandte Chemie</i> , 2016, 128, 1472-1476. | 2.0 | 39 |
| 27 | BEEP: A Python library for Battery Evaluation and Early Prediction. <i>SoftwareX</i> , 2020, 11, 100506. | 2.6 | 29 |
| 28 | Electroreduction of Methanediol on Copper. <i>Catalysis Letters</i> , 2013, 143, 631-635. | 2.6 | 21 |
| 29 | Two-dimensional forms of robust CO2 reduction photocatalysts. <i>Npj 2D Materials and Applications</i> , 2020, 4, . | 7.9 | 20 |
| 30 | The Materials Research Platform: Defining the Requirements from User Stories. <i>Matter</i> , 2019, 1, 1433-1438. | 10.0 | 19 |
| 31 | Direct Water Decomposition on Transition Metal Surfaces: Structural Dependence and Catalytic Screening. <i>Catalysis Letters</i> , 2016, 146, 718-724. | 2.6 | 18 |
| 32 | Toward autonomous materials research: Recent progress and future challenges. <i>Applied Physics Reviews</i> , 2022, 9, . | 11.3 | 17 |
| 33 | Automated Adsorption Workflow for Semiconductor Surfaces and the Application to Zinc Telluride. <i>Journal of Chemical Information and Modeling</i> , 2021, 61, 3908-3916. | 5.4 | 11 |
| 34 | Agents for sequential learning using multiple-fidelity data. <i>Scientific Reports</i> , 2022, 12, 4694. | 3.3 | 9 |
| 35 | Polymer Structure Predictor (PSP): A Python Toolkit for Predicting Atomic-Level Structural Models for a Range of Polymer Geometries. <i>Journal of Chemical Theory and Computation</i> , 2022, 18, 2737-2748. | 5.3 | 7 |
| 36 | Novel inorganic crystal structures predicted using autonomous simulation agents. <i>Scientific Data</i> , 2022, 9, . | 5.3 | 7 |

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|----|---|-----|-----------|
| 37 | MaterialNet: A web-based graph explorer for materials science data. Journal of Open Source Software, 2020, 5, 2105. | 4.6 | 5 |