

Sergio Posada-PÃ©rez

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Effect of Aqueous Electrolytes on LiCoO_2 Surfaces: Role of Proton Adsorption on Oxygen Vacancy Formation. <i>Journal of Physical Chemistry C</i> , 2022, 126, 110-119.	3.1	7
2	The importance of the bite angle of metal(III) salen catalysts in the sequestration of CO_2 with epoxides in mild conditions. <i>Green Chemical Engineering</i> , 2022, 3, 180-187.	6.3	18
3	Exploring cocatalyst type effect on the Ziegler-Natta catalyzed ethylene polymerizations: experimental and DFT studies. <i>Journal of Polymer Research</i> , 2022, 29, .	2.4	19
4	The Importance of Li-Ion Nature and Stacking in Layered Cathode Materials for Aqueous Ion Batteries: From Bulk to Surface Models. <i>ECS Meeting Abstracts</i> , 2021, MA2021-01, 99-99.	0.0	0
5	Influence of Stacking on H^+ Intercalation in Layered CoO_2 ($\text{A} = \text{Tj}$) Investigation. <i>Chemistry of Materials</i> , 2021, 33, 6942-6954.	0.784314	15
6	Are Protons Involve in the Fading Capacity in Li Ion Aqueous Batteries? Shedding Light By Means of First Principle Computations. <i>ECS Meeting Abstracts</i> , 2020, MA2020-01, 223-223.	0.0	0
7	Shedding Light about the Role of Proton Intercalation in Layered Cathode Materials By Means of First Principle Computations. <i>ECS Meeting Abstracts</i> , 2020, MA2020-02, 68-68.	0.0	0
8	Kinetic Monte Carlo Simulations Unveil Synergic Effects at Work on Bifunctional Catalysts. <i>ACS Catalysis</i> , 2019, 9, 9117-9126.	11.2	30
9	CO_2 interaction with violarite (FeNi_2S_4) surfaces: a dispersion-corrected DFT study. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 20439-20446.	2.8	15
10	Highly active Au/MoC and $\text{Au}/\text{Mo}_2\text{C}$ catalysts for the low-temperature water gas shift reaction: effects of the carbide metal/carbon ratio on the catalyst performance. <i>Catalysis Science and Technology</i> , 2017, 7, 5332-5342.	4.1	39
11	Adsorption and dissociation of molecular hydrogen on orthorhombic Mo_2C and cubic MoC (001) surfaces. <i>Surface Science</i> , 2017, 656, 24-32.	1.9	50
12	Highly Active Au/MoC and Cu/MoC Catalysts for the Conversion of CO_2 : The Metal/C Ratio as a Key Factor Defining Activity, Selectivity, and Stability. <i>Journal of the American Chemical Society</i> , 2016, 138, 8269-8278.	13.7	140
13	The conversion of CO_2 to methanol on orthorhombic Mo_2C and $\text{Cu}/\text{Mo}_2\text{C}$ catalysts: mechanism for admetal induced change in the selectivity and activity. <i>Catalysis Science and Technology</i> , 2016, 6, 6766-6777.	4.1	101
14	Methane capture at room temperature: adsorption on cubic MoC and orthorhombic Mo_2C molybdenum carbide (001) surfaces. <i>RSC Advances</i> , 2015, 5, 33737-33746.	3.6	18
15	Structure and electronic properties of Cu nanoclusters supported on $\text{Mo}_2\text{C}(001)$ and $\text{MoC}(001)$ surfaces. <i>Journal of Chemical Physics</i> , 2015, 143, 114704.	3.0	25
16	Fundamentals of Methanol Synthesis on Metal Carbide Based Catalysts: Activation of CO_2 and H_2 . <i>Topics in Catalysis</i> , 2015, 58, 159-173.	2.8	64
17	The bending machine: CO_2 activation and hydrogenation on $\text{MoC}(001)$ and $\text{Mo}_2\text{C}(001)$ surfaces. <i>Physical Chemistry Chemical Physics</i> , 2014, 16, 14912-14921.	2.8	175