

# MarÃ-a F Juarez

## List of Publications by Year in descending order

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Version: 2024-02-01

26  
papers

723  
citations

840776

11  
h-index

552781

26  
g-index

27  
all docs

27  
docs citations

27  
times ranked

1228  
citing authors

#	ARTICLE	IF	CITATIONS
1	Composition and Electronic Structure of Mn <sub>3</sub> O <sub>4</sub> and Co <sub>3</sub> O <sub>4</sub> Cathodes in Zinc-Air Batteries: A DFT Study. <i>Journal of Physical Chemistry C</i> , 2022, 126, 2561-2572.	3.1	3
2	Interactions of ions across carbon nanotubes. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 10603-10608.	2.8	4
3	Role of the Partial Charge Transfer on the Chloride Adlayers on Au(100). <i>ChemElectroChem</i> , 2020, 7, 4269-4282.	3.4	10
4	Interaction between chloride ions mediated by carbon nanotubes: a chemical attraction. <i>Journal of Solid State Electrochemistry</i> , 2020, 24, 3207-3214.	2.5	5
5	Sulfate, Bisulfate, and Hydrogen Co-adsorption on Pt(111) and Au(111) in an Electrochemical Environment. <i>Frontiers in Chemistry</i> , 2020, 8, 634.	3.6	43
6	Hydrogen Oxidation in Alkaline Media: the Bifunctional Mechanism for Water Formation. <i>Electrocatalysis</i> , 2019, 10, 584-590.	3.0	7
7	Tuning the rate of an outer-sphere electron transfer by changing the electronic structure of carbon nanotubes. <i>Journal of Electroanalytical Chemistry</i> , 2019, 847, 113186.	3.8	10
8	An Unusual Exchange Mechanism in the Tafel Reaction on Pt(110) (1 $\bar{1}$ -1) Surfaces. <i>ChemElectroChem</i> , 2019, 6, 3279-3284.	3.4	4
9	Why are trace amounts of chloride so highly surface-active?. <i>Journal of Electroanalytical Chemistry</i> , 2019, 847, 113128.	3.8	2
10	The initial stage of OH adsorption on Ni(111). <i>Journal of Electroanalytical Chemistry</i> , 2019, 832, 137-141.	3.8	7
11	Defying Coulomb's law: A lattice-induced attraction between lithium ions. <i>Carbon</i> , 2018, 139, 808-812.	10.3	10
12	Oxygen Reduction in Alkaline Media—a Discussion. <i>Electrocatalysis</i> , 2017, 8, 554-564.	3.0	17
13	Interaction of Hydrogen with Au Modified by Pd and Rh in View of Electrochemical Applications. <i>Computation</i> , 2016, 4, 26.	2.0	6
14	On the Energetics of Ions in Carbon and Gold Nanotubes. <i>ChemPhysChem</i> , 2016, 17, 78-85.	2.1	19
15	Combined ab initio and XPS Investigations of the Electronic Interactions of L-Cysteine Adsorbed on GaAs(1 0 0). <i>ChemistrySelect</i> , 2016, 1, 3623-3634.	1.5	1
16	A scenario for oxygen reduction in alkaline media. <i>Nano Energy</i> , 2016, 29, 362-368.	16.0	15
17	A scenario for oxygen reduction in alkaline media. <i>Nano Energy</i> , 2016, 26, 558-564.	16.0	20
18	Oxygen Reduction on Ag(100) in Alkaline Solutions—A Theoretical Study. <i>ChemPhysChem</i> , 2016, 17, 500-505.	2.1	12

#	ARTICLE	IF	CITATIONS
19	Nanotubes for charge storage “ towards an atomistic model. <i>Electrochimica Acta</i> , 2015, 162, 11-16.	5.2	31
20	Role of surface contaminants, functionalities, defects and electronic structure: general discussion. <i>Faraday Discussions</i> , 2014, 172, 365-395.	3.2	1
21	Carbon electrodes for energy storage: general discussion. <i>Faraday Discussions</i> , 2014, 172, 239-260.	3.2	11
22	Spontaneous formation of metallic nanostructures on highly oriented pyrolytic graphite (HOPG): an ab initio and experimental study. <i>Faraday Discussions</i> , 2014, 172, 327-347.	3.2	14
23	Screening of ions in carbon and gold nanotubes “ A theoretical study. <i>Electrochemistry Communications</i> , 2014, 45, 48-51.	4.7	34
24	Volcano plots in hydrogen electrocatalysis “ uses and abuses. <i>Beilstein Journal of Nanotechnology</i> , 2014, 5, 846-854.	2.8	410
25	Electronic Anisotropy at Vicinal Ag(111) Surfaces: Work Function Changes Induced by Steps and Hydrogen Adsorption. <i>Journal of Physical Chemistry C</i> , 2013, 117, 4606-4618.	3.1	23
26	The role of the organic layer functionalization in the formation of silicon/organic layer/metal junctions with coinage metals. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 21411.	2.8	4