

# Vladimir Tripkovic

## List of Publications by Year in descending order

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36  
papers

4,655  
citations

236833

25  
h-index

414303

32  
g-index

38  
all docs

38  
docs citations

38  
times ranked

6506  
citing authors

#	ARTICLE	IF	CITATIONS
1	Modeling the Electrochemical Hydrogen Oxidation and Evolution Reactions on the Basis of Density Functional Theory Calculations. <i>Journal of Physical Chemistry C</i> , 2010, 114, 18182-18197.	1.5	990
2	Tuning the activity of Pt alloy electrocatalysts by means of the lanthanide contraction. <i>Science</i> , 2016, 352, 73-76.	6.0	783
3	The oxygen reduction reaction mechanism on Pt(111) from density functional theory calculations. <i>Electrochimica Acta</i> , 2010, 55, 7975-7981.	2.6	491
4	Universal transition state scaling relations for (de)hydrogenation over transition metals. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 20760.	1.3	363
5	Modeling the electrified solid-liquid interface. <i>Chemical Physics Letters</i> , 2008, 466, 68-71.	1.2	349
6	Electrochemical CO <sub>2</sub> and CO Reduction on Metal-Functionalized Porphyrin-like Graphene. <i>Journal of Physical Chemistry C</i> , 2013, 117, 9187-9195.	1.5	260
7	The Pt(111)/Electrolyte Interface under Oxygen Reduction Reaction Conditions: An Electrochemical Impedance Spectroscopy Study. <i>Langmuir</i> , 2011, 27, 2058-2066.	1.6	170
8	pH in atomic scale simulations of electrochemical interfaces. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 10321.	1.3	127
9	Standard hydrogen electrode and potential of zero charge in density functional calculations. <i>Physical Review B</i> , 2011, 84, .	1.1	118
10	Intermetallic Alloys as CO Electroreduction Catalysts-Role of Isolated Active Sites. <i>ACS Catalysis</i> , 2014, 4, 2268-2273.	5.5	101
11	On the pH dependence of electrochemical proton transfer barriers. <i>Catalysis Today</i> , 2016, 262, 36-40.	2.2	91
12	Oxygen Reduction Reaction on Pt Overlayers Deposited onto a Gold Film: Ligand, Strain, and Ensemble Effect. <i>ACS Catalysis</i> , 2016, 6, 671-676.	5.5	79
13	Density functional theory based screening of ternary alkali-transition metal borohydrides: A computational material design project. <i>Journal of Chemical Physics</i> , 2009, 131, 014101.	1.2	77
14	The Influence of Particle Shape and Size on the Activity of Platinum Nanoparticles for Oxygen Reduction Reaction: A Density Functional Theory Study. <i>Catalysis Letters</i> , 2014, 144, 380-388.	1.4	66
15	Potential- and Rate-Determining Step for Oxygen Reduction on Pt(111). <i>Journal of Physical Chemistry C</i> , 2017, 121, 26785-26793.	1.5	56
16	Avoiding pitfalls in the modeling of electrochemical interfaces. <i>Chemical Physics Letters</i> , 2013, 555, 145-148.	1.2	50
17	From 3D to 2D Co and Ni Oxyhydroxide Catalysts: Elucidation of the Active Site and Influence of Doping on the Oxygen Evolution Activity. <i>ACS Catalysis</i> , 2017, 7, 8558-8571.	5.5	50
18	Insight into the Effect of Sn on CO and Formic Acid Oxidation at PtSn Catalysts. <i>Journal of Physical Chemistry C</i> , 2014, 118, 278-289.	1.5	48

#	ARTICLE	IF	CITATIONS
19	Comparative DFT+U and HSE Study of the Oxygen Evolution Electrocatalysis on Perovskite Oxides. <i>Journal of Physical Chemistry C</i> , 2018, 122, 1135-1147.	1.5	46
20	Metal Oxide-Supported Platinum Overlayers as Proton-Exchange Membrane Fuel Cell Cathodes. <i>ChemCatChem</i> , 2012, 4, 228-235.	1.8	44
21	First principles investigation of zinc-anode dissolution in zinc-air batteries. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 6416.	1.3	44
22	Thermodynamic assessment of the oxygen reduction activity in aqueous solutions. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 29381-29388.	1.3	43
23	First principles investigation of the activity of thin film Pt, Pd and Au surface alloys for oxygen reduction. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 11647-11657.	1.3	41
24	Computational Screening of Doped $\text{MnO}_2$ Catalysts for the Oxygen Evolution Reaction. <i>ChemSusChem</i> , 2018, 11, 629-637.	3.6	40
25	Comparison between the Oxygen Reduction Reaction Activity of $\text{Pd}_5\text{Ce}$ and $\text{Pt}_5\text{Ce}$ : The Importance of Crystal Structure. <i>ACS Catalysis</i> , 2015, 5, 6032-6040.	5.5	21
26	Trends for Methane Oxidation at Solid Oxide Fuel Cell Conditions. <i>Journal of the Electrochemical Society</i> , 2009, 156, B1447.	1.3	16
27	Modeling of the symmetry factor of electrochemical proton discharge via the Volmer reaction. <i>Catalysis Today</i> , 2013, 202, 168-174.	2.2	16
28	Platinum redispersion on metal oxides in low temperature fuel cells. <i>Physical Chemistry Chemical Physics</i> , 2013, 15, 3279.	1.3	10
29	First principles study of (Cd, Hg, In, Tl, Sn, Pb, As, Sb, Bi, Se) modified Pt(111), Pt(100) and Pt(211) electrodes as CO oxidation catalysts. <i>Electrochimica Acta</i> , 2015, 168, 370-378.	2.6	7
30	Formic Acid Oxidation at Platinum-Bismuth Clusters. <i>Journal of the Electrochemical Society</i> , 2014, 161, H547-H554.	1.3	5
31	Theoretical Study of the Structural Stability and the Electronic Properties of $\text{Al}_m\text{H}_n$ Clusters. <i>Journal of Computational and Theoretical Nanoscience</i> , 2011, 8, 609-615.	0.4	0
32	Exploring the Lanthanide Contraction to Tune the Activity and Stability of Pt. <i>ECS Meeting Abstracts</i> , 2016, , .	0.0	0
33	Tailoring the Performance of Oxides for the Oxygen Evolution Reaction. <i>ECS Meeting Abstracts</i> , 2016, , .	0.0	0
34	(Invited) Genetic Algorithms and DFT for Accelerated Design of Nanoalloys. <i>ECS Meeting Abstracts</i> , 2016, , .	0.0	0
35	Understanding and Tailoring the Performance of Transition Metal Oxides for the Oxygen Evolution Reaction. <i>ECS Meeting Abstracts</i> , 2017, , .	0.0	0
36	Understanding and Tailoring the Performance of Transition Metal Oxides for the Oxygen Evolution Reaction. <i>ECS Meeting Abstracts</i> , 2018, , .	0.0	0