Vladimir Tripkovic

List of Publications by Year
in descending order

36	4,655	236833	414303
papers	citations	25 h-index	32 g-index

38	38	38	6506
all docs	docs citations	times ranked	citing authors

#	Article	IF	Citations
1	Computational Screening of Doped Î±â€MnO₂ Catalysts for the Oxygen Evolution Reaction. ChemSusChem, 2018, 11, 629-637.	3.6	40
2	Comparative DFT+U and HSE Study of the Oxygen Evolution Electrocatalysis on Perovskite Oxides. Journal of Physical Chemistry C, 2018, 122, 1135-1147.	1.5	46
3	Understanding and Tailoring the Performance of Transition Metal Oxides for the Oxygen Evolution Reaction. ECS Meeting Abstracts, 2018, , .	0.0	0
4	Thermodynamic assessment of the oxygen reduction activity in aqueous solutions. Physical Chemistry Chemical Physics, 2017, 19, 29381-29388.	1.3	43
5	From 3D to 2D Co and Ni Oxyhydroxide Catalysts: Elucidation of the Active Site and Influence of Doping on the Oxygen Evolution Activity. ACS Catalysis, 2017, 7, 8558-8571.	5.5	50
6	Potential- and Rate-Determining Step for Oxygen Reduction on Pt(111). Journal of Physical Chemistry C, 2017, 121, 26785-26793.	1.5	56
7	Understanding and Tailoring the Performance of Transition Metal Oxides for the Oxygen Evolution Reaction. ECS Meeting Abstracts, 2017, , .	0.0	0
8	Tuning the activity of Pt alloy electrocatalysts by means of the lanthanide contraction. Science, 2016, 352, 73-76.	6.0	783
9	On the pH dependence of electrochemical proton transfer barriers. Catalysis Today, 2016, 262, 36-40.	2.2	91
10	Oxygen Reduction Reaction on Pt Overlayers Deposited onto a Gold Film: Ligand, Strain, and Ensemble Effect. ACS Catalysis, 2016, 6, 671-676.	5.5	79
11	Exploring the Lanthanide Contraction to Tune the Activity and Stability of Pt. ECS Meeting Abstracts, 2016, , .	0.0	0
12	Tailoring the Performance of Oxides for the Oxygen Evolution Reaction. ECS Meeting Abstracts, 2016, , .	0.0	0
13	(Invited) Genetic Algorithms and DFT for Accelerated Design of Nanoalloys. ECS Meeting Abstracts, 2016, , .	0.0	0
14	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. Electrochimica Acta, 2015, 168, 370-378.	2.6	7
15	First principles investigation of the activity of thin film Pt, Pd and Au surface alloys for oxygen reduction. Physical Chemistry Chemical Physics, 2015, 17, 11647-11657.	1.3	41
16	Comparison between the Oxygen Reduction Reaction Activity of Pd₅Ce and Pt₅Ce: The Importance of Crystal Structure. ACS Catalysis, 2015, 5, 6032-6040.	5.5	21
17	Formic Acid Oxidation at Platinum-Bismuth Clusters. Journal of the Electrochemical Society, 2014, 161, H547-H554.	1.3	5
18	The Influence of Particle Shape and Size on the Activity of Platinum Nanoparticles for Oxygen Reduction Reaction: A Density Functional Theory Study. Catalysis Letters, 2014, 144, 380-388.	1.4	66

#	Article	IF	Citations
19	Insight into the Effect of Sn on CO and Formic Acid Oxidation at PtSn Catalysts. Journal of Physical Chemistry C, 2014, 118, 278-289.	1.5	48
20	Intermetallic Alloys as CO Electroreduction Catalystsâ€"Role of Isolated Active Sites. ACS Catalysis, 2014, 4, 2268-2273.	5.5	101
21	Platinum redispersion on metal oxides in low temperature fuel cells. Physical Chemistry Chemical Physics, 2013, 15, 3279.	1.3	10
22	Modeling of the symmetry factor of electrochemical proton discharge via the Volmer reaction. Catalysis Today, 2013, 202, 168-174.	2.2	16
23	Avoiding pitfalls in the modeling of electrochemical interfaces. Chemical Physics Letters, 2013, 555, 145-148.	1.2	50
24	Electrochemical CO₂ and CO Reduction on Metal-Functionalized Porphyrin-like Graphene. Journal of Physical Chemistry C, 2013, 117, 9187-9195.	1.5	260
25	First principles investigation of zinc-anode dissolution in zincâ€"air batteries. Physical Chemistry Chemical Physics, 2013, 15, 6416.	1.3	44
26	pH in atomic scale simulations of electrochemical interfaces. Physical Chemistry Chemical Physics, 2013, 15, 10321.	1.3	127
27	Metal Oxideâ€Supported Platinum Overlayers as Protonâ€Exchange Membrane Fuel Cell Cathodes. ChemCatChem, 2012, 4, 228-235.	1.8	44
28	Universal transition state scaling relations for (de)hydrogenation over transition metals. Physical Chemistry Chemical Physics, 2011, 13, 20760.	1.3	363
29	Theoretical Study of the Structural Stability and the Electronic Properties of Al_mH_n Clusters. Journal of Computational and Theoretical Nanoscience, 2011, 8, 609-615.	0.4	0
30	The Pt(111)/Electrolyte Interface under Oxygen Reduction Reaction Conditions: An Electrochemical Impedance Spectroscopy Study. Langmuir, 2011, 27, 2058-2066.	1.6	170
31	Standard hydrogen electrode and potential of zero charge in density functional calculations. Physical Review B, 2011, 84, .	1.1	118
32	The oxygen reduction reaction mechanism on Pt(111) from density functional theory calculations. Electrochimica Acta, 2010, 55, 7975-7981.	2.6	491
33	Modeling the Electrochemical Hydrogen Oxidation and Evolution Reactions on the Basis of Density Functional Theory Calculations. Journal of Physical Chemistry C, 2010, 114, 18182-18197.	1.5	990
34	Trends for Methane Oxidation at Solid Oxide Fuel Cell Conditions. Journal of the Electrochemical Society, 2009, 156, B1447.	1.3	16
35	Density functional theory based screening of ternary alkali-transition metal borohydrides: A computational material design project. Journal of Chemical Physics, 2009, 131, 014101.	1.2	77
36	Modeling the electrified solidâ€"liquid interface. Chemical Physics Letters, 2008, 466, 68-71.	1.2	349