

# Ismail-Can OÄuz

## List of Publications by Year in descending order

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Version: 2024-02-01

11  
papers

704  
citations

1477746

6  
h-index

1473754

9  
g-index

11  
all docs

11  
docs citations

11  
times ranked

1045  
citing authors

#	ARTICLE	IF	CITATIONS
1	Identification of durable and non-durable Fe <sub>Nx</sub> sites in Fe-N-C materials for proton exchange membrane fuel cells. <i>Nature Catalysis</i> , 2021, 4, 10-19.	16.1	368
2	Assessing the performances of different continuum solvation models for the calculation of hydration energies of molecules, polymers and surfaces: a comparison between the SMD, VASPsol and FDPB models. <i>Theoretical Chemistry Accounts</i> , 2021, 140, 1.	0.5	13
3	Generalizing Continuum Solvation in Crystal to Nonaqueous Solvents: Implementation, Parametrization, and Application to Molecules and Surfaces. <i>Journal of Chemical Theory and Computation</i> , 2021, 17, 6432-6448.	2.3	2
4	P-block single-metal-site tin/nitrogen-doped carbon fuel cell cathode catalyst for oxygen reduction reaction. <i>Nature Materials</i> , 2020, 19, 1215-1223.	13.3	278
5	Predicting the Activity of Nano-Transition-Metal DeNox Catalysts. <i>Journal of Physical Chemistry C</i> , 2019, 123, 20314-20318.	1.5	11
6	Theoretical Prediction of the Distribution of Spin Moment on Metal-N-C Catalyst Embedded in Truncated Graphene Sheets. <i>ECS Meeting Abstracts</i> , 2019, , .	0.0	0
7	Can Density Functional Theory Predict Mössbauer Spectra in Pyrolyzed Fe-N-C Catalysts?. <i>ECS Meeting Abstracts</i> , 2019, , .	0.0	0
8	The effect of Pd ensemble structure on the O <sub>2</sub> dissociation and CO oxidation mechanisms on Au-Pd(100) surface alloys. <i>Journal of Chemical Physics</i> , 2018, 148, 024701.	1.2	15
9	Equilibrium Au-Pd(100) Surface Structures under CO Pressure: Energetic Stabilities and Phase Diagrams. <i>Journal of Physical Chemistry C</i> , 2018, 122, 18922-18932.	1.5	2
10	On the structural and electronic properties of Ir-silicide nanowires on Si(001) surface. <i>Journal of Applied Physics</i> , 2016, 120, .	1.1	8
11	Investigation of finite-size effects in chemical bonding of AuPd nanoalloys. <i>Journal of Chemical Physics</i> , 2015, 143, 144309.	1.2	7