

# Sateesh Bandaru

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/5778185/publications.pdf>

Version: 2024-02-01

10

papers

106

citations

1478505

6

h-index

1372567

10

g-index

10

all docs

10

docs citations

10

times ranked

240

citing authors

#	ARTICLE	IF	CITATIONS
1	Facet-Dependent Bactericidal Activity of Ag <sub>3</sub> PO <sub>4</sub> Nanostructures against Gram-Positive/Negative Bacteria. ACS Omega, 2022, 7, 16616-16628.	3.5	5
2	Sulfur-doped wood-derived porous carbon for optimizing electromagnetic response performance. Nanoscale, 2021, 13, 16084-16093.	5.6	6
3	Structural and Electronic Properties of MgO/TiO <sub>2</sub> Interfaces: A First-Principles Molecular-Simulation Study. Journal of Physical Chemistry C, 2021, 125, 10795-10802.	3.1	2
4	Oxygen-evolution reactions (OER) on transition-metal-doped Fe <sub>3</sub> Co(PO <sub>4</sub> ) <sub>4</sub> iron-phosphate surfaces: a first-principles study. Catalysis Science and Technology, 2021, 11, 4619-4626.	4.1	4
5	Novel Superstructure-Phase Two-Dimensional Material 1T-VSe <sub>2</sub> at High Pressure. Journal of Physical Chemistry Letters, 2020, 11, 380-386.	4.6	17
6	First-principles studies on $\text{Fe}_{2}\text{O}_3$ surface slabs and mechanistic elucidation of a g-C <sub>3</sub> N <sub>4</sub> / $\text{Fe}_{2}\text{O}_3$ heterojunction. Catalysis Science and Technology, 2020, 10, 1376-1384.	4.1	20
7	Ternary structure reveals mechanism of a membrane diacylglycerol kinase. Nature Communications, 2015, 6, 10140.	12.8	30
8	Implicit and explicit solvent models for modeling a bifunctional arene ruthenium hydrogen storage catalyst: A classical and ab initio molecular simulation study. Journal of Computational Chemistry, 2014, 35, 683-691.	3.3	6
9	Towards the design of novel boron- and nitrogen-substituted ammonia-borane and bifunctional arene ruthenium catalysts for hydrogen storage. Journal of Computational Chemistry, 2014, 35, 891-903.	3.3	7
10	Effect of space linkers in dinuclear copper cryptates on the efficiency of atmospheric CO <sub>2</sub> uptake: a DFT study. Catalysis Science and Technology, 2013, 3, 2234.	4.1	9