

# Sambath Baskaran

## List of Publications by Year in descending order

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Version: 2024-02-01

34  
papers

648  
citations

623188

14  
h-index

580395

25  
g-index

35  
all docs

35  
docs citations

35  
times ranked

725  
citing authors

#	ARTICLE	IF	CITATIONS
1	Mo <sub>2</sub> CS <sub>2</sub> -MXene supported single-atom catalysts for efficient and selective CO <sub>2</sub> electrochemical reduction. <i>Applied Surface Science</i> , 2022, 592, 153339.	3.1	20
2	Identifying Key Descriptors for the Single-Atom Catalyzed CO Oxidation. <i>CCS Chemistry</i> , 2022, 4, 3296-3308.	4.6	25
3	Phosphorene Supported Single-Atom Catalysts for CO Oxidation: A Computational Study. <i>ChemPhysChem</i> , 2021, 22, 378-385.	1.0	12
4	A study on the interaction of Nile blue with Uracils: A spectroscopic and computational approach. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2021, 246, 119011.	2.0	5
5	Non-noble metal single-atom catalyst of Co <sub>1</sub> /MXene (Mo <sub>2</sub> CS <sub>2</sub> ) for CO oxidation. <i>Science China Materials</i> , 2021, 64, 651-663.	3.5	44
6	Profuse Surface Activation of Ir-Dispersed Titanium Nitride Bifunctional Electrocatalysts. <i>Advanced Energy and Sustainability Research</i> , 2021, 2, 2000054.	2.8	5
7	Unveiling the In Situ Generation of a Monovalent Fe(I) Site in the Single-Fe-Atom Catalyst for Electrochemical CO <sub>2</sub> Reduction. <i>ACS Catalysis</i> , 2021, 11, 7292-7301.	5.5	51
8	Cobalt Oxide on a Nanoporous TUD-1 Catalyst for Methylene Blue Dye Interaction DFT Studies and Degradation. <i>Symmetry</i> , 2021, 13, 1754.	1.1	1
9	Chromium Single-Atom Catalyst with Graphyne Support: A Theoretical Study of NO Oxidation and Reduction. <i>ACS Catalysis</i> , 2020, 10, 11951-11961.	5.5	49
10	Non-noble metal single-atom catalysts with phosphotungstic acid (PTA) support: A theoretical study of ethylene epoxidation. <i>Science China Materials</i> , 2020, 63, 1003-1014.	3.5	41
11	Construction of frustrated Lewis pair from nitride and phosphine for the activation and cleavage of molecular hydrogen. <i>Applied Organometallic Chemistry</i> , 2020, 34, e5811.	1.7	0
12	Catalytic mechanism and bonding analyses of Au-Pd single atom alloy (SAA): CO oxidation reaction. <i>Science China Materials</i> , 2020, 63, 993-1002.	3.5	23
13	Spectroscopic investigation and computational studies on the interaction of Acriflavine with various estrogens. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2019, 206, 622-629.	2.0	5
14	NiCoFe oxide amorphous nanoheterostructures for oxygen evolution reaction. <i>International Journal of Hydrogen Energy</i> , 2019, 44, 22991-23001.	3.8	39
15	Exploring the potential of novel transition metal complexes derived from ONO donor type ligand: a quantum chemical study. <i>Journal of Molecular Modeling</i> , 2019, 25, 284.	0.8	3
16	Inkjet-printed phosphorescent Iridium(III) complex based paper sensor for highly selective detection of Hg <sup>2+</sup> . <i>Dyes and Pigments</i> , 2019, 163, 176-182.	2.0	22
17	Development of paper-based chemosensor for the detection of mercury ions using mono- and tetra-sulfur bearing phenanthridines. <i>New Journal of Chemistry</i> , 2018, 42, 8530-8536.	1.4	25
18	Experimental and theoretical studies on the corrosion inhibition of vitamins " Thiamine hydrochloride or biotin in corrosion of mild steel in aqueous chloride environment. <i>Egyptian Journal of Petroleum</i> , 2018, 27, 371-381.	1.2	25

#	ARTICLE	IF	CITATIONS
19	Possibility of reducing the coordinated dinitrogen into ammonia and hydrazine using [Ru-L] (L =) Tj ETQq1 1 0.784314 rgBT /Overlock 10	0.7	10
20	Hydroxylamine synthesis by oxygen insertion into $\text{Re}\xi\zeta\text{NH}<\text{sub}>2</\text{sub}>$ bond via Baeyerâ€“Villiger oxidation: a Theoretical study. Journal of Physical Organic Chemistry, 2015, 28, 690-694.	0.9	0
21	Ionic and Neutral Halfâ€“Sandwich Guanidinatoruthenium(II) Complexes and Their Solution Behavior. European Journal of Inorganic Chemistry, 2015, 2015, 3182-3194.	1.0	13
22	Functionalization of $\text{N}_2$ to $\text{NH}_3$ via direct $\text{N}\hat{\%}\text{o}j\text{N}$ bond cleavage using $\text{M(III)(NMe}_2)_3$ (M=W/Mo): A theoretical study. Journal of Chemical Sciences, 2015, 127, 83-94.	0.7	4
23	Slow hydrolysis of an organozirconium complex: The first polyoxometallic heptanuclear zirconium oxide. Journal of Organometallic Chemistry, 2015, 775, 76-79.	0.8	2
24	Lessons learned and lessons to be learned for developing homogeneous transition metal complexes catalyzed reduction of $\text{N}_2$ to ammonia. Journal of Organometallic Chemistry, 2014, 752, 44-58.	0.8	45
25	Molecular and electronic structure analysis of some novel copper and zinc complexes of hypervalent carbon based ligand: DFT studies. Journal of Organometallic Chemistry, 2014, 752, 123-132.	0.8	2
26	Ammonia and hydrazine synthesis from $[\text{N}_2\text{-W}\{(\text{NHCH}_2\text{CH}_2)_3\text{N}\}]$ and $[\text{AH}]^+[\text{BH}]^{\text{â}^{\text{~}}}$ using Sivasankar catalytic cycle: DFT studies. Computational and Theoretical Chemistry, 2014, 1027, 73-78.	1.1	9
27	Calix[2]bispyrrolylarenes: New Expanded Calix[4]pyrroles for Fluorometric Sensing of Anions via Extended $\text{I}\hat{\text{e}}\text{-Conjugation}$ . Organic Letters, 2013, 15, 306-309.	2.4	34
28	Reduction of $\text{N}_2$ by $\text{H}_2$ to $\text{NH}_3$ and $\text{N}_2\text{H}_4$ using $[\text{MoL}]$ (L=triamidoamine) and organic co-catalysts: A theoretical approach. Journal of Molecular Catalysis A, 2013, 370, 140-144.	4.8	14
29	Hydrogenation of dinitrogen to ammonia in $[\text{WF}(\text{PH}_2(\text{CH}_2)_2\text{PH}_2)_2\text{N}_2]$ using $\text{H}_2$ : Insights from DFT calculations. New Journal of Chemistry, 2012, 36, 562.	1.4	14
30	Colorimetric Sensing of Fluoride Ion by New Expanded Calix[4]pyrrole through Anionâ€“ $\text{I}\hat{\text{e}}$ Interaction. Organic Letters, 2012, 14, 548-551.	2.4	72
31	A possibility of functionalizing the dinitrogen in a Chatt complex by $\text{H}_2$ : Density functional studies. Polyhedron, 2012, 31, 676-681.	1.0	13
32	A $[\text{Fe}(\text{CB}<\text{sub}>6</\text{sub}>)]$ platform for binding of small molecules: Insights from DFT calculations. Journal of Computational Chemistry, 2012, 33, 1047-1054.	1.5	4
33	Understanding the stability, electronic and molecular structure of some copper(III) complexes containing alkyl and non alkyl ligands: Insights from DFT calculations. Journal of Organometallic Chemistry, 2011, 696, 2627-2634.	0.8	12
34	Functionalization of Dinitrogen Using a Historically Significant Ru Complex: A New Life for an Old Complex. European Journal of Inorganic Chemistry, 2010, 2010, 4716-4719.	1.0	5