

Nishamol Kuriakose

List of Publications by Year in descending order

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Version: 2024-02-01

14
papers

246
citations

1307594

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1125743

13
g-index

15
all docs

15
docs citations

15
times ranked

255
citing authors

#	ARTICLE	IF	CITATIONS
1	Rhodium Complexes Bearing PAIP Pincer Ligands. <i>Journal of the American Chemical Society</i> , 2018, 140, 7070-7073.	13.7	96
2	CO ₂ capture, activation and dissociation on the Ti ₂ C surface and Ti ₂ C MXene: the role of surface structure. <i>Physical Chemistry Chemical Physics</i> , 2020, 22, 14599-14612.	2.8	30
3	Characterization of Rh-Al Bond in Rh(PAIP) (PAIP = Pincer-type Diphosphino-Alumanyl Ligand) in Comparison with Rh(L)(PMe ₃) ₂ (L = AlMe ₂), <i>Tj ETQq1 1 0.784314 rgBT / Overlock 10 Tf 50 66</i>	4.0	27
4	A Theoretical Study of Metal-Metal Cooperativity in the Homogeneous Water Gas Shift Reaction. <i>Inorganic Chemistry</i> , 2012, 51, 377-385.	4.0	25
5	New insights into small molecule activation by acyclic silylenes: a computational investigation. <i>Dalton Transactions</i> , 2014, 43, 2194-2201.	3.3	20
6	Asymmetric transfer hydrogenation of imines in water/methanol co-solvent system and mechanistic investigation by DFT study. <i>RSC Advances</i> , 2014, 4, 46351-46356.	3.6	18
7	Ethene Dimerization and Hydrogenation over a Zeolite-Supported Rh(I)-Carbonyl Complex: Mechanistic Insights from DFT Modeling. <i>ACS Catalysis</i> , 2018, 8, 9836-9846.	11.2	14
8	C-C coupling at a zeolite-supported Rh(<i>scpi</i>) complex. DFT search for the mechanism. <i>Catalysis Science and Technology</i> , 2019, 9, 2781-2793.	4.1	8
9	Can Molecular Cages Be Effective at Small Molecule Activation? A Computational Investigation. <i>Inorganic Chemistry</i> , 2013, 52, 4238-4243.	4.0	2
10	Can substituted allenes be highly efficient leaving groups in catalytic processes? A computational investigation. <i>Journal of Computational Chemistry</i> , 2015, 36, 795-804.	3.3	2
11	CH ₄ activation and C-C coupling on the Ti ₂ C(100) surface in the presence of intrinsic C-vacancies: is excess good?. <i>Journal of Materials Chemistry A</i> , 2021, 9, 23703-23713.	10.3	2
12	Can main group systems act as superior catalysts for dihydrogen generation reactions? A computational investigation. <i>Dalton Transactions</i> , 2016, 45, 5968-5977.	3.3	1
13	Coverage dependent CO ₂ activation on Ti ₂ C(111) surface: Effect of intrinsic subsurface Carbon vacancies. <i>Surface Science</i> , 2021, 706, 121798.	1.9	1
14	Modeling the effect of ligands and solvation on hydrolysis variants in the Pd(II)-Catalyzed hydroxycarbonylation of pentenoic acids. <i>Journal of Organometallic Chemistry</i> , 2020, 914, 121221.	1.8	0