

# Riddhish Pandharkar

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9137593/publications.pdf>

Version: 2024-02-01

10  
papers

162  
citations

1307594

7  
h-index

1372567

10  
g-index

17  
all docs

17  
docs citations

17  
times ranked

142  
citing authors

#	ARTICLE	IF	CITATIONS
1	Insights into the Structure–Activity Relationships in Metal–Organic Framework-Supported Nickel Catalysts for Ethylene Hydrogenation. <i>ACS Catalysis</i> , 2020, 10, 8995-9005.	11.2	40
2	A New Mixing of Nonlocal Exchange and Nonlocal Correlation with Multiconfiguration Pair-Density Functional Theory. <i>Journal of Physical Chemistry Letters</i> , 2020, 11, 10158-10163.	4.6	21
3	Excited States of Crystalline Point Defects with Multireference Density Matrix Embedding Theory. <i>Journal of Physical Chemistry Letters</i> , 2021, 12, 11688-11694.	4.6	20
4	Variational Localized Active Space Self-Consistent Field Method. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 4923-4937.	5.3	19
5	Electronic structure of strongly correlated systems: recent developments in multiconfiguration pair-density functional theory and multiconfiguration nonclassical-energy functional theory. <i>Chemical Science</i> , 2022, 13, 7685-7706.	7.4	18
6	Isomerization and Selective Hydrogenation of Propyne: Screening of Metal–Organic Frameworks Modified by Atomic Layer Deposition. <i>Journal of the American Chemical Society</i> , 2020, 142, 20380-20389.	13.7	15
7	Spin-State Ordering in Metal-Based Compounds Using the Localized Active Space Self-Consistent Field Method. <i>Journal of Physical Chemistry Letters</i> , 2019, 10, 5507-5513.	4.6	11
8	Localized Active Space Pair-Density Functional Theory. <i>Journal of Chemical Theory and Computation</i> , 2021, 17, 2843-2851.	5.3	6
9	Investigating the effect of metal nuclearity on activity for ethylene hydrogenation by metal-organic-framework-supported oxy-Ni(II) catalysts. <i>Journal of Catalysis</i> , 2022, 407, 162-173.	6.2	5
10	Factors Affecting the Mechanism of 1,3-Butadiene Polymerization at Open Metal Sites in Co-MFU-4l. <i>Organometallics</i> , 2022, 41, 169-177.	2.3	3