

Yongjin Lee

List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1516937/publications.pdf>

Version: 2024-02-01

34
papers

1,703
citations

394421

19
h-index

395702

33
g-index

34
all docs

34
docs citations

34
times ranked

2048
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|------|-----------|
| 1 | Deciphering van der Waals interaction between polypropylene and carbonated fly ash from experimental and molecular simulation. <i>Journal of Hazardous Materials</i> , 2022, 421, 126725. | 12.4 | 5 |
| 2 | In Silico Generation of a Topologically Diverse Zeolite-Templated Carbon Library. <i>Crystal Growth and Design</i> , 2022, 22, 123-130. | 3.0 | 3 |
| 3 | Control over interpenetration for boosting methane storage capacity in metal-organic frameworks. <i>Journal of Materials Chemistry A</i> , 2021, 9, 24857-24862. | 10.3 | 14 |
| 4 | Machine Learning-Driven Discovery of Metal-Organic Frameworks for Efficient CO ₂ Capture in Humid Condition. <i>ACS Sustainable Chemistry and Engineering</i> , 2021, 9, 2872-2879. | 6.7 | 34 |
| 5 | Coating the Right Polymer: Achieving Ideal Metal-Organic Framework Particle Dispersibility in Polymer Matrixes Using a Coordinative Crosslinking Surface Modification Method. <i>Angewandte Chemie</i> , 2021, 133, 14257-14264. | 2.0 | 14 |
| 6 | Coating the Right Polymer: Achieving Ideal Metal-Organic Framework Particle Dispersibility in Polymer Matrixes Using a Coordinative Crosslinking Surface Modification Method. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 14138-14145. | 13.8 | 48 |
| 7 | Physicochemical Understanding of the Impact of Pore Environment and Species of Adsorbates on Adsorption Behaviour. <i>Angewandte Chemie</i> , 2021, 133, 20667-20673. | 2.0 | 1 |
| 8 | Physicochemical Understanding of the Impact of Pore Environment and Species of Adsorbates on Adsorption Behaviour. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 20504-20510. | 13.8 | 8 |
| 9 | Machine Learning-based approach for Tailor-Made design of ionic Liquids: Application to CO ₂ capture. <i>Separation and Purification Technology</i> , 2021, 275, 119117. | 7.9 | 17 |
| 10 | <i>In Situ</i> Mapping and Local Negative Uptake Behavior of Adsorbates in Individual Pores of Metal-Organic Frameworks. <i>Journal of the American Chemical Society</i> , 2021, 143, 20747-20757. | 13.7 | 5 |
| 11 | A computational study to design zeolite-templated carbon materials with high performance for CO ₂ /N ₂ separation. <i>Microporous and Mesoporous Materials</i> , 2020, 295, 109947. | 4.4 | 12 |
| 12 | Machine Learning Enabled Tailor-Made Design of Application-Specific Metal-Organic Frameworks. <i>ACS Applied Materials & Interfaces</i> , 2020, 12, 734-743. | 8.0 | 42 |
| 13 | Enhancing the Gas Separation Selectivity of Mixed-Matrix Membranes Using a Dual-Interfacial Engineering Approach. <i>Journal of the American Chemical Society</i> , 2020, 142, 18503-18512. | 13.7 | 86 |
| 14 | Tuning Metal-Organic Framework Nanocrystal Shape through Facet-Dependent Coordination. <i>Nano Letters</i> , 2020, 20, 1774-1780. | 9.1 | 52 |
| 15 | Understanding the diversity of the metal-organic framework ecosystem. <i>Nature Communications</i> , 2020, 11, 4068. | 12.8 | 282 |
| 16 | Robust Metal-Triazolate Frameworks for CO ₂ Capture from Flue Gas. <i>Journal of the American Chemical Society</i> , 2020, 142, 2750-2754. | 13.7 | 159 |
| 17 | Engineering plasticization resistant gas separation membranes using metal-organic nanocapsules. <i>Chemical Science</i> , 2020, 11, 4687-4694. | 7.4 | 22 |
| 18 | Tracking and Visualization of Functional Domains in Stratified Metal-Organic Frameworks Using Gold Nanoparticles. <i>ACS Central Science</i> , 2020, 6, 247-253. | 11.3 | 13 |

| # | ARTICLE | IF | CITATIONS |
|----|--|------|-----------|
| 19 | Machine Learning Prediction on Properties of Nanoporous Materials Utilizing Pore Geometry Barcodes. <i>Journal of Chemical Information and Modeling</i> , 2019, 59, 4636-4644. | 5.4 | 29 |
| 20 | Understanding Adsorption Behavior of Periodic Mesoporous Organosilica Having a Heterogeneous Chemical Environment: Selective Coverage and Interpenetration of Adsorbates inside the Channel Wall. <i>Journal of Physical Chemistry C</i> , 2019, 123, 24884-24889. | 3.1 | 6 |
| 21 | A generalizable method for the construction of MOF@polymer functional composites through surface-initiated atom transfer radical polymerization. <i>Chemical Science</i> , 2019, 10, 1816-1822. | 7.4 | 75 |
| 22 | General Way To Construct Micro- and Mesoporous Metal-Organic Framework-Based Porous Liquids. <i>Journal of the American Chemical Society</i> , 2019, 141, 19708-19714. | 13.7 | 111 |
| 23 | Strong thermal conductivity dependence on arsenic-vacancy complex formation in arsenic-doped silicon. <i>Journal of Applied Physics</i> , 2019, 126, 195104. | 2.5 | 1 |
| 24 | High-Throughput Screening Approach for Nanoporous Materials Genome Using Topological Data Analysis: Application to Zeolites. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 4427-4437. | 5.3 | 53 |
| 25 | Generating carbon schwarzites via zeolite-templating. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018, 115, E8116-E8124. | 7.1 | 88 |
| 26 | Quantifying similarity of pore-geometry in nanoporous materials. <i>Nature Communications</i> , 2017, 8, 15396. | 12.8 | 98 |
| 27 | Engineering of Pore Geometry for Ultrahigh Capacity Methane Storage in Mesoporous Metal-Organic Frameworks. <i>Journal of the American Chemical Society</i> , 2017, 139, 13300-13303. | 13.7 | 140 |
| 28 | Computational development of the nanoporous materials genome. <i>Nature Reviews Materials</i> , 2017, 2, . | 48.7 | 123 |
| 29 | What is the thermal conductivity limit of silicon germanium alloys?. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 19544-19548. | 2.8 | 18 |
| 30 | Fundamental insight into control of thermal conductivity in silicon-germanium alloy nanowires. <i>Materials Research Society Symposia Proceedings</i> , 2014, 1707, 31. | 0.1 | 0 |
| 31 | Microsegregation effects on the thermal conductivity of silicon-germanium alloys. <i>Journal of Applied Physics</i> , 2013, 114, 174910. | 2.5 | 12 |
| 32 | Mechanism of thermal conductivity suppression in doped silicon studied with nonequilibrium molecular dynamics. <i>Physical Review B</i> , 2012, 86, . | 3.2 | 30 |
| 33 | Force-matching-based parameterization of the Stillinger-Weber potential for thermal conduction in silicon. <i>Physical Review B</i> , 2012, 85, . | 3.2 | 37 |
| 34 | Effects of vacancy defects on thermal conductivity in crystalline silicon: A nonequilibrium molecular dynamics study. <i>Physical Review B</i> , 2011, 83, . | 3.2 | 65 |