

# Kevin Maik Jablonka

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/9524357/publications.pdf>

Version: 2024-02-01

14  
papers

1,092  
citations

840585

11  
h-index

1125617

13  
g-index

27  
all docs

27  
docs citations

27  
times ranked

886  
citing authors

| #  | ARTICLE  | IF   | CITATIONS |
|----|--|------|-----------|
| 1  | Big-Data Science in Porous Materials: Materials Genomics and Machine Learning. <i>Chemical Reviews</i> , 2020, 120, 8066-8129.   | 23.0 | 284       |
| 2  | Understanding the diversity of the metal-organic framework ecosystem. <i>Nature Communications</i> , 2020, 11, 4068.   | 5.8  | 282       |
| 3  | The Role of Machine Learning in the Understanding and Design of Materials. <i>Journal of the American Chemical Society</i> , 2020, 142, 20273-20287.   | 6.6  | 179       |
| 4  | Bias free multiobjective active learning for materials design and discovery. <i>Nature Communications</i> , 2021, 12, 2312.  | 5.8  | 78        |
| 5  | Charge Separation and Charge Carrier Mobility in Photocatalytic Metal-Organic Frameworks. <i>Advanced Functional Materials</i> , 2020, 30, 2003792.  | 7.8  | 64        |
| 6  | Diversifying Databases of Metal Organic Frameworks for High-Throughput Computational Screening. <i>ACS Applied Materials &amp; Interfaces</i> , 2021, 13, 61004-61014.                             | 4.0  | 50        |
| 7  | Using collective knowledge to assign oxidation states of metal cations in metal-organic frameworks. <i>Nature Chemistry</i> , 2021, 13, 771-777.   | 6.6  | 35        |
| 8  | Making the collective knowledge of chemistry open and machine actionable. <i>Nature Chemistry</i> , 2022, 14, 365-376.   | 6.6  | 34        |
| 9  | Applicability of Tail Corrections in the Molecular Simulations of Porous Materials. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 5635-5641.                                       | 2.3  | 30        |
| 10 | A data-driven perspective on the colours of metal-organic frameworks. <i>Chemical Science</i> , 2021, 12, 3587-3598.   | 3.7  | 16        |
| 11 | Making Molecules Vibrate: Interactive Web Environment for the Teaching of Infrared Spectroscopy. <i>Journal of Chemical Education</i> , 2022, 99, 561-569.   | 1.1  | 13        |
| 12 | Characterization of Chemisorbed Species and Active Adsorption Sites in Mg-Al Mixed Metal Oxides for High-Temperature CO <sub>2</sub> Capture. <i>Chemistry of Materials</i> , 2022, 34, 3893-3901. | 3.2  | 10        |
| 13 | Data-Driven Matching of Experimental Crystal Structures and Gas Adsorption Isotherms of Metal-Organic Frameworks. <i>Journal of Chemical &amp; Engineering Data</i> , 2022, 67, 1743-1756.         | 1.0  | 6         |
| 14 | Grundlagen der Thermodynamik für Studierende der Chemie. <i>Essentials</i> , 2017, , .   | 0.1  | 0         |