

# 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	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
2	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
3	Making the collective knowledge of chemistry open and machine actionable. <i>Nature Chemistry</i> , 2022, 14, 365-376.	6.6	34
4	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
5	Bias free multiobjective active learning for materials design and discovery. <i>Nature Communications</i> , 2021, 12, 2312.	5.8	78
6	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
7	A data-driven perspective on the colours of metal-organic frameworks. <i>Chemical Science</i> , 2021, 12, 3587-3598.	3.7	16
8	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
9	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
10	Understanding the diversity of the metal-organic framework ecosystem. <i>Nature Communications</i> , 2020, 11, 4068.	5.8	282
11	Charge Separation and Charge Carrier Mobility in Photocatalytic Metal-Organic Frameworks. <i>Advanced Functional Materials</i> , 2020, 30, 2003792.	7.8	64
12	Big-Data Science in Porous Materials: Materials Genomics and Machine Learning. <i>Chemical Reviews</i> , 2020, 120, 8066-8129.	23.0	284
13	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
14	Grundlagen der Thermodynamik für Studierende der Chemie. <i>Essentials</i> , 2017, , .	0.1	0