

## List of Publications by Year in descending order

Source: <https://exaly.com/author-pdf/1149451/publications.pdf>

Version: 2024-02-01

12  
papers

1,078  
citations

949033

11  
h-index

1336881

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g-index

12  
all docs

12  
docs citations

12  
times ranked

1437  
citing authors

#	ARTICLE	IF	CITATIONS
1	Efficient implementation of atom-density representations. Journal of Chemical Physics, 2021, 154, 114109.	1.2	32
2	Physics-Inspired Structural Representations for Molecules and Materials. Chemical Reviews, 2021, 121, 9759-9815.	23.0	247
3	Optimal radial basis for density-based atomic representations. Journal of Chemical Physics, 2021, 155, 104106.	1.2	17
4	Atom-density representations for machine learning. Journal of Chemical Physics, 2019, 150, 154110.	1.2	120
5	Data Science Based Mg Corrosion Engineering. Frontiers in Materials, 2019, 6, .	1.2	34
6	Machine Learning at the Atomic Scale. Chimia, 2019, 73, 972.	0.3	4
7	Fast and Accurate Uncertainty Estimation in Chemical Machine Learning. Journal of Chemical Theory and Computation, 2019, 15, 906-915.	2.3	102
8	Machine learning for the structureâ€“energyâ€“property landscapes of molecular crystals. Chemical Science, 2018, 9, 1289-1300.	3.7	153
9	Feature optimization for atomistic machine learning yields a data-driven construction of the periodic table of the elements. Physical Chemistry Chemical Physics, 2018, 20, 29661-29668.	1.3	88
10	Chemical shifts in molecular solids by machine learning. Nature Communications, 2018, 9, 4501.	5.8	170
11	Mapping and classifying molecules from a high-throughput structural database. Journal of Cheminformatics, 2017, 9, 6.	2.8	28
12	The GBS code for tokamak scrape-off layer simulations. Journal of Computational Physics, 2016, 315, 388-408.	1.9	83