## **Eleftherios Lambros**

## List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

10<br/>papers104<br/>citations7<br/>h-index10<br/>g-index12<br/>ext. papers182<br/>ext. citations6.7<br/>avg, IF3.33<br/>L-index

#	Paper	IF	Citations
10	Density functional theory of waterwith the machine-learned DM21 functional <i>Journal of Chemical Physics</i> , <b>2022</b> , 156, 161103	3.9	2
9	Anomalies and Local Structure of Liquid Water from Boiling to the Supercooled Regime as Predicted by the Many-Body MB-pol Model <i>Journal of Physical Chemistry Letters</i> , <b>2022</b> , 3652-3658	6.4	4
8	Elevating density functional theory to chemical accuracy for water simulations through a density-corrected many-body formalism. <i>Nature Communications</i> , <b>2021</b> , 12, 6359	17.4	9
7	Highly Accurate Many-Body Potentials for Simulations of NO in Water: Benchmarks, Development, and Validation. <i>Journal of Chemical Theory and Computation</i> , <b>2021</b> , 17, 3931-3945	6.4	5
6	Assessing the Accuracy of the SCAN Functional for Water through a Many-Body Analysis of the Adiabatic Connection Formula. <i>Journal of Chemical Theory and Computation</i> , <b>2021</b> , 17, 3739-3749	6.4	7
5	General Many-Body Framework for Data-Driven Potentials with Arbitrary Quantum Mechanical Accuracy: Water as a Case Study. <i>Journal of Chemical Theory and Computation</i> , <b>2021</b> , 17, 5635-5650	6.4	10
4	A Many-Body, Fully Polarizable Approach to QM/MM Simulations. <i>Journal of Chemical Theory and Computation</i> , <b>2020</b> , 16, 7462-7472	6.4	7
3	How good are polarizable and flexible models for water: Insights from a many-body perspective Journal of Chemical Physics, <b>2020</b> , 153, 060901	3.9	14
2	Low-order many-body interactions determine the local structure of liquid water. <i>Chemical Science</i> , <b>2019</b> , 10, 8211-8218	9.4	29
1	Modeling Membrane Protein-Ligand Binding Interactions: The Human Purinergic Platelet Receptor. <i>Journal of Physical Chemistry B</i> , <b>2016</b> , 120, 12293-12304	3.4	14