

# Thomas Dannenhoffer-Lafage

## List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	A Data-Driven Hydrophobicity Scale for Predicting Liquid-Liquid Phase Separation of Proteins. <i>Journal of Physical Chemistry B</i> , 2021, 125, 4046-4056.	2.6	71
2	Reactive Coarse-Grained Molecular Dynamics. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 2541-2549.	5.3	11
3	Compatible observable decompositions for coarse-grained representations of real molecular systems. <i>Journal of Chemical Physics</i> , 2019, 151, 134115.	3.0	16
4	Systematic Coarse-Grained Lipid Force Fields with Semiexplicit Solvation via Virtual Sites. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 2087-2100.	5.3	26
5	Multiconfigurational Coarse-Grained Molecular Dynamics. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 3306-3315.	5.3	22
6	Extending the range and physical accuracy of coarse-grained models: Order parameter dependent interactions. <i>Journal of Chemical Physics</i> , 2017, 147, 044113.	3.0	53
7	Coarse-Grained Directed Simulation. <i>Journal of Chemical Theory and Computation</i> , 2017, 13, 4593-4603.	5.3	11
8	A Direct Method for Incorporating Experimental Data into Multiscale Coarse-Grained Models. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 2144-2153.	5.3	29