

# Daniel Graf

## List of Publications by Year in descending order

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8  
papers

118  
citations

1478505

6  
h-index

1588992

8  
g-index

8  
all docs

8  
docs citations

8  
times ranked

46  
citing authors

#	ARTICLE	IF	CITATIONS
1	Benchmarking the Accuracy of the Direct Random Phase Approximation and ĩf-Functionals for NMR Shieldings. <i>Journal of Chemical Theory and Computation</i> , 2022, 18, 192-205.	5.3	3
2	Lagrangian-Based Minimal-Overhead Batching Scheme for the Efficient Integral-Direct Evaluation of the RPA Correlation Energy. <i>Journal of Chemical Theory and Computation</i> , 2021, 17, 5623-5634.	5.3	7
3	Efficient Reduced-Scaling Second-Order MÅller-Plisset Perturbation Theory with Cholesky-Decomposed Densities and an Attenuated Coulomb Metric. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 6856-6868.	5.3	14
4	Range-Separated Density-Functional Theory in Combination with the Random Phase Approximation: An Accuracy Benchmark. <i>Journal of Chemical Theory and Computation</i> , 2020, 16, 2985-2994.	5.3	10
5	A range-separated generalized Kohn-Sham method including a long-range nonlocal random phase approximation correlation potential. <i>Journal of Chemical Physics</i> , 2020, 153, 244118.	3.0	7
6	Low-Scaling Self-Consistent Minimization of a Density Matrix Based Random Phase Approximation Method in the Atomic Orbital Space. <i>Journal of Chemical Theory and Computation</i> , 2019, 15, 4468-4477.	5.3	17
7	Accurate and Efficient Parallel Implementation of an Effective Linear-Scaling Direct Random Phase Approximation Method. <i>Journal of Chemical Theory and Computation</i> , 2018, 14, 2505-2515.	5.3	35
8	Efficient calculation of beyond RPA correlation energies in the dielectric matrix formalism. <i>Journal of Chemical Physics</i> , 2018, 148, 204104.	3.0	25