Janus J Eriksen

List of Publications by Year
in descending order

31 papers	2,304 citations	516710 16 h-index	477307 29 g-index
31 all docs	31 docs citations	31 times ranked	2392 citing authors

#	Article	IF	Citations
1	Electronic excitations through the prism of mean-field decomposition techniques. Journal of Chemical Physics, 2022, 156, 061101.	3.0	4
2	Incremental treatments of the full configuration interaction problem. Wiley Interdisciplinary Reviews: Computational Molecular Science, 2021, 11, e1525.	14.6	10
3	Decomposed Mean-Field Simulations of Local Properties in Condensed Phases. Journal of Physical Chemistry Letters, 2021, 12, 6048-6055.	4.6	4
4	The Shape of Full Configuration Interaction to Come. Journal of Physical Chemistry Letters, 2021, 12, 418-432.	4.6	29
5	Foreword: Prof. Gauss Festschrift. Molecular Physics, 2020, 118, e1817247.	1.7	0
6	Ground and excited state first-order properties in many-body expanded full configuration interaction theory. Journal of Chemical Physics, 2020, 153, 154107.	3.0	14
7	The Ground State Electronic Energy of Benzene. Journal of Physical Chemistry Letters, 2020, 11, 8922-8929.	4.6	90
8	Recent developments in the P<scp>y</scp>SCF program package. Journal of Chemical Physics, 2020, 153, 024109.	3.0	388
9	Mean-field density matrix decompositions. Journal of Chemical Physics, 2020, 153, 214109.	3.0	10
10	Many-Body Expanded Full Configuration Interaction. II. Strongly Correlated Regime. Journal of Chemical Theory and Computation, 2019, 15, 4873-4884.	5.3	38
11	Generalized Many-Body Expanded Full Configuration Interaction Theory. Journal of Physical Chemistry Letters, 2019, 10, 7910-7915.	4.6	30
12	Many-Body Expanded Full Configuration Interaction. I. Weakly Correlated Regime. Journal of Chemical Theory and Computation, 2018, 14, 5180-5191.	5.3	45
13	Efficient and portable acceleration of quantum chemical many-body methods in mixed floating point precision using OpenACC compiler directives. Molecular Physics, 2017, 115, 2086-2101.	1.7	18
14	Virtual Orbital Many-Body Expansions: A Possible Route towards the Full Configuration Interaction Limit. Journal of Physical Chemistry Letters, 2017, 8, 4633-4639.	4.6	60
15	Massively parallel and linear-scaling algorithm for second-order MÃ¸llerâ€“Plesset perturbation theory applied to the study of supramolecular wires. Computer Physics Communications, 2017, 212, 152-160.	7.5	16
16	Incrementally accelerating the RI-MP2 correlated method of electronic structure theory using OpenACC compiler directives. , 2017, , 241-265.		1
17	Convergence of coupled cluster perturbation theory. Journal of Chemical Physics, 2016, 145, 224104.	3.0	6
18	Assessment of the accuracy of coupled cluster perturbation theory for open-shell systems. I. Triples expansions. Journal of Chemical Physics, 2016, 144, 194102.	3.0	16

#	Article	IF	Citations
19	A view on coupled cluster perturbation theory using a bivariational Lagrangian formulation. Journal of Chemical Physics, 2016, 144, 064103.	3.0	18
20	Assessment of the accuracy of coupled cluster perturbation theory for open-shell systems. II. Quadruples expansions. Journal of Chemical Physics, 2016, 144, 194103.	3.0	6
21	Communication: The performance of non-iterative coupled cluster quadruples models. Journal of Chemical Physics, 2015, 143, 041101.	3.0	15
22	Linear-Scaling Coupled Cluster with Perturbative Triple Excitations: The Divideâ€“Expandâ€“Consolidate CCSD(T) Model. Journal of Chemical Theory and Computation, 2015, 11, 2984-2993.	5.3	77
23	On the convergence of perturbative coupled cluster triples expansions: Error cancellations in the CCSD(T) model and the importance of amplitude relaxation. Journal of Chemical Physics, 2015, 142, 014102.	3.0	23
24	The same number of optimized parameters scheme for determining intermolecular interaction energies. Journal of Chemical Physics, 2015, 142, 114116.	3.0	12
25	A Lagrangian framework for deriving triples and quadruples corrections to the CCSD energy. Journal of Chemical Physics, 2014, 140, 064108.	3.0	40
26	Equation-of-motion coupled cluster perturbation theory revisited. Journal of Chemical Physics, 2014, 140, 174114.	3.0	17
27	The Second-Order Polarization Propagator Approximation (SOPPA) method coupled to the polarizable continuum model. Computational and Theoretical Chemistry, 2014, 1040-1041, 54-60.	2.5	9
28	The <scp>D</scp>alton quantum chemistry program system. Wiley Interdisciplinary Reviews: Computational Molecular Science, 2014, 4, 269-284.	14.6	1,166
29	Failures of TDDFT in describing the lowest intramolecular charge-transfer excitation in para-nitroaniline. Molecular Physics, 2013, 111, 1235-1248.	1.7	79
30	On the importance of excited state dynamic response electron correlation in polarizable embedding methods. Journal of Computational Chemistry, 2012, 33, 2012-2022.	3.3	38
31	Computational protocols for prediction of solute NMR relative chemical shifts. A case study of <scp>L</scp>â€tryptophan in aqueous solution. Journal of Computational Chemistry, 2011, 32, 2853-2864.	3.3	25