Lucas O Wagner

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9 papers 335 ph-index 9 g-index

9 as 383 as 3.9 avg, IF L-index

| # | Paper | IF | Citations |
|---|--|-----|-----------|
| 9 | DFT in a nutshell. International Journal of Quantum Chemistry, 2013, 113, 96-101 | 2.1 | 103 |
| 8 | Reference electronic structure calculations in one dimension. <i>Physical Chemistry Chemical Physics</i> , 2012 , 14, 8581-90 | 3.6 | 52 |
| 7 | Guaranteed convergence of the Kohn-Sham equations. <i>Physical Review Letters</i> , 2013 , 111, 093003 | 7.4 | 35 |
| 6 | Kohn-Sham calculations with the exact functional. <i>Physical Review B</i> , 2014 , 90, | 3.3 | 29 |
| 5 | Exchange-correlation functionals from the strong interaction limit of DFT: applications to model chemical systems. <i>Physical Chemistry Chemical Physics</i> , 2014 , 16, 14551-8 | 3.6 | 26 |
| 4 | Hydrogen Molecule Dissociation Curve with Functionals Based on the Strictly Correlated Regime. <i>Journal of Chemical Theory and Computation</i> , 2015 , 11, 3153-62 | 6.4 | 24 |
| 3 | Interpolated energy densities, correlation indicators and lower bounds from approximations to the strong coupling limit of DFT. <i>Physical Chemistry Chemical Physics</i> , 2017 , 19, 6169-6183 | 3.6 | 23 |
| 2 | One-dimensional mimicking of electronic structure: The case for exponentials. <i>Physical Review B</i> , 2015 , 91, | 3.3 | 22 |
| 1 | Electron avoidance: A nonlocal radius for strong correlation. <i>Physical Review A</i> , 2014 , 90, | 2.6 | 21 |