

# Towards an exact description of electronic wavefunction

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Citation Report

#	ARTICLE	IF	CITATIONS
3	Incrementally Corrected Periodic Local MP2 Calculations: I. The Cohesive Energy of Molecular Crystals. <i>Journal of Chemical Theory and Computation</i> , 2013, 9, 5590-5598.	2.3	45
4	Quantum Monte Carlo with coupled-cluster wave functions. <i>Physical Review B</i> , 2013, 88, .	1.1	24
5	Wave function methods for fractional electrons. <i>Journal of Chemical Physics</i> , 2013, 139, 074107.	1.2	19
6	Explicitly correlated plane waves: Accelerating convergence in periodic wavefunction expansions. <i>Journal of Chemical Physics</i> , 2013, 139, 084112.	1.2	62
7	Computing molecular correlation energies with guaranteed precision. <i>Journal of Chemical Physics</i> , 2013, 139, 114106.	1.2	31
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14	Linear-scaling explicitly correlated treatment of solids: Periodic local MP2-F12 method. <i>Journal of Chemical Physics</i> , 2013, 139, 194101.	1.2	42
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20	Generalizing the self-healing diffusion Monte Carlo approach to finite temperature: A path for the optimization of low-energy many-body bases. <i>Journal of Chemical Physics</i> , 2014, 140, 074103.	1.2	2
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24	Sign problem in full configuration interaction quantum Monte Carlo: Linear and sublinear representation regimes for the exact wave function. <i>Physical Review B</i> , 2014, 90, .	1.1	31
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