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| **M** | Bates, Jefferson  | UC Irvine | SOSEX and Beyond: Perturbative Corrections to RPA |  |
| **M** | Booth, George  | Princeton University  | Full Configuration Interaction Quantum Monte Carlo |  |
| **M** | Burow, Asbjorn  | UC Irvine | Investigation of pi-pi stacking using RPA correlation |  |
| **M** | Chakraborty, Arindam  | Syracuse University  | Development of electron-hole density functional theory for investigation of optical properties of quantum dots" |  |
| **M** | Eliav, Ephraim  | Tel Aviv University  | Intermediate Hamiltonian Hilbert-space coupled cluster method: theory and pilot applications |  |
| **M** | Eshuis, Henk  | UC Irvine | Molecular correlation energies from the random phase approximation: fast implementation, applications and basis set convergence |  |
| **M** | Ettenhuber, Patrick  | Aarhus University  | Integral direct and memory conservative CCSD residual algorithm |  |
| **M** | Gora, Urszula  | University of Silesia  | SAMBA of water clusters |  |
| **M** | Henderson, Tom  | Rice Univeristy  | CC Theory on Projected HF wavefunctions |  |
| **M** | Hoeyvik, Ida-Marie  | Aarhus University  | Local Hartree-Fock orbitals for large molecular systems using trust-region optimization |  |
| **M** | Hua, Shugui  | Nanjing University  | Cooperativity in Long α- and 310-Helical Polyalanines: Both Electrostatic and van der Waals Interactions are Essential |  |
| **M** | Jacob, Christoph  | Karlsruhe Institute of Tech  | Spin in Density-Functional Theory |  |
| **M** | Jacobson, Leif  | Yale University  | Construction of Diabatic Hamiltonians for Non-Adiabatic Dynamics from the Single Particle Greens Function |  |
| **M** | Jensen, Stig Rune  | University of Tromsø  | Multiwavelets in chemistry |  |
| **M** | Knizia, Gerald  | Princeton University  | The atomic valence active space |  |
| **W** | Kristensen, Kasper  | Aarhus University  | Large-scale MP2 calcuations using the Divide-Expand-Consolidate method: Energy and density for insulin |  |
| **W** | Li, Wei  | Nanjing University  | A refined cluster-in-molecule local correlation approach for predicting the relative energies of large systems |  |
| **W** | Lyakh (Liakh), Dmitry  | University of Florida  | Does the disconnected nature of Lambda-equations present an obstacle for a linear-scaling implementation of analytic gradients of the coupled-cluster energy? |  |
| **W** | Meyer, Wilfried  | University of Kaiserslautern  | Triple substitutions in Coupled pair and Coupled Cluster Theories |  |
| **W** | Paiboonvorachat, Nattapong  | University of Oxford  | Electronic Structure of Porphyrin from DMRG calculation in HF space |  |
| **W** | Pham, Tuan Anh  | University of California Davis  | Calculations of quasiparticle energies from a spectral decomposition of the static dielectric matrix |  |
| **W** | Plessow, Philipp  | BASF SE  | Quasi-Second-Order Reaction Path Optimization |  |
| **W** | Podeszwa, Rafal  | University of Silesia  | Dispersion energies for nano-scale systems from coupled density response functions |  |
| **W** | Sharma, Sandeep | Princeton University  | DMRG: Exact calculations and challenging transition metal clusters |  |
| **W** | Staroverov, Viktor  | University of Western Ontario  | Advances in methodology and applications of model Kohn-Sham potentials |  |
| **W** | Vazquez Mayagoitia, Alvaro  | Argonne National Laboratory  | Modeling of Molecular Properties with Learning Machine |  |
| **W** | Watson, Mark | Princeton University  | Excited states of butadiene to chemical accuracy: reconciling theory with experiment |  |
| **W** | Yanai, Takeshi  | Institute for Molecular Science  | Canonical transcorrelated theory with projected Slater-type geminals |  |
| **W** | Yang, Jun | Princeton University  | The orbital-specific-virtual local coupled-cluster method |  |
| **W** | Yun, Jeonghun  | Pohang University of S & T  | Design of high-resolution DNA sequencing from molecular dynamics and quantum transport simulations |  |
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