Self-interaction-free electric dipole polarizabilities for a Fermi-LÃ¶wdin self-interaction correction

Physical Review A

100,

DOI: 10.1103/physreva.100.012505

Citation Report

#	Article	IF	Citations
1	Fermi-LÃ¶wdin orbital self-interaction correction using the strongly constrained and appropriately normed meta-GGA functional. Journal of Chemical Physics, 2019, 151, 154105.	1.2	38
2	A step in the direction of resolving the paradox of Perdew-Zunger self-interaction correction. Journal of Chemical Physics, 2019, 151, 214108.	1.2	56
3	Importance of self-interaction-error removal in density functional calculations on water cluster anions. Physical Chemistry Chemical Physics, 2020, 22, 3789-3799.	1.3	32
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5	Assessing the effect of regularization on the molecular properties predicted by SCAN and self-interaction corrected SCAN meta-GGA. Physical Chemistry Chemical Physics, 2020, 22, 18060-18070.	1.3	6
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8	PyFLOSIC: Python-based Fermiâ€“LÃ¶wdin orbital self-interaction correction. Journal of Chemical Physics, 2020, 153, 084104.	1.2	17
9	Improvements in the orbitalwise scaling down of Perdewâ€“Zunger self-interaction correction in many-electron regions. Journal of Chemical Physics, 2020, 152, 174112.	1.2	23
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15	Self-interaction correction in waterâ€“ion clusters. Journal of Chemical Physics, 2021, 154, 094302.	1.2	16
16	Static dipole polarizabilities of polyacenes using self-interaction-corrected density functional approximations. Journal of Chemical Physics, 2021, 154, 114305.	1.2	12
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19	General Many-Body Framework for Data-Driven Potentials with Arbitrary Quantum Mechanical Accuracy: Water as a Case Study. Journal of Chemical Theory and Computation, 2021, 17, 5635-5650.	2.3	28
20	How well do self-interaction corrections repair the overestimation of static polarizabilities in density functional calculations?. Physical Chemistry Chemical Physics, 2021, 23, 18678-18685.	1.3	14
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27	Downward quantum learning from element 118: Automated generation of Fermiâ€“LÃ¶wdin orbitals for all atoms. Journal of Chemical Physics, 2023, 158, .	1.2	5