

Self-consistent molecular orbital methods. XX. A basis

Journal of Chemical Physics

72, 650-654

DOI: 10.1063/1.438955

Citation Report

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5339	Understanding Conjugation and Hyperconjugation from Electronic Delocalization Measures. <i>Journal of Physical Chemistry A</i> , 2011, 115, 13104-13113.	1.1	30
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5341	Fourier transform infrared spectroscopic and theoretical study of water interactions with glycine and its N-methylated derivatives. <i>Journal of Chemical Physics</i> , 2011, 134, 115104.	1.2	32
5342	QM/MM calculation of protein magnetic shielding tensors with generalized hybrid-orbital method: A GIAO approach. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 14490.	1.3	3
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8915	Small Molecule Activation by Constrained Phosphorus Compounds: Insights from Theory. <i>Inorganic Chemistry</i> , 2016, 55, 558-565.	1.9	27
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10206	Computational screening and molecular design of anthracene-based semiconductors. <i>Organic Electronics</i> , 2018, 61, 87-95.	1.4	5
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10287	Structure and Magnetization Dynamics of Dy ^{III} -Fe and Dy ^{III} -Ru Bonded Complexes. <i>Angewandte Chemie - International Edition</i> , 2018, 57, 8144-8148.	7.2	38
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13441	Coordination-Induced Bond Weakening. <i>Chemical Reviews</i> , 2022, 122, 13447-13477.	23.0	22
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