

# Yuxiang Mo

## List of Publications by Year in descending order

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docs citations

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214  
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#	ARTICLE	IF	CITATIONS
1	First-principles study of the binding energy between nanostructures and its scaling with system size. Physical Review B, 2018, 97, .	3.2	13
2	Accurate lattice geometrical parameters and bulk moduli from a semilocal density functional. AIP Advances, 2018, 8, .	1.3	15
3	Assessment of the Tao-Mo nonempirical semilocal density functional in applications to solids and surfaces. Physical Review B, 2017, 95, .	3.2	37
4	Accurate excitation energies of molecules and oligomers from a semilocal density functional. Journal of Chemical Physics, 2017, 146, 234102.	3.0	11
5	Performance of a nonempirical exchange functional from density matrix expansion: comparative study with different correlations. Physical Chemistry Chemical Physics, 2017, 19, 21707-21713.	2.8	20
6	Energetic Study of Clusters and Reaction Barrier Heights from Efficient Semilocal Density Functionals. Computation, 2017, 5, 27.	2.0	7
7	Performance of a nonempirical density functional on molecules and hydrogen-bonded complexes. Journal of Chemical Physics, 2016, 145, 234306.	3.0	25
8	Accurate Semilocal Density Functional for Condensed-Matter Physics and Quantum Chemistry. Physical Review Letters, 2016, 117, 073001.	7.8	124
9	Accurate van der Waals coefficients between fullerenes and fullerene-alkali atoms and clusters: Modified single-frequency approximation. Physical Review B, 2016, 94, .	3.2	12