J Ulises Reveles

List of Publications by Year in descending order

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	840776		940533	
16	782	11	16	
papers	citations	h-index	g-index	
16	16	16	694	
all docs	docs citations	times ranked	citing authors	

#	Article	IF	CITATIONS
1	Designer magnetic superatoms. Nature Chemistry, 2009, 1, 310-315.	13.6	223
2	Multiple valence superatoms. Proceedings of the National Academy of Sciences of the United States of America, 2006, 103, 18405-18410.	7.1	197
3	Geometry optimization in density functional methods. Journal of Computational Chemistry, 2004, 25, 1109-1116.	3.3	131
4	Ti ₄ - and Ni ₄ -Doped Defective Graphene Nanoplatelets as Efficient Materials for Hydrogen Storage. Journal of Physical Chemistry C, 2016, 120, 5001-5009.	3.1	68
5	H ₂ O Nucleation around Au ⁺ . Journal of the American Chemical Society, 2007, 129, 15565-15571.	13.7	30
6	Highly efficient (Cs8V) superatom-based spin-polarizer. Applied Physics Letters, 2009, 95, .	3.3	26
7	Magnetism of electrons in atoms and superatoms. Journal of Applied Physics, 2012, 112, 064313.	2.5	17
8	Hydrogen storage in bimetallic Ti–Al sub-nanoclusters supported on graphene. Physical Chemistry Chemical Physics, 2017, 19, 21174-21184.	2.8	16
9	Structural changes of Pd13 upon charging and oxidation/reduction. Journal of Chemical Physics, 2012, 136, 114505.	3.0	13
10	Influence of Interligand Interactions and Core-Charge Distribution on Gold Cluster Stability: Enthalpy Versus Entropy. Journal of Physical Chemistry C, 2019, 123, 24899-24911.	3.1	13
11	Electronic and Structural Properties of C ₆₀ and Sc ₃ N@C ₈₀ Supported on Graphene Nanoflakes. Journal of Physical Chemistry C, 2016, 120, 26083-26092.	3.1	11
12	Ion Mobility Spectrometry Characterization of the Intermediate Hydrogen-Containing Gold Cluster Au ₇ (PPh ₃) ₇ H ₅ ²⁺ . Journal of Physical Chemistry Letters, 2021, 12, 2502-2508.	4.6	11
13	Equivalent delocalized internal coordinates. Computational and Theoretical Chemistry, 2006, 762, 171-178.	1.5	10
14	Water inhibits CO oxidation on gold cations in the gas phase. Structures and binding energies of the sequential addition of CO, H ₂ O, O ₂ , and N ₂ onto Au ⁺ . Physical Chemistry Chemical Physics, 2016, 18, 28606-28616.	2.8	7
15	Electronic and Structural Study of Zn _{<i>x</i>} S _{<i>x</i>} [<i>x</i> [<i>x [<i>x [<i>x] [<i>x] [<i>x] [<i>x] [<i>x]]]</i></i></i></i></i></i></i>	2.5	6
16	Ab Initio Molecular Dynamics Investigation of the Electronic and Structural Stability of Anionic O ₂ _{3€"} (H ₂ O) _{<i>n</i>} , <i>n</i> = 1â€"16 Clusters. Journal of Physical Chemistry A, 2019, 123, 7528-7535.	2.5	3