

# Nasim Hassani

## List of Publications by Year in descending order

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20  
papers

316  
citations

1163117

8  
h-index

888059

17  
g-index

21  
all docs

21  
docs citations

21  
times ranked

157  
citing authors

#	ARTICLE	IF	CITATIONS
1	Catalytic properties of cyclo-carbon clusters: An investigation on $\sigma$ - $\pi$ activation and CO oxidation. Surface Science, 2022, 720, 122050.	1.9	6
2	Trilayer Metal-Organic Frameworks as Multifunctional Electrocatalysts for Energy Conversion and Storage Applications. Journal of the American Chemical Society, 2022, 144, 3411-3428.	13.7	142
3	Theoretical investigation of the interaction between the metal phthalocyanine [MPC] <sub>a</sub> (M = Sc, Ti, and Tj) and CO. Journal of Physical Chemistry C, 2021, 125, 110086.	1.2	1
4	Theoretical perspective on CO oxidation over small cobalt oxide clusters. Chemical Physics Letters, 2021, 767, 138361.	2.6	8
5	Oscillation in the electrical conductivity of a thick graphene oxide membrane. Journal of Applied Physics, 2021, 129, 235105.	2.5	1
6	The inhibition performance of quinoa seed on corrosion behavior of carbon steel in the HCl solution; theoretical and experimental evaluations. Journal of Molecular Liquids, 2021, 335, 116183.	4.9	26
7	Breakdown of Universal Scaling for Nanometer-Sized Bubbles in Graphene. Nano Letters, 2021, 21, 8103-8110.	9.1	23
8	The interaction between atomic-scale pores and particles. Journal of Physics Condensed Matter, 2021, 34, .	1.8	0
9	Evaluating gas permeance through graphene nanopores and porous 2D-membranes: A generalized approach. Carbon Trends, 2021, 5, 100086.	3.0	2
10	Gas Permeability and Selectivity of a Porous WS <sub>2</sub> Monolayer. Journal of Physical Chemistry C, 2021, 125, 25055-25066.	3.1	11
11	CO oxidation by linear oxocarbon chains O-C-O (n=10, x=1, 2): A theoretical study. Chemical Physics, 2020, 530, 110652.	1.9	3
12	Kinetics and Mechanism of the NH <sub>3</sub> + SO <sub>3</sub> Reaction: A Theoretical Approach. Journal of Physical Chemistry A, 2020, 124, 6585-6600.	2.5	3
13	Gas flow through atomic-scale apertures. Science Advances, 2020, 6, .	10.3	22
14	Exploring the adsorption and sensing behavior of the M-N <sub>x</sub> -B <sub>36</sub> -x (M=Fe, Ni, and Cu; x=0, 3) bowl-shaped structures upon CO, NO, O <sub>2</sub> , and N <sub>2</sub> molecules: A first-principles study. Physica E: Low-Dimensional Systems and Nanostructures, 2020, 124, 114242.	2.7	3
15	The reaction mechanism of the hydration of ethylene over the Corrole (M=Al, and Ga) complexes: A theoretical approach. Computational and Theoretical Chemistry, 2020, 1177, 112766.	2.5	3
16	C <sub>20</sub> fullerene and its boron- and nitrogen-doped counterparts as an efficient catalyst for CO oxidation. Molecular Physics, 2020, 118, e1766708.	1.7	3
17	The interplay between structural perfectness and CO oxidation catalysis on aluminum, phosphorous and silicon complexes of corroles. Physical Chemistry Chemical Physics, 2019, 21, 7661-7674.	2.8	9
18	Cu <sub>2</sub> O/TiO <sub>2</sub> nanoparticles as visible light photocatalysts concerning C(sp <sup>2</sup> )=P bond formation. Catalysis Science and Technology, 2018, 8, 4044-4051.	4.1	41

#	ARTICLE	IF	CITATIONS
19	Catalytic activity of corrole complexes with post-transition elements for the oxidation of carbon monoxide: a first-principles study. <i>New Journal of Chemistry</i> , 2018, 42, 12632-12643.	2.8	7
20	NO oxidation catalyzed by Ir <sub>4</sub> -based nanoclusters: the role of alloying on the catalytic activity. <i>Theoretical Chemistry Accounts</i> , 2017, 136, 1.	1.4	1