

# Maedeh Kamel

## List of Publications by Year in descending order

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Version: 2024-02-01

7  
papers

197  
citations

1478505

6  
h-index

1872680

6  
g-index

7  
all docs

7  
docs citations

7  
times ranked

137  
citing authors

#	ARTICLE	IF	CITATIONS
1	Assessment of the adsorption mechanism of Flutamide anticancer drug on the functionalized single-walled carbon nanotube surface as a drug delivery vehicle: An alternative theoretical approach based on DFT and MD. <i>Applied Surface Science</i> , 2018, 434, 492-503.	6.1	87
2	Theoretical study of solvent and co-solvent effects on the interaction of Flutamide anticancer drug with Carbon nanotube as a drug delivery system. <i>Journal of Molecular Liquids</i> , 2017, 248, 490-500.	4.9	60
3	Theoretical elucidation of the amino acid interaction with graphene and functionalized graphene nanosheets: insights from DFT calculation and MD simulation. <i>Amino Acids</i> , 2020, 52, 1465-1478.	2.7	22
4	Density functional theory study towards investigating the adsorption properties of the $\hat{\text{I}}^3\text{-Fe}_2\text{O}_3$ nanoparticles as a nanocarrier for delivery of Flutamide anticancer drug. <i>Adsorption</i> , 2020, 26, 925-939.	3.0	12
5	Selective adsorption and dissociation of NO, NO <sub>2</sub> , and N <sub>2</sub> O molecules on Si-doped haeckelite boron nitride nanotube: an investigation for sensitive molecular sensors and catalysts. <i>Journal of Molecular Modeling</i> , 2022, 28, 6.	1.8	9
6	Probing and comparison of graphene, boron nitride and boron carbide nanosheets for Flutamide adsorption: A DFT computational study. <i>Journal of Molecular Liquids</i> , 2021, 343, 117487.	4.9	7
7	Understanding the role of hydrogen bonds in destruction of DNA by screening interactions of Flutamide anticancer drug with nucleotides bases: DFT perspective, MD simulation and free energy calculation. <i>Adsorption</i> , 2020, 26, 491-508.	3.0	0