Ahmed A Elhenawy

List of Publications by Year in descending order

Source: https://exaly.com/author-pdf/8566414/publications.pdf

Version: 2024-02-01

516710 642732 32 614 16 23 citations g-index h-index papers 33 33 33 408 docs citations times ranked citing authors all docs

#	Article	IF	Citations
1	Synthesis, Biological Evaluation, and Molecular Docking of New Benzimidazole-1,2,3-Triazole Hybrids as Antibacterial and Antitumor Agents. Polycyclic Aromatic Compounds, 2023, 43, 3380-3391.	2.6	16
2	Synthesis of 1,4-dihydropyrano[2,3-c]pyrazole derivatives and exploring molecular and cytotoxic properties based on DFT and molecular docking studies. Journal of Molecular Structure, 2022, 1249, 131555.	3.6	18
3	Experimental and theoretical investigation for 6-Morpholinosulfonylquinoxalin-2(1H)-one and its haydrazone derivate: Synthesis, characterization, tautomerization and antimicrobial evaluation. Journal of Molecular Structure, 2022, 1247, 131314.	3.6	34
4	Therapeutic strategies for Covid-19 based on molecular docking and dynamic studies to the ACE-2 receptors, Furin, and viral spike proteins. Journal of Biomolecular Structure and Dynamics, 2022, 40, 13291-13309.	3. 5	24
5	Synthesis, antiproliferative, docking and DFT studies of benzimidazole derivatives as EGFR inhibitors. Journal of Molecular Structure, 2022, 1253, 132265.	3.6	29
6	Investigation of the Anticancer Effect of \hat{l}_{\pm} -Aminophosphonates and Arylidine Derivatives of 3-Acetyl-1-aminoquinolin-2(1H)-one on the DMBA Model of Breast Cancer in Albino Rats with In Silico Prediction of Their Thymidylate Synthase Inhibitory Effect. Molecules, 2022, 27, 756.	3.8	10
7	Discovery of novel rigid analogs of 2-naphthol with potent anticancer activity through multi-target topoisomerase I & Damp; II and tyrosine kinase receptor EGFR & DEGFR-2 inhibition mechanism. Chemico-Biological Interactions, 2022, 355, 109838.	4.0	9
8	Theoretical and experimental evaluation of the anticorrosion properties of new Coumarin's derivatives. AEJ - Alexandria Engineering Journal, 2022, 61, 6937-6948.	6.4	8
9	Myco-Synthesized Molluscicidal and Larvicidal Selenium Nanoparticles: A New Strategy to Control Biomphalaria alexandrina Snails and Larvae of Schistosoma mansoni with an In Silico Study on Induced Oxidative Stress. Journal of Fungi (Basel, Switzerland), 2022, 8, 262.	3.5	27
10	Synthesis, cytotoxic activity, crystal structure, DFT, molecular docking study of some heterocyclic compounds incorporating benzo[f]chromene moieties. Journal of Molecular Structure, 2022, 1260, 132829.	3.6	5
11	Discovery Potent of Thiazolidinedione Derivatives as Antioxidant, α-Amylase Inhibitor, and Antidiabetic Agent. Biomedicines, 2022, 10, 24.	3.2	20
12	Extensive Study of DFT-Quantum Calculations Based QSAR Modeling of Fused 1,2,4-Triazine Derivatives Revealed Potent CYP1A1 Inhibitors. Journal of Computational Biophysics and Chemistry, 2022, 21, 741-758.	1.7	3
13	The Crystal Structure of 2-Amino-4-(2,3-Dichlorophenyl)-6-Methoxy-4H-Benzo[h]chromene-3-Carbonitrile: Antitumor and Tyrosine Kinase Receptor Inhibition Mechanism Studies. Crystals, 2022, 12, 737.	2.2	5
14	The Crystal Structure of 3-Amino-1-(4-Chlorophenyl)-9-Methoxy-1H-Benzo[f]Chromene-2-Carbonitrile: Antimicrobial Activity and Docking Studies. Crystals, 2022, 12, 982.	2.2	5
15	New carbazole-based organic dyes with different acceptors for dye-sensitized solar cells: Synthesis, characterization, dssc fabrications and density functional theory studies. Journal of Molecular Structure, 2021, 1225, 129297.	3.6	52
16	Synthesis and characterization of MnII and CoII complexes with poly (vinyl alcohol-nicotinic acid) for photocatalytic degradation of Indigo carmine dye. Inorganic Chemistry Communication, 2021, 124, 108360.	3.9	12
17	Synthesis, Cytotoxic Activity, Crystal Structure, DFT Studies and Molecular Docking of 3-Amino-1-(2,5-dichlorophenyl)-8-methoxy-1H-benzo[f]chromene-2-carbonitrile. Crystals, 2021, 11, 184.	2.2	27
18	Nα-1, 3-Benzenedicarbonyl-Bis-(Amino Acid) and Dipeptide Candidates: Synthesis, Cytotoxic, Antimicrobial and Molecular Docking Investigation. Drug Design, Development and Therapy, 2021, Volume 15, 1315-1332.	4.3	11

#	ARTICLE	IF	Citations
19	Design, synthesis and molecular docking studies of thymol based 1,2,3-triazole hybrids as thymidylate synthase inhibitors and apoptosis inducers against breast cancer cells. Bioorganic and Medicinal Chemistry, 2021, 38, 116136.	3.0	33
20	Preparation, Antimicrobial Activity and Docking Study of Vanadium Mixed Ligand Complexes Containing 4-Amino-5-hydrazinyl-4H-1,2,4-triazole-3-thiol and Aminophenol Derivatives. Processes, 2021, 9, 1008.	2.8	11
21	Naproxen Based 1,3,4-Oxadiazole Derivatives as EGFR Inhibitors: Design, Synthesis, Anticancer, and Computational Studies. Pharmaceuticals, 2021, 14, 870.	3.8	32
22	Synthesis and Biological Evaluation of 1,2,3-Triazole Tethered Thymol-1,3,4-Oxadiazole Derivatives as Anticancer and Antimicrobial Agents. Pharmaceuticals, 2021, 14, 866.	3.8	30
23	Structural and Spectroscopic Characteristics of Nill and Cull Complexes with Poly (Vinyl) Tj ETQq1 1 0.784314 rg 2021, 11, 1244.	gBT /Overl 2.2	ock 10 Tf 50 5 2
24	Glucosinolates with their Hydrolysis Products from Two Cruciferous Plants with Study of Antidiabetic Activity Based on Molecular Docking. Biomedical and Pharmacology Journal, 2021, 14, 2047-2062.	0.5	0
25	Novel coumarin amino acid derivatives: Design, synthesis, docking, absorption, distribution, metabolism, elimination, toxicity (ADMET), quantitative structure–activity relationship (QSAR) and anticancer studies. Materials Express, 2020, 10, 1375-1394.	0.5	8
26	Naproxenylamino acid derivatives: Design, synthesis, docking, QSAR and anti-inflammatory and analgesic activity. Biomedicine and Pharmacotherapy, 2019, 116, 109024.	5.6	11
27	<p>Synthesis, comparative docking, and pharmacological activity of naproxen amino acid derivatives as possible anti-inflammatory and analgesic agents</p> . Drug Design, Development and Therapy, 2019, Volume 13, 1773-1790.	4.3	30
28	Synthesis, molecular properties and comparative docking and QSAR of new 2-(7-hydroxy-2-oxo-2H-chromen-4-yl)acetic acid derivatives as possible anticancer agents. Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy, 2019, 218, 248-262.	3.9	26
29	Understanding the adsorption performance of two glycine derivatives as novel and environmentally safe anti-corrosion agents for copper in chloride solutions: experimental, DFT, and MC studies. RSC Advances, 2019, 9, 42120-42131.	3.6	47
30	Synthesis and characterization of some arylhydrazone ligand and its metal complexes and their potential application as flame retardant and antimicrobial additives in polyurethane for surface coating. Journal of Organometallic Chemistry, 2015, 791, 99-106.	1.8	35
31	Synthesis, characterization, biological activity, molecular modeling and docking studies of complexes 4-(4-hydroxy)-3-(2-pyrazine-2-carbonyl)hydrazonomethylphenyl-diazen-yl-benzenesulfonamide with manganese(II), cobalt(II), nickel(II), zinc(II) and cadmium(II). Journal of Molecular Structure, 2015, 1084, 352-367.	3.6	25
32	DFT, characterization and investigation of vibrational spectroscopy of 4-(4-hydroxy)-3-(2-pyrazine-2-carbonyl)hydrazonomethylphenyl-diazen-yl-benzenesulfonamide and its copper(II) complex. Journal of Molecular Structure, 2014, 1067, 94-103.	3.6	9