

Shazia Parveen

List of Publications by Year in descending order

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#	ARTICLE	IF	CITATIONS
1	Synthesis, Docking and Density Functional Theory Approaches on 1,3-Bis-3-(4-Chlorophenyl)-2,3-Dihydroquinazolin-4(1H)-on-2-Thioxopropane toward the Discovery of Dual Kinase Inhibitor. <i>Polycyclic Aromatic Compounds</i> , 2022, 42, 3736-3747.	1.4	10
2	2-(Alkylthio)-3-(Naphthalen-1-yl)Quinazolin-4(3H)-Ones: Ultrasonic Synthesis, DFT and Molecular Docking Aspects. <i>Polycyclic Aromatic Compounds</i> , 2022, 42, 4034-4048.	1.4	15
3	New quinoline-based BODIPYs as EGFR/VEGFR-2 inhibitors: Molecular docking, DFT and in vitro cytotoxicity on HeLa cells. <i>Journal of Molecular Structure</i> , 2022, 1247, 131312.	1.8	8
4	Furanone-functionalized benzothiazole derivatives: synthesis, in vitro cytotoxicity, ADME, and molecular docking studies. <i>Zeitschrift Fur Naturforschung - Section B Journal of Chemical Sciences</i> , 2022, 77, 41-53.	0.3	1
5	Platinum-based Cancer Chemotherapeutics: Recent Trends and Future Perspectives. <i>Current Chinese Science</i> , 2022, 2, 275-293.	0.2	2
6	UV-selective organic absorbers for the cosensitization of greenhouse-integrated dye-sensitized solar cells: synthesis and computational study. <i>RSC Advances</i> , 2022, 12, 11420-11435.	1.7	8
7	Molecular docking, DFT and antimicrobial studies of Cu(II) complex as topoisomerase I inhibitor. <i>Journal of Biomolecular Structure and Dynamics</i> , 2021, 39, 2092-2105.	2.0	13
8	Synthesis, in vitro cytotoxicity, ADME, and molecular docking studies of benzimidazole-bearing furanone derivatives. <i>Journal of the Chinese Chemical Society</i> , 2021, 68, 362-373.	0.8	12
9	Biophysical and theoretical investigation of benzo[c]coumarin functionalized Schiff base with human serum albumin. <i>Chemical Papers</i> , 2021, 75, 2339-2351.	1.0	8
10	Bioactive Tryptophan-Based Copper Complex with Auxiliary $\hat{2}$ -Carboline Spectacle Potential on Human Breast Cancer Cells: In Vitro and In Vivo Studies. <i>Molecules</i> , 2021, 26, 1606.	1.7	6
11	Phytoremediation: In situ Alternative for Pollutant Removal from Contaminated Natural Media: A Brief Review. <i>Biointerface Research in Applied Chemistry</i> , 2021, 12, 4945-4960.	1.0	4
12	Synthesis and Computational Characterization of Organic UV-Dyes for Cosensitization of Transparent Dye-Sensitized Solar Cells. <i>Molecules</i> , 2021, 26, 7336.	1.7	4
13	N-alkyl 2-pyridone versus O-alkyl 2-pyridol: Ultrasonic synthesis, DFT, docking studies and their antimicrobial evaluation. <i>Journal of Molecular Structure</i> , 2020, 1199, 126926.	1.8	21
14	A new chiral boron-dipyrromethene (BODIPY)-based fluorescent probe: molecular docking, DFT, antibacterial and antioxidant approaches. <i>Journal of Biomolecular Structure and Dynamics</i> , 2020, 38, 5429-5442.	2.0	34
15	Synthesis, Molecular Docking, and DFT Calculation of a Half-Strapped BODIPY as Potential EGFR Inhibitor**. <i>ChemistrySelect</i> , 2020, 5, 13163-13173.	0.7	5
16	Recent advances in anticancer ruthenium Schiff base complexes. <i>Applied Organometallic Chemistry</i> , 2020, 34, e5687.	1.7	49
17	Recent advancements in the anticancer potentials of phenylorganotin(IV) complexes. <i>Inorganica Chimica Acta</i> , 2020, 505, 119464.	1.2	18
18	Computational and molecular docking approaches of a New axially chiral BODIPY fluorescent dye. <i>Journal of Photochemistry and Photobiology A: Chemistry</i> , 2020, 395, 112508.	2.0	15

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19	Potential Exploration of Recent FDA-Approved Anticancer Drugs Against Models of SARS-CoV-2's Main Protease and Spike Glycoprotein: A Computational Study. <i>Biointerface Research in Applied Chemistry</i> , 2020, 11, 10059-10073.	1.0	3
20	Clinical developments of antitumor polymer therapeutics. <i>RSC Advances</i> , 2019, 9, 24699-24721.	1.7	47
21	Molecular Interaction of Amino Acid-Based Gemini Surfactant with Human Serum Albumin: Tensiometric, Spectroscopic, and Molecular Docking Study. <i>ACS Omega</i> , 2019, 4, 22152-22160.	1.6	14
22	Enantiomeric in vitro DNA binding, pBR322 DNA cleavage and molecular docking studies of chiral l- and d-ternary copper(II) complexes of histidine and picolinic acid. <i>Journal of Photochemistry and Photobiology B: Biology</i> , 2014, 130, 170-178.	1.7	40
23	Zinc(II) complexes of Pro-Gly and Pro-Leu dipeptides: Synthesis, characterization, in vitro DNA binding and cleavage studies. <i>Journal of Photochemistry and Photobiology B: Biology</i> , 2013, 126, 78-86.	1.7	11
24	Design and synthesis of (S)- and (R)-enantiomers of [4-(2-hydroxy-1-phenylethylimino)pent-2-yl]dimethyltin(IV) and 2,2-dimethyl-4-phenyl-1,3,2-oxazastannolidine: in vitro antitumor activity against human tumor cell lines and in vivo assay of (S)-enantiomers. <i>Dalton Transactions</i> , 2013, 42, 3390-3401.	1.6	35
25	Synthesis, characterization, biological studies (DNA binding, cleavage, antibacterial and) Tj ETQq1 1 0.784314 rgBT /Overlock 10 Tf 50 5 Photochemistry and Photobiology B: Biology, 2012, 114, 15-26.	1.7	123
26	Enantiomeric recognition of chiral L- and D-penicillamine Zinc(II) complexes: DNA binding behavior and cleavage studies. <i>RSC Advances</i> , 2012, 2, 6354.	1.7	24
27	Synthesis, characterization of Cu(II) and Zn(II) complexes of proline-glycine and proline-leucine tetrapeptides: In vitro DNA binding and cleavage studies. <i>Inorganica Chimica Acta</i> , 2012, 388, 1-10.	1.2	29
28	Molecular drug design, synthesis and crystal structure determination of Cu(II)-Sn(IV) heterobimetallic core: DNA binding and cleavage studies. <i>European Journal of Medicinal Chemistry</i> , 2012, 49, 141-150.	2.6	70
29	De novo design, synthesis and spectroscopic characterization of chiral benzimidazole-derived amino acid Zn(II) complexes: Development of tryptophan-derived specific hydrolytic DNA artificial nuclease agent. <i>Spectrochimica Acta - Part A: Molecular and Biomolecular Spectroscopy</i> , 2012, 85, 53-60.	2.0	36
30	In vitro binding studies of organotin(IV) complexes of 1,2-bis(1H-benzimidazol-2-yl)ethane-1,2-diol with CT-DNA and nucleotides (5'-GMP and 5'-TMP): Effect of the ancillary ligand on the binding propensity. <i>Journal of Organometallic Chemistry</i> , 2011, 696, 3836-3845.	0.8	9