

Samuel Kojo Kwofie

List of Publications by Year in descending order

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Version: 2024-02-01

44
papers

545
citations

687363

13
h-index

713466

21
g-index

44
all docs

44
docs citations

44
times ranked

553
citing authors

#	ARTICLE	IF	CITATIONS
1	HCVpro: Hepatitis C virus protein interaction database. <i>Infection, Genetics and Evolution</i> , 2011, 11, 1971-1977.	2.3	76
2	DDPC: Dragon Database of Genes associated with Prostate Cancer. <i>Nucleic Acids Research</i> , 2011, 39, D980-D985.	14.5	38
3	Cheminformatics-Based Identification of Potential Novel Anti-SARS-CoV-2 Natural Compounds of African Origin. <i>Molecules</i> , 2021, 26, 406.	3.8	35
4	Pharmacoinformatics-based identification of potential bioactive compounds against Ebola virus protein VP24. <i>Computers in Biology and Medicine</i> , 2019, 113, 103414.	7.0	32
5	Molecular Docking and Dynamics Simulation Studies Predict Munc18b as a Target of Mycolactone: A Plausible Mechanism for Granule Exocytosis Impairment in Buruli Ulcer Pathogenesis. <i>Toxins</i> , 2019, 11, 181.	3.4	30
6	In Silico Screening of Isocitrate Lyase for Novel Anti-Buruli Ulcer Natural Products Originating from Africa. <i>Molecules</i> , 2018, 23, 1550.	3.8	29
7	Computational Identification of Potential Anti-Inflammatory Natural Compounds Targeting the p38 Mitogen-Activated Protein Kinase (MAPK): Implications for COVID-19-Induced Cytokine Storm. <i>Biomolecules</i> , 2021, 11, 653.	4.0	25
8	Chitosan Composites Synthesized Using Acetic Acid and Tetraethylorthosilicate Respond Differently to Methylene Blue Adsorption. <i>Polymers</i> , 2018, 10, 466.	4.5	24
9	Multitargeted Effects of Vitexin and Isovitexin on Diabetes Mellitus and Its Complications. <i>Scientific World Journal, The</i> , 2021, 2021, 1-20.	2.1	24
10	A Molecular Modeling Approach to Identify Potential Antileishmanial Compounds Against the Cell Division Cycle (cdc)-2-Related Kinase 12 (CRK12) Receptor of <i>Leishmania donovani</i> . <i>Biomolecules</i> , 2021, 11, 458.	4.0	19
11	Therapeutic Anti-Depressant Potential of Microbial GABA Produced by <i>Lactobacillus rhamnosus</i> Strains for GABAergic Signaling Restoration and Inhibition of Addiction-Induced HPA Axis Hyperactivity. <i>Current Issues in Molecular Biology</i> , 2022, 44, 1434-1451.	2.4	19
12	Ergonomics assessment of locally fabricated passenger seats in trotro vehicles in Accra, Ghana. <i>Journal of Transport and Health</i> , 2017, 6, 167-176.	2.2	17
13	Molecular Docking Simulation Studies Identifies Potential Natural Product Derived-Antiwoibachial Compounds as Filaricides against Onchocerciasis. <i>Biomedicines</i> , 2021, 9, 1682.	3.2	17
14	Leishmanicidal Potential of Hardwickiic Acid Isolated From <i>Croton sylvaticus</i> . <i>Frontiers in Pharmacology</i> , 2020, 11, 753.	3.5	16
15	Dragon exploratory system on Hepatitis C Virus (DESHCV). <i>Infection, Genetics and Evolution</i> , 2011, 11, 734-739.	2.3	13
16	The Search for Putative Hits in Combating Leishmaniasis: The Contributions of Natural Products Over the Last Decade. <i>Natural Products and Bioprospecting</i> , 2021, 11, 489-544.	4.3	13
17	Outwitting an Old Neglected Nemesis: A Review on Leveraging Integrated Data-Driven Approaches to Aid in Unraveling of Leishmanicides of Therapeutic Potential. <i>Current Topics in Medicinal Chemistry</i> , 2020, 20, 349-366.	2.1	13
18	Computational Study on Potential Novel Anti-Ebola Virus Protein VP35 Natural Compounds. <i>Biomedicines</i> , 2021, 9, 1796.	3.2	11

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19	Molecular Structure-Based Screening of the Constituents of <i>Calotropis procera</i> Identifies Potential Inhibitors of Diabetes Mellitus Target Alpha Glucosidase. <i>Current Issues in Molecular Biology</i> , 2022, 44, 963-987.	2.4	11
20	The emerging paradigm of calcium homeostasis as a new therapeutic target for protozoan parasites. <i>Medicinal Research Reviews</i> , 2022, 42, 56-82.	10.5	8
21	Homology Modeling, de Novo Design of Ligands, and Molecular Docking Identify Potential Inhibitors of <i>Leishmania donovani</i> 24-Sterol Methyltransferase. <i>Frontiers in Cellular and Infection Microbiology</i> , 2022, 12, .	3.9	8
22	Molecular Informatics Studies of the Iron-Dependent Regulator (ideR) Reveal Potential Novel Anti- <i>Mycobacterium ulcerans</i> Natural Product-Derived Compounds. <i>Molecules</i> , 2019, 24, 2299.	3.8	7
23	Antimalarials: Review of Plasmepsins as Drug Targets and HIV Protease Inhibitors Interactions. <i>Current Topics in Medicinal Chemistry</i> , 2019, 18, 2022-2028.	2.1	7
24	Review of Atypical Organometallic Compounds as Antimalarial Drugs. <i>Journal of Chemistry</i> , 2020, 2020, 1-9.	1.9	7
25	Unravelling the myth surrounding sterol biosynthesis as plausible target for drug design against leishmaniasis. <i>Journal of Parasitic Diseases</i> , 2021, 45, 1152-1171.	1.0	6
26	Molecular docking and dynamics simulations studies of OmpATb identifies four potential novel natural product-derived anti- <i>Mycobacterium tuberculosis</i> compounds. <i>Computers in Biology and Medicine</i> , 2020, 122, 103811.	7.0	6
27	Electrochemical Response of <i>Saccharomyces cerevisiae</i> Corresponds to Cell Viability upon Exposure to <i>Dioclea reflexa</i> Seed Extracts and Antifungal Drugs. <i>Biosensors</i> , 2019, 9, 45.	4.7	5
28	Pharmacophore-Guided Identification of Natural Products as Potential Inhibitors of <i>Mycobacterium ulcerans</i> Cystathionine β -Synthase MetB. <i>Computation</i> , 2021, 9, 32.	2.0	4
29	Molecular modelling and de novo fragment-based design of potential inhibitors of beta-tubulin gene of <i>Necator americanus</i> from natural products. <i>Informatics in Medicine Unlocked</i> , 2021, 26, 100734.	3.4	4
30	Antileishmanial effects of <i>Sargassum vulgare</i> products and prediction of trypanothione reductase inhibition by fucosterol. <i>Future Drug Discovery</i> , 2020, 2, .	2.1	3
31	Investigating the Conformation of S100 β Protein Under Physiological Parameters Using Computational Modeling: A Clue for Rational Drug Design. <i>Open Biomedical Engineering Journal</i> , 2018, 12, 36-50.	0.5	3
32	Development of a proteochemometric-based support vector machine model for predicting bioactive molecules of tubulin receptors. <i>Molecular Diversity</i> , 2022, 26, 2231-2242.	3.9	3
33	Density Functional Theory-Based Studies Predict Carbon Nanotubes as Effective Mycolactone Inhibitors. <i>Molecules</i> , 2022, 27, 4440.	3.8	3
34	A plasmonic photo-thermal probe for thermoablation of post-operative breast cancer cells. <i>Cogent Engineering</i> , 2017, 4, 1331966.	2.2	2
35	Identification of novel potential inhibitors of varicella-zoster virus thymidine kinase from ethnopharmacologic relevant plants through an in-silico approach. <i>Journal of Biomolecular Structure and Dynamics</i> , 2021, , 1-16.	3.5	2
36	Surveys, Serologies, and Sequences Reveal History of Iatrogenic Transmission of HIV-1. <i>Journal of Infectious Diseases</i> , 2016, 214, 341-343.	4.0	1

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37	Therapeutic potential of HIV-1 entry inhibitor peptidomimetics. <i>Experimental Biology and Medicine</i> , 2021, 246, 1060-1068.	2.4	1
38	Prediction of antischistosomal small molecules using machine learning in the era of big data. <i>Molecular Diversity</i> , 2021, , 1.	3.9	1
39	Dual Inhibition of Human N-Myristoyltransferase Subtypes Halts Common Cold Pathogenesis: Atomistic Perspectives from the Case of IMP1088**. <i>Chemistry and Biodiversity</i> , 2021, , .	2.1	1
40	Computer-aided identification of potential inhibitors against <i>Necator americanus</i> glutathione S-transferase 3. <i>Informatics in Medicine Unlocked</i> , 2022, 30, 100957.	3.4	1
41	Electrochemical Response of Cells Using Bioactive Plant Isolates. , 0, , .		0
42	Investigating the Conformation of S100Î² Protein Under Physiological Parameters Using Computational Modeling: A Clue for Rational Drug Design. <i>Open Biomedical Engineering Journal</i> , 2018, 12, 73-73.	0.5	0
43	Deep Neural Networks predict inhibitors of <i>Schistosoma mansoni</i> thioredoxin glutathione reductase (SmTGR). <i>Journal of Computational Biophysics and Chemistry</i> , 0, , .	1.7	0
44	Molecular Informatics of Trypanothione Reductase of <i>Leishmania major</i> Reveals Novel Chromen-2-One Analogues as Potential Leishmanicides. , 0, , .		0