ViÅ;nja Stepanić

List of Publications by Year in descending order

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#	Article	IF	CITATIONS
1	Identification of 37 Heterogeneous Drug Candidates for Treatment of COVID-19 via a Rational Transcriptomics-Based Drug Repurposing Approach. Pharmaceuticals, 2021, 14, 87.	1.7	5
2	Involvement of NRF2 in Breast Cancer and Possible Therapeutical Role of Polyphenols and Melatonin. Molecules, 2021, 26, 1853.	1.7	31
3	Crowdsourced mapping of unexplored target space of kinase inhibitors. Nature Communications, 2021, 12, 3307.	5.8	41
4	Comprehensive machine learning based study of the chemical space of herbicides. Scientific Reports, 2021, 11, 11479.	1.6	18
5	Green One-Pot Synthesis of Coumarin-Hydroxybenzohydrazide Hybrids and Their Antioxidant Potency. Antioxidants, 2021, 10, 1106.	2.2	31
6	Antioxidative potential of ferulic acid phenoxyl radical. Phytochemistry, 2020, 170, 112218.	1.4	40
7	Modulators of Oxidative Stress: Chemical and Pharmacological Aspects. Antioxidants, 2020, 9, 657.	2.2	8
8	The Antioxidant and Antiproliferative Activities of 1,2,3-Triazolyl-L-Ascorbic Acid Derivatives. International Journal of Molecular Sciences, 2019, 20, 4735.	1.8	15
9	Antitumor and antiviral activities of 4-substituted 1,2,3-triazolyl-2,3-dibenzyl-L-ascorbic acid derivatives. European Journal of Medicinal Chemistry, 2019, 184, 111739.	2.6	25
10	Curcumin and its Potential for Systemic Targeting of Inflamm-Aging and Metabolic Reprogramming in Cancer. International Journal of Molecular Sciences, 2019, 20, 1180.	1.8	19
11	Antioxidant Activities of Alkyl Substituted Pyrazine Derivatives of Chalcones—In Vitro and In Silico Study. Antioxidants, 2019, 8, 90.	2.2	31
12	Effects of conjugation metabolism on radical scavenging and transport properties of quercetin – In silico study. Journal of Molecular Graphics and Modelling, 2019, 86, 278-285.	1.3	5
13	The Influence of InÂVivo Metabolic Modifications on ADMET Properties of Green Tea Catechins–In Silico Analysis. Journal of Pharmaceutical Sciences, 2018, 107, 2957-2964.	1.6	3
14	European contribution to the study of ROS: A summary of the findings and prospects for the future from the COST action BM1203 (EU-ROS). Redox Biology, 2017, 13, 94-162.	3.9	242
15	Free radical scavenging potency of quercetin catecholic colonic metabolites: Thermodynamics of 2H+/2eâ^' processes. Food Chemistry, 2017, 218, 144-151.	4.2	83
16	Synthesis and SAR Study of Novel Amidino 2-substituted Benzimidazoles as Potential Antibacterial Agents. Croatica Chemica Acta, 2017, 90, .	0.1	7
17	Synthesis, inÂvitro anticancer and antibacterial activities and in silico studies of new 4-substituted 1,2,3-triazole–coumarin hybrids. European Journal of Medicinal Chemistry, 2016, 124, 794-808. 	2.6	110
18	Synthesis and evaluation of antibacterial and antioxidant activity of novel 2-phenyl-quinoline analogs derivatized at position 4 with aromatically substituted 4H-1,2,4-triazoles. Journal of Enzyme Inhibition and Medicinal Chemistry, 2016, 31, 104-110.	2.5	19

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19	The 2H+/2eâ^' free radical scavenging mechanisms of uric acid: thermodynamics of NH bond cleavage. Computational and Theoretical Chemistry, 2016, 1077, 2-10.	1.1	22
20	Synthesis and structure–activity relationship of amidine derivatives of 3,4-ethylenedioxythiophene as novel antibacterial agents. European Journal of Medicinal Chemistry, 2015, 90, 68-81.	2.6	25
21	Selected Attributes of Polyphenols in Targeting Oxidative Stress in Cancer. Current Topics in Medicinal Chemistry, 2015, 15, 496-509.	1.0	56
22	Towards an improved prediction of the free radical scavenging potency of flavonoids: The significance of double PCET mechanisms. Food Chemistry, 2014, 152, 578-585.	4.2	54
23	Correlation between 13C NMR chemical shifts and antiradical activity of flavonoids. Monatshefte Für Chemie, 2014, 145, 457-463.	0.9	6
24	Synthesis and in vitro antiproliferative evaluation of novel N-alkylated 6-isobutyl- and propyl pyrimidine derivatives. Bioorganic and Medicinal Chemistry Letters, 2014, 24, 2913-2917.	1.0	11
25	Influence of different free radicals on scavenging potency of gallic acid. Journal of Molecular Modeling, 2014, 20, 2345.	0.8	38
26	PM6 study of free radical scavenging mechanisms of flavonoids: why does O–H bond dissociation enthalpy effectively represent free radical scavenging activity?. Journal of Molecular Modeling, 2013, 19, 2593-2603.	0.8	75
27	Study of lipophilicity and membrane partition of 4-hydroxycoumarins by HPLC and PCA. Journal of Pharmaceutical and Biomedical Analysis, 2013, 76, 104-111.	1.4	4
28	Structure–property relationship for cellular accumulation of macrolones in human polymorphonuclear leukocytes (PMNs). European Journal of Pharmaceutical Sciences, 2013, 49, 206-219.	1.9	7
29	Bond dissociation free energy as a general parameter for flavonoid radical scavenging activity. Food Chemistry, 2013, 141, 1562-1570.	4.2	78
30	Antitumor Mechanisms of Amino Acid Hydroxyurea Derivatives in the Metastatic Colon Cancer Model. International Journal of Molecular Sciences, 2013, 14, 23654-23671.	1.8	1
31	Free radical scavenging activity of morin 2′-Oâ^' phenoxide anion. Food Chemistry, 2012, 135, 2070-2077.	4.2	45
32	Synthesis and antiproliferative evaluation of some new amidino-substituted bis-benzothiazolyl-pyridines and pyrazine. European Journal of Medicinal Chemistry, 2012, 55, 108-116.	2.6	23
33	Tebrophen — An Old Polyphenol Drug with Anticancer Potential â€. Molecules, 2012, 17, 7864-7886.	1.7	2
34	PM6 and DFT study of free radical scavenging activity of morin. Food Chemistry, 2012, 134, 1754-1760.	4.2	97
35	Physicochemical profile of macrolides and their comparison with small molecules. European Journal of Medicinal Chemistry, 2012, 47, 462-472.	2.6	33
36	An efficient and convenient microwave-assisted chemical synthesis of (thio)xanthones with additional in vitro and in silico characterization. Bioorganic and Medicinal Chemistry, 2012, 20, 3180-3185.	1.4	19

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37	Modeling Cellular Pharmacokinetics of 14- and 15-Membered Macrolides with Physicochemical Properties. Journal of Medicinal Chemistry, 2011, 54, 719-733.	2.9	40
38	6-Imino-2-thioxo-pyrimidinones as a new class of dipeptidyl peptidase IV inhibitors. Medicinal Chemistry Research, 2011, 20, 339-345.	1.1	2
39	Structure and vibrational spectra of conjugated acids of trans- and cis-azobenzene. Journal of Molecular Structure, 2001, 569, 89-109.	1.8	17
40	Ground and excited states of isodiazene – an ab initio study. Chemical Physics, 2000, 254, 151-168.	0.9	12