

Damir Vukicevic

List of Publications by Year in descending order

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81
papers

1,848
citations

361296
20
h-index

289141
40
g-index

84
all docs

84
docs citations

84
times ranked

1146
citing authors

#	ARTICLE	IF	CITATIONS
1	Evaluating topological ordering in directed acyclic graphs. <i>Electronic Journal of Graph Theory and Applications</i> , 2021, 9, 567.	0.2	1
2	A few examples and counterexamples in spectral graph theory. <i>Discussiones Mathematicae - Graph Theory</i> , 2020, 40, 637.	0.2	2
3	Optimizing the diagnostic capacity for COVID-19 PCR testing for low resource and high demand settings: The development of information-dependent pooling protocol. <i>Journal of Global Health</i> , 2020, 10, 020515.	1.2	7
4	Selection and redesign for high selectivity of membrane-active antimicrobial peptides from a dedicated sequence/function database. <i>Biochimica Et Biophysica Acta - Biomembranes</i> , 2019, 1861, 827-834.	1.4	22
5	One-alpha weighted network descriptors. <i>Rad Hrvatske Akademije Znanosti I Umjetnosti MatematiĀke Znanosti</i> , 2019, Knj. 538, 58, 31-49.	0.6	0
6	Exponential generalised network descriptors. <i>Advances in Mathematics of Communications</i> , 2019, 13, 405-420.	0.4	0
7	Global forcing number for maximal matchings. <i>Discrete Mathematics</i> , 2018, 341, 801-809.	0.4	4
8	Antibacterial Activity Affected by the Conformational Flexibility in Glycineâ€“Lysine Based Î±-Helical Antimicrobial Peptides. <i>Journal of Medicinal Chemistry</i> , 2018, 61, 2924-2936.	2.9	48
9	On indices of Wiener and anti-Wiener type. <i>Discrete Applied Mathematics</i> , 2018, 251, 290-298.	0.5	1
10	Multicoloring of graphs to secure a secret. <i>Rad Hrvatske Akademije Znanosti I Umjetnosti, Matematicke Znanosti</i> , 2018, 534, 1-22.	0.4	1
11	Tools for Designing Amphipathic Helical Antimicrobial Peptides. <i>Methods in Molecular Biology</i> , 2017, 1548, 23-34.	0.4	10
12	Relative edge betweenness centrality. <i>Ars Mathematica Contemporanea</i> , 2017, 12, 261-270.	0.3	5
13	A measure for a balanced workload and its extremal values. <i>Discrete Applied Mathematics</i> , 2016, 200, 59-66.	0.5	2
14	Predicting the Minimal Inhibitory Concentration for Antimicrobial Peptides with Rana-Box Domain. <i>Journal of Chemical Information and Modeling</i> , 2015, 55, 2275-2287.	2.5	17
15	Community structure in networks: Girvan-Newman algorithm improvement. , 2014, , .		28
16	Compression ratio of Wiener index in 2-d rectangular and polygonal lattices. <i>Ars Mathematica Contemporanea</i> , 2014, 7, 1-12.	0.3	8
17	Network descriptors based on betweenness centrality and transmission and their extremal values. <i>Discrete Applied Mathematics</i> , 2013, 161, 2678-2686.	0.5	11
18	Generalised network descriptors. <i>Glasnik Matematicki</i> , 2013, 48, 211-230.	0.1	0

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19	Limitations of Pauling Bond Order Concept. <i>Polycyclic Aromatic Compounds</i> , 2012, 32, 36-47.	1.4	3
20	Improving the Selectivity of Antimicrobial Peptides from Anuran Skin. <i>Journal of Chemical Information and Modeling</i> , 2012, 52, 3341-3351.	2.5	30
21	Computer search for trees with minimal ABC index. <i>Applied Mathematics and Computation</i> , 2012, 219, 767-772.	1.4	25
22	Electron currents in polycyclic conjugated hydrocarbons: Coronene and its isomers having five and seven member rings. <i>International Journal of Quantum Chemistry</i> , 2012, 112, 972-985.	1.0	14
23	Electron currents in larger fully aromatic benzenoids. <i>International Journal of Quantum Chemistry</i> , 2012, 112, 2456-2462.	1.0	3
24	Conjugated circuits currents in hexabenzocoronene and its derivatives formed by joining proximal carbons. <i>Journal of Computational Chemistry</i> , 2012, 33, 1111-1122.	1.5	7
25	On the vertex degree indices of connected graphs. <i>Chemical Physics Letters</i> , 2011, 512, 283-286.	1.2	24
26	Bond additive modeling 10. Upper and lower bounds of bond incident degree indices of catacondensed fluoranthenes. <i>Chemical Physics Letters</i> , 2011, 515, 186-189.	1.2	13
27	Using size for bounding expressions of graph invariants. <i>Annals of Operations Research</i> , 2011, 188, 415-427.	2.6	0
28	Knowledge-based computational methods for identifying or designing novel, non-homologous antimicrobial peptides. <i>European Biophysics Journal</i> , 2011, 40, 371-385.	1.2	50
29	Topological efficiency of C66 fullerene. <i>Chemical Physics Letters</i> , 2011, 501, 442-445.	1.2	44
30	Applications of Perfect Matchings in Chemistry. , 2011, , 463-482.		1
31	Topological Determination of ¹³ C-NMR Spectra of C66 Fullerenes. <i>Carbon Materials</i> , 2011, , 205-216.	0.2	3
32	Which generalized Randić indices are suitable measures of molecular branching?. <i>Discrete Applied Mathematics</i> , 2010, 158, 2056-2065.	0.5	4
33	A note on the Estrada-Hatano communicability algorithm for detecting community structure in complex networks. <i>Applied Mathematics and Computation</i> , 2010, 217, 3516-3521.	1.4	4
34	Augmented Zagreb index. <i>Journal of Mathematical Chemistry</i> , 2010, 48, 370-380.	0.7	299
35	One-two descriptor. <i>Journal of Mathematical Chemistry</i> , 2010, 48, 395-400.	0.7	1
36	On the number of Kekulé structures of fluoranthene congeners. <i>Journal of the Serbian Chemical Society</i> , 2010, 75, 1093-1098.	0.4	5

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37	On the extremal values of ratios of number of paths. <i>Ars Mathematica Contemporanea</i> , 2010, 3, 215-235.	0.3	4
38	Master Connectivity Index and Master Connectivity Polynomial. <i>Current Computer-Aided Drug Design</i> , 2010, 6, 235-239.	0.8	1
39	Note on the comparison of the first and second normalized zagreb eccentricity indices. <i>Acta Chimica Slovenica</i> , 2010, 57, 524-8.	0.2	39
40	Design of selective peptide antibiotics by using the sequence moment concept. <i>Nature Precedings</i> , 2009, , ,	0.1	0
41	Design of selective peptide antibiotics by using the sequence moment concept. <i>Nature Precedings</i> , 2009, , ,	0.1	0
42	On Wiener-type polynomials of thorn graphs. <i>Journal of Chemometrics</i> , 2009, 23, 600-604.	0.7	7
43	Variable neighborhood search for extremal graphs. 23. On the Randić index and the chromatic number. <i>Discrete Mathematics</i> , 2009, 309, 4228-4234.	0.4	31
44	On Kekulé structures count. <i>Journal of Mathematical Chemistry</i> , 2009, 45, 279-286.	0.7	1
45	On decompositions of leapfrog fullerenes. <i>Journal of Mathematical Chemistry</i> , 2009, 45, 406-416.	0.7	2
46	On the anti-Kekulé number of leapfrog fullerenes. <i>Journal of Mathematical Chemistry</i> , 2009, 45, 431-441.	0.7	4
47	Tubercular fulleroids. <i>Journal of Mathematical Chemistry</i> , 2009, 45, 513-524.	0.7	1
48	On the path-Zagreb matrix. <i>Journal of Mathematical Chemistry</i> , 2009, 45, 538-543.	0.7	7
49	On functionalized fullerenes C ₆₀ X _n . <i>Journal of Mathematical Chemistry</i> , 2009, 45, 557-562.	0.7	1
50	Topological index based on the ratios of geometrical and arithmetical means of end-vertex degrees of edges. <i>Journal of Mathematical Chemistry</i> , 2009, 46, 1369-1376.	0.7	478
51	Graphical representation of proteins as four-color maps and their numerical characterization. <i>Journal of Molecular Graphics and Modelling</i> , 2009, 27, 637-641.	1.3	45
52	Atom-bond connectivity index of trees. <i>Discrete Applied Mathematics</i> , 2009, 157, 2828-2835.	0.5	106
53	Computational Design of Highly Selective Antimicrobial Peptides. <i>Journal of Chemical Information and Modeling</i> , 2009, 49, 2873-2882.	2.5	79
54	On the edge degrees of trees. <i>Glasnik Matematički</i> , 2009, 44, 259-266.	0.1	7

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55	On the anti-Kekulé number and anti-forcing number of cata-condensed benzenoids. <i>Journal of Mathematical Chemistry</i> , 2008, 43, 719-726.	0.7	26
56	Partition of π -electrons between faces of polyhedral carbon aggregates. <i>Journal of Mathematical Chemistry</i> , 2008, 43, 773-779.	0.7	8
57	Note on ordering and complexity of Platonic and Archimedean polyhedra based on solid angles. <i>Journal of Mathematical Chemistry</i> , 2008, 44, 725-730.	0.7	4
58	Binary coding of algebraic Kekulé structures of catacondensed benzenoid graphs. <i>Applied Mathematics Letters</i> , 2008, 21, 712-716.	1.5	2
59	Statistical investigation of new topological indices based on the molecular path code. <i>Chemical Physics Letters</i> , 2008, 464, 155-159.	1.2	4
60	The Anti-Kekule number of the infinite triangular, rectangular and hexagonal grids. <i>Glasnik Matematički</i> , 2008, 43, 243-252.	0.1	7
61	Numerical Kekulé Structures of Fullerenes and Partitioning of π -Electrons to Pentagonal and Hexagonal Rings. <i>Journal of Chemical Information and Modeling</i> , 2007, 47, 897-904.	2.5	22
62	Linear regression model of DNA sequences and its application. <i>Journal of Computational Chemistry</i> , 2007, 28, 1434-1445.	1.5	9
63	Computing the bipartite edge frustration of fullerene graphs. <i>Discrete Applied Mathematics</i> , 2007, 155, 1294-1301.	0.5	35
64	Mathematical studies of Kekulé structures. <i>Structural Chemistry</i> , 2007, 18, 807-812.	1.0	4
65	Kekulé Structure Count in Corazulenic Fullerenes. <i>Journal of Nanoscience and Nanotechnology</i> , 2007, 7, 1321-1328.	0.9	3
66	Vindicating the Pauling-bond-order concept. <i>Chemical Physics Letters</i> , 2006, 427, 418-420.	1.2	16
67	On the complexity of Archimedean solids. <i>Journal of Mathematical Chemistry</i> , 2006, 39, 119-132.	0.7	9
68	On Acyclic Molecular Graphs with Prescribed Numbers of Edges that Connect Vertices with given Degrees. <i>Journal of Mathematical Chemistry</i> , 2006, 40, 155-178.	0.7	5
69	On Kekulé structures of buckminsterfullerene. <i>Chemical Physics Letters</i> , 2005, 401, 446-450.	1.2	31
70	An efficient method to enumerate topologically distinct clusters of hydrogen-bonding in water molecules. <i>Chemical Physics Letters</i> , 2005, 416, 212-214.	1.2	13
71	Characterization of distribution of π -electrons amongst benzenoid rings for Randić's algebraic Kekulé structures. <i>Journal of Mathematical Chemistry</i> , 2005, 37, 163-170.	0.7	4
72	Canonical Labeling of Proteome Maps. <i>Journal of Proteome Research</i> , 2005, 4, 1347-1352.	1.8	18

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73	Paths and walks in acyclic structures: Kenograms versus plerograms. <i>Arkivoc</i> , 2005, 2005, 33-44.	0.3	4
74	Partitioning of π -electrons in Rings of Polycyclic Conjugated Hydrocarbons. Part 4. Benzenoids with More Than One Geometric Kekulé Structure Corresponding to the Same Algebraic Kekulé Structure. <i>Journal of Mathematical Chemistry</i> , 2004, 36, 271-279.	0.7	30
75	Algebraic Kekule Structures of Benzenoid Hydrocarbons.. <i>ChemInform</i> , 2004, 35, no.	0.1	0
76	Algebraic Kekulé Structures of Benzenoid Hydrocarbons. <i>Journal of Chemical Information and Computer Sciences</i> , 2004, 44, 296-299.	2.8	34
77	Almost All Trees and Chemical Trees Have Equiseparable Mates. <i>Journal of Computer Chemistry Japan</i> , 2004, 3, 109-112.	0.0	2
78	Decomposition of complete graphs into factors of diameter two and three. <i>Discussiones Mathematicae - Graph Theory</i> , 2003, 23, 37.	0.2	0
79	Mix-decomposition of the complete graph into directed factors of diameter 2 and undirected factors of diameter 3. <i>Glasnik Matematički</i> , 2003, 38, 211-232.	0.1	0
80	Bond Additive Modeling 4. QSPR and QSAR Studies of the Variable Adriatic Indices. <i>Croatica Chemica Acta</i> , 0, , 87-91.	0.1	30
81	Bond Additive Modeling 5. Mathematical Properties of the Variable Sum Exdeg Index. <i>Croatica Chemica Acta</i> , 0, , 93-101.	0.1	17