

Ross King

List of Publications by Year in descending order

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Version: 2024-02-01

32
papers

2,810
citations

331670

21
h-index

414414

32
g-index

32
all docs

32
docs citations

32
times ranked

2830
citing authors

#	ARTICLE	IF	CITATIONS
1	Testing the reproducibility and robustness of the cancer biology literature by robot. <i>Journal of the Royal Society Interface</i> , 2022, 19, 20210821.	3.4	2
2	A simple spatial extension to the extended connectivity interaction features for binding affinity prediction. <i>Royal Society Open Science</i> , 2022, 9, 211745.	2.4	2
3	Transformational machine learning: Learning how to learn from many related scientific problems. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2021, 118, .	7.1	11
4	Predicting rice phenotypes with meta and multi-target learning. <i>Machine Learning</i> , 2020, 109, 2195-2212.	5.4	3
5	Closed-loop cycles of experiment design, execution, and learning accelerate systems biology model development in yeast. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2019, 116, 18142-18147.	7.1	19
6	Plasmodium dihydrofolate reductase is a second enzyme target for the antimalarial action of triclosan. <i>Scientific Reports</i> , 2018, 8, 1038.	3.3	31
7	Computing exponentially faster: implementing a non-deterministic universal Turing machine using DNA. <i>Journal of the Royal Society Interface</i> , 2017, 14, 20160990.	3.4	30
8	Cheaper faster drug development validated by the repositioning of drugs against neglected tropical diseases. <i>Journal of the Royal Society Interface</i> , 2015, 12, 20141289.	3.4	97
9	The Use of Weighted Graphs for Large-Scale Genome Analysis. <i>PLoS ONE</i> , 2014, 9, e89618.	2.5	4
10	Merits of random forests emerge in evaluation of chemometric classifiers by external validation. <i>Analytica Chimica Acta</i> , 2013, 801, 22-33.	5.4	34
11	Yeast-based automated high-throughput screens to identify anti-parasitic lead compounds. <i>Open Biology</i> , 2013, 3, 120158.	3.6	32
12	On the formalization and reuse of scientific research. <i>Journal of the Royal Society Interface</i> , 2011, 8, 1440-1448.	3.4	15
13	Functional Expression of Parasite Drug Targets and Their Human Orthologs in Yeast. <i>PLoS Neglected Tropical Diseases</i> , 2011, 5, e1320.	3.0	29
14	Towards Robot Scientists for autonomous scientific discovery. <i>Automated Experimentation</i> , 2010, 2, 1.	2.0	101
15	Enhancement of Plant Metabolite Fingerprinting by Machine Learning \hat{A} . <i>Plant Physiology</i> , 2010, 153, 1506-1520.	4.8	24
16	The Automation of Science. <i>Science</i> , 2009, 324, 85-89.	12.6	458
17	Finding Motifs in Protein Secondary Structure for Use in Function Prediction. <i>Journal of Computational Biology</i> , 2006, 13, 719-731.	1.6	14
18	An ontology of scientific experiments. <i>Journal of the Royal Society Interface</i> , 2006, 3, 795-803.	3.4	147

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19	Quantitative pharmacophore models with inductive logic programming. <i>Machine Learning</i> , 2006, 64, 65-90.	5.4	12
20	Are the current ontologies in biology good ontologies?. <i>Nature Biotechnology</i> , 2005, 23, 1095-1098.	17.5	66
21	Hierarchical metabolomics demonstrates substantial compositional similarity between genetically modified and conventional potato crops. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2005, 102, 14458-14462.	7.1	367
22	Functional genomic hypothesis generation and experimentation by a robot scientist. <i>Nature</i> , 2004, 427, 247-252.	27.8	481
23	Intelligent software for laboratory automation. <i>Trends in Biotechnology</i> , 2004, 22, 440-445.	9.3	15
24	New Approach to Pharmacophore Mapping and QSAR Analysis Using Inductive Logic Programming. Application to Thermolysin Inhibitors and Glycogen Phosphorylase b Inhibitors. <i>Journal of Medicinal Chemistry</i> , 2002, 45, 399-409.	6.4	57
25	Modeling quantitative structure-property relationships in calculated reaction pathways using a new 3D quantum topological representation. <i>Analytica Chimica Acta</i> , 2001, 446, 3-13.	5.4	15
26	On the optimization of classes for the assignment of unidentified reading frames in functional genomics programmes: the need for machine learning. <i>Trends in Biotechnology</i> , 2000, 18, 93-98.	9.3	78
27	Structure-activity relationships derived by machine learning: the use of atoms and their bond connectivities to predict mutagenicity by inductive logic programming.. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 1996, 93, 438-442.	7.1	211
28	STATLOG: COMPARISON OF CLASSIFICATION ALGORITHMS ON LARGE REAL-WORLD PROBLEMS. <i>Applied Artificial Intelligence</i> , 1995, 9, 289-333.	3.2	234
29	Relating chemical activity to structure: An examination of ILP successes. <i>New Generation Computing</i> , 1995, 13, 411-433.	3.3	65
30	COMPARISON OF ARTIFICIAL INTELLIGENCE METHODS FOR MODELING PHARMACEUTICAL QSARS. <i>Applied Artificial Intelligence</i> , 1995, 9, 213-233.	3.2	30
31	New approaches to QSAR: Neural networks and machine learning. <i>Journal of Computer - Aided Molecular Design</i> , 1993, 1, 279-290.	1.0	26
32	Machine learning approach for the prediction of protein secondary structure. <i>Journal of Molecular Biology</i> , 1990, 216, 441-457.	4.2	100