Samik Ghosh

List of Publications by Year in descending order

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361413 254184 3,312 54 20 43 citations h-index g-index papers 57 57 57 5254 docs citations times ranked citing authors all docs

| # | Article | IF | CITATIONS |
|----|--|-----|-----------|
| 1 | Al-driven laboratory workflows enable operation in the age of social distancing. SLAS Technology, 2022, 27, 195-203. | 1.9 | 8 |
| 2 | Toxicity Analysis of Pentachlorophenol Data with a Bioinformatics Tool Set. Methods in Molecular Biology, 2022, 2486, 105-125. | 0.9 | 1 |
| 3 | Artificial intelligence-based computational framework for drug-target prioritization and inference of novel repositionable drugs for Alzheimer's disease. Alzheimer's Research and Therapy, 2021, 13, 92. | 6.2 | 29 |
| 4 | A Geometric Clustering Tool (AGCT) to robustly unravel the inner cluster structures of time-series gene expressions. PLoS ONE, 2020, 15, e0233755. | 2.5 | 3 |
| 5 | Computational Systems Biology Applications. , 2019, , 66-73. | | 0 |
| 6 | Computational Systems Biology. , 2019, , 789-795. | | 4 |
| 7 | Lenvatinib plus anti-PD-1 antibody combination treatment activates CD8+ T cells through reduction of tumor-associated macrophage and activation of the interferon pathway. PLoS ONE, 2019, 14, e0212513. | 2.5 | 294 |
| 8 | Multi-dimensional computational pipeline for large-scale deep screening of compound effect assessment: an in silico case study on ageing-related compounds. Npj Systems Biology and Applications, 2019, 5, 42. | 3.0 | 3 |
| 9 | Sequence homology in eukaryotes (SHOE): interactive visual tool for promoter analysis. BMC Genomics, 2018, 19, 715. | 2.8 | 7 |
| 10 | Precise Temporal Profiling of Signaling Complexes in Primary Cells Using SWATH Mass Spectrometry. Cell Reports, 2017, 18, 3219-3226. | 6.4 | 28 |
| 11 | Supporting evidence-based analysis for modified risk tobacco products through a toxicology data-sharing infrastructure. F1000Research, 2017, 6, 12. | 1.6 | 7 |
| 12 | Supporting evidence-based analysis for modified risk tobacco products through a toxicology data-sharing infrastructure. F1000Research, 2017, 6, 12. | 1.6 | 10 |
| 13 | A machine learning approach for the identification of key markers involved in brain development from single-cell transcriptomic data. BMC Genomics, 2016, 17, 1025. | 2.8 | 35 |
| 14 | Network analyses based on comprehensive molecular interaction maps reveal robust control structures in yeast stress response pathways. Npj Systems Biology and Applications, 2016, 2, 15018. | 3.0 | 27 |
| 15 | Digital health revolution: perfect storm or perfect opportunity for pharmaceutical R&D?. Drug Discovery Today, 2016, 21, 900-911. | 6.4 | 42 |
| 16 | Elucidation of the molecular mechanisms underlying adverse reactions associated with a kinase inhibitor using systems toxicology. Npj Systems Biology and Applications, 2015, 1, 15005. | 3.0 | 16 |
| 17 | Application of machine leaning approaches in drug target identification and network pharmacology. , 2015, , . | | 0 |
| 18 | Oscillation of cAMP and Ca2+ in cardiac myocytes: a systems biology approach. Journal of Physiological Sciences, 2015, 65, 195-200. | 2.1 | 15 |

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|----|--|------|-----------|
| 19 | Weaving Knowledge into Biological Pathways in a Collaborative Manner. Methods in Pharmacology and Toxicology, 2015, , 181-208. | 0.2 | 1 |
| 20 | Software Platform for Systems Biology. Drug Delivery System, 2014, 29, 386-396. | 0.0 | 0 |
| 21 | A Versatile Platform for Multilevel Modeling of Physiological Systems: SBML-PHML Hybrid Modeling and Simulation. Advanced Biomedical Engineering, 2014, 3, 50-58. | 0.6 | 15 |
| 22 | Identification of drug-target modules in the human protein–protein interaction network. Artificial Life and Robotics, 2014, 19, 406-413. | 1.2 | 4 |
| 23 | Integrating Pathways of Parkinson's Disease in a Molecular Interaction Map. Molecular Neurobiology, 2014, 49, 88-102. | 4.0 | 231 |
| 24 | Modeling and Simulation Using CellDesigner. Methods in Molecular Biology, 2014, 1164, 121-145. | 0.9 | 60 |
| 25 | A comprehensive map of the influenza A virus replication cycle. BMC Systems Biology, 2013, 7, 97. | 3.0 | 97 |
| 26 | Toward an integrated software platform for systems pharmacology. Biopharmaceutics and Drug Disposition, 2013, 34, 508-526. | 1.9 | 18 |
| 27 | Software Platform for Metabolic Network Reconstruction of Mycobacterium tuberculosis., 2013,, 21-35. | | 1 |
| 28 | Harnessing Diversity towards the Reconstructing of Large Scale Gene Regulatory Networks. PLoS Computational Biology, 2013, 9, e1003361. | 3.2 | 32 |
| 29 | A versatile platform for multilevel modeling of physiological systems: Template/instance framework for large-scale modeling and simulation., 2013, 2013, 5529-32. | | 7 |
| 30 | Combining Machine Learning Systems and Multiple Docking Simulation Packages to Improve Docking Prediction Reliability for Network Pharmacology. PLoS ONE, 2013, 8, e83922. | 2.5 | 268 |
| 31 | CTen: a web-based platform for identifying enriched cell types from heterogeneous microarray data. BMC Genomics, 2012, 13, 460. | 2.8 | 113 |
| 32 | AlzPathway: a comprehensive map of signaling pathways of Alzheimer's disease. BMC Systems Biology, 2012, 6, 52. | 3.0 | 114 |
| 33 | Software for systems biology: from tools to integrated platforms. Nature Reviews Genetics, 2011, 12, 821-832. | 16.3 | 228 |
| 34 | Social engineering for virtual 'big science' in systems biology. Nature Chemical Biology, 2011, 7, 323-326. | 8.0 | 35 |
| 35 | Discrete diffusion models to study the effects of Mg2+ concentration on the PhoPQ signal transduction system. BMC Genomics, 2010, 11, S3. | 2.8 | 3 |
| 36 | Connecting the dots: role of standardization and technology sharing in biological simulation. Drug Discovery Today, 2010, 15, 1024-1031. | 6.4 | 6 |

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|----|---|------|-----------|
| 37 | Payao: a community platform for SBML pathway model curation. Bioinformatics, 2010, 26, 1381-1383. | 4.1 | 50 |
| 38 | A comprehensive molecular interaction map of the budding yeast cell cycle. Molecular Systems Biology, 2010, 6, 415. | 7.2 | 62 |
| 39 | A comprehensive map of the mTOR signaling network. Molecular Systems Biology, 2010, 6, 453. | 7.2 | 201 |
| 40 | Consistent design schematics for biological systems: standardization of representation in biological engineering. Journal of the Royal Society Interface, 2009, 6, S393-404. | 3.4 | 15 |
| 41 | The Systems Biology Graphical Notation. Nature Biotechnology, 2009, 27, 735-741. | 17.5 | 828 |
| 42 | Parametric modeling of protein–DNA binding kinetics: A discrete event based simulation approach. Discrete Applied Mathematics, 2009, 157, 2395-2415. | 0.9 | 2 |
| 43 | Construction and analysis of a modular model of caspase activation in apoptosis. Theoretical Biology and Medical Modelling, 2008, 5, 26. | 2.1 | 89 |
| 44 | Holding time estimation for reactions in stochastic event-based simulation of complex biological systems. Simulation Modelling Practice and Theory, 2008, 16, 1615-1639. | 3.8 | 4 |
| 45 | Modeling protein-DNA binding time in Stochastic Discrete Event Simulation of Biological Processes. , 2007, , . | | 5 |
| 46 | Modeling the Stochastic Dynamics of Gene Expression in Single Cells: A Birth and Death Markov Chain Analysis. , 2007, , . | | 5 |
| 47 | Channel Assignment Strategies for Multiradio Wireless Mesh Networks: Issues and Solutions. , 2007, 45, 86-95. | | 250 |
| 48 | Towards Optimal Virtual Patients: An Online Adaptive Control Approach. Annual International Conference of the IEEE Engineering in Medicine and Biology Society, 2007, 2007, 3292-5. | 0.5 | 3 |
| 49 | Enabling real-time fleet route planning and execution in a pervasive transportation environment. , 2007, , . | | 3 |
| 50 | "MeshUp― Self-organizing mesh-based topologies for next generation radio access networks. Ad Hoc Networks, 2007, 5, 652-679. | 5.5 | 6 |
| 51 | Revisiting the Optimal Partitioning of Zones in Next Generation Cellular Networks: A Network Capacity Impact Perspective. , 2007, , 1011-1023. | | 0 |
| 52 | A Computationally Fast and Parametric Model to Estimate Protein-Ligand Docking Time for Stochastic Event Based Simulation. , 2007, , 14-41. | | 3 |
| 53 | A markov model based analysis of stochastic biochemical systems. Computational Systems Bioinformatics / Life Sciences Society Computational Systems Bioinformatics Conference, 2007, 6, 121-32. | 0.4 | 3 |
| 54 | A Stochastic model to estimate the time taken for Protein-Ligand Docking. , 2006, , . | | 8 |