

Susan M Mniszewski

List of Publications by Year in descending order

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Version: 2024-02-01

38
papers

812
citations

759233

12
h-index

610901

24
g-index

40
all docs

40
docs citations

40
times ranked

851
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 1 | Sampling electronic structure quadratic unconstrained binary optimization problems (QUBOs) with Ocean and Mukai solvers. PLoS ONE, 2022, 17, e0263849. | 2.5 | 5 |
| 2 | Cabana: A Performance Portable Library for Particle-Based Simulations. Journal of Open Source Software, 2022, 7, 4115. | 4.6 | 7 |
| 3 | Controlled precision QUBO-based algorithm to compute eigenvectors of symmetric matrices. PLoS ONE, 2022, 17, e0267954. | 2.5 | 3 |
| 4 | Toward a QUBO-Based Density Matrix Electronic Structure Method. Journal of Chemical Theory and Computation, 2022, 18, 4177-4185. | 5.3 | 2 |
| 5 | Quantum Perturbation Theory Using Tensor Cores and a Deep Neural Network. Journal of Chemical Theory and Computation, 2022, 18, 4255-4268. | 5.3 | 7 |
| 6 | A QUBO formulation for top- k , eigencentality nodes. PLoS ONE, 2022, 17, e0271292. | 2.5 | 2 |
| 7 | Reduction of the molecular hamiltonian matrix using quantum community detection. Scientific Reports, 2021, 11, 4099. | 3.3 | 11 |
| 8 | Multilevel Combinatorial Optimization across Quantum Architectures. ACM Transactions on Quantum Computing, 2021, 2, 1-29. | 4.3 | 30 |
| 9 | Mixed Precision Fermi-Operator Expansion on Tensor Cores from a Machine Learning Perspective. Journal of Chemical Theory and Computation, 2021, 17, 2256-2265. | 5.3 | 9 |
| 10 | Enabling particle applications for exascale computing platforms. International Journal of High Performance Computing Applications, 2021, 35, 572-597. | 3.7 | 15 |
| 11 | Computing molecular excited states on a D-Wave quantum annealer. Scientific Reports, 2021, 11, 18796. | 3.3 | 16 |
| 12 | Quantum-Based Molecular Dynamics Simulations Using Tensor Cores. Journal of Chemical Theory and Computation, 2021, 17, 6180-6192. | 5.3 | 12 |
| 13 | Detecting multiple communities using quantum annealing on the D-Wave system. PLoS ONE, 2020, 15, e0227538. | 2.5 | 40 |
| 14 | Quantum isomer search. PLoS ONE, 2020, 15, e0226787. | 2.5 | 10 |
| 15 | Using Graph Partitioning for Scalable Distributed Quantum Molecular Dynamics. Algorithms, 2019, 12, 187. | 2.1 | 7 |
| 16 | A Hybrid Approach for Solving Optimization Problems on Small Quantum Computers. Computer, 2019, 52, 18-26. | 1.1 | 38 |
| 17 | Linear Scaling Pseudo Fermi-Operator Expansion for Fractional Occupation. Journal of Chemical Theory and Computation, 2019, 15, 190-200. | 5.3 | 5 |
| 18 | Graph Partitioning as Quadratic Unconstrained Binary Optimization (QUBO) on Spiking Neuromorphic Hardware. , 2019, , . | | 10 |

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 19 | The basic matrix library (BML) for quantum chemistry. <i>Journal of Supercomputing</i> , 2018, 74, 6201-6219. | 3.6 | 12 |
| 20 | Graph Partitioning using Quantum Annealing on the D-Wave System. , 2017, , . | | 99 |
| 21 | Graph-based linear scaling electronic structure theory. <i>Journal of Chemical Physics</i> , 2016, 144, 234101. | 3.0 | 29 |
| 22 | Discrete event performance prediction of speculatively parallel temperature-accelerated dynamics. <i>Simulation</i> , 2016, 92, 1065-1086. | 1.8 | 11 |
| 23 | Recursive Factorization of the Inverse Overlap Matrix in Linear-Scaling Quantum Molecular Dynamics Simulations. <i>Journal of Chemical Theory and Computation</i> , 2016, 12, 3063-3073. | 5.3 | 19 |
| 24 | Integrating predictive analytics into a spatiotemporal epidemic simulation. , 2015, , . | | 15 |
| 25 | TADSim. <i>ACM Transactions on Modeling and Computer Simulation</i> , 2015, 25, 1-26. | 0.8 | 12 |
| 26 | An Evaluation of Threaded Models for a Classical MD Proxy Application. , 2014, , . | | 7 |
| 27 | Optimizing human activity patterns using global sensitivity analysis. <i>Computational and Mathematical Organization Theory</i> , 2014, 20, 394-416. | 2.0 | 8 |
| 28 | Understanding the Impact of Face Mask Usage Through Epidemic Simulation of Large Social Networks. <i>Intelligent Systems Reference Library</i> , 2014, , 97-115. | 1.2 | 17 |
| 29 | Molecular dynamics simulations of detonation on the roadrunner supercomputer. , 2012, , . | | 0 |
| 30 | Self-consistent tight-binding molecular dynamics simulations of shock-induced reactions in hydrocarbons. , 2012, , . | | 2 |
| 31 | Subband coding for large-scale scientific simulation data using JPEG 2000. , 2012, , . | | 2 |
| 32 | Revisiting wavelet compression for large-scale climate data using JPEG 2000 and ensuring data precision. , 2011, , . | | 43 |
| 33 | Real-world hydrologic assessment of a fully-distributed hydrological model in a parallel computing environment. <i>Journal of Hydrology</i> , 2011, 409, 483-496. | 5.4 | 95 |
| 34 | Designing systems for large-scale, discrete-event simulations: Experiences with the FastTrans parallel microsimulator. , 2009, , . | | 13 |
| 35 | Pandemic simulation of antivirals+school closures: buying time until strain-specific vaccine is available. <i>Computational and Mathematical Organization Theory</i> , 2008, 14, 209-221. | 2.0 | 27 |
| 36 | Semi-empirical power-law scaling of new infection rate to model epidemic dynamics with inhomogeneous mixing. <i>Mathematical Biosciences</i> , 2006, 203, 301-318. | 1.9 | 49 |

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 37 | A categorization approach to automated ontological function annotation. Protein Science, 2006, 15, 1544-1549. | 7.6 | 59 |
| 38 | The Gene Ontology Categorizer. Bioinformatics, 2004, 20, i169-i177. | 4.1 | 64 |