

David Doty

List of Publications by Year in descending order

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Version: 2024-02-01

58
papers

1,094
citations

687220

13
h-index

501076

28
g-index

66
all docs

66
docs citations

66
times ranked

534
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|--|------|-----------|
| 1 | Diverse and robust molecular algorithms using reprogrammable DNA self-assembly. <i>Nature</i> , 2019, 567, 366-372. | 13.7 | 198 |
| 2 | Theory of algorithmic self-assembly. <i>Communications of the ACM</i> , 2012, 55, 78-88. | 3.3 | 100 |
| 3 | Deterministic function computation with chemical reaction networks. <i>Natural Computing</i> , 2014, 13, 517-534. | 1.8 | 88 |
| 4 | The Tile Assembly Model is Intrinsically Universal. , 2012, , . | | 77 |
| 5 | Limitations of self-assembly at temperature 1. <i>Theoretical Computer Science</i> , 2011, 412, 145-158. | 0.5 | 49 |
| 6 | Randomized Self-Assembly for Exact Shapes. <i>SIAM Journal on Computing</i> , 2010, 39, 3521-3552. | 0.8 | 33 |
| 7 | Rate-independent computation in continuous chemical reaction networks. , 2014, , . | | 32 |
| 8 | Stable leader election in population protocols requires linear time. <i>Distributed Computing</i> , 2018, 31, 257-271. | 0.7 | 32 |
| 9 | Parallelism and Time in Hierarchical Self-Assembly. , 2012, , . | | 31 |
| 10 | Strong Fault-Tolerance for Self-Assembly with Fuzzy Temperature. , 2010, , . | | 27 |
| 11 | Negative Interactions in Irreversible Self-assembly. <i>Algorithmica</i> , 2013, 66, 153-172. | 1.0 | 26 |
| 12 | Timing in chemical reaction networks. , 2014, , . | | 26 |
| 13 | Speed Faults in Computation by Chemical Reaction Networks. <i>Lecture Notes in Computer Science</i> , 2014, , 16-30. | 1.0 | 22 |
| 14 | Leaderless deterministic chemical reaction networks. <i>Natural Computing</i> , 2015, 14, 213-223. | 1.8 | 21 |
| 15 | Stable Leader Election in Population Protocols Requires Linear Time. <i>Lecture Notes in Computer Science</i> , 2015, , 602-616. | 1.0 | 21 |
| 16 | Deterministic Function Computation with Chemical Reaction Networks. <i>Lecture Notes in Computer Science</i> , 2012, , 25-42. | 1.0 | 20 |
| 17 | Scalable, Time-Responsive, Digital, Energy-Efficient Molecular Circuits Using DNA Strand Displacement. <i>Lecture Notes in Computer Science</i> , 2011, , 25-36. | 1.0 | 18 |
| 18 | Probability 1 computation with chemical reaction networks. <i>Natural Computing</i> , 2016, 15, 245-261. | 1.8 | 17 |

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|----|--|-----|-----------|
| 19 | Randomized Self-Assembly for Exact Shapes. , 2009, , . | | 16 |
| 20 | Negative Interactions in Irreversible Self-assembly. Lecture Notes in Computer Science, 2011, , 37-48. | 1.0 | 13 |
| 21 | Title is missing!. Theory of Computing, 2013, 9, 1-29. | 0.3 | 13 |
| 22 | Finite-state dimension and real arithmetic. Information and Computation, 2007, 205, 1640-1651. | 0.5 | 12 |
| 23 | Fast Algorithmic Self-assembly of Simple Shapes Using Random Agitation. Lecture Notes in Computer Science, 2014, , 20-36. | 1.0 | 12 |
| 24 | A Domain-Specific Language for Programming in the Tile Assembly Model. Lecture Notes in Computer Science, 2009, , 25-34. | 1.0 | 12 |
| 25 | Efficient Size Estimation and Impossibility of Termination in Uniform Dense Population Protocols. , 2019, , . | | 12 |
| 26 | Constructive Dimension and Turing Degrees. Theory of Computing Systems, 2009, 45, 740-755. | 0.7 | 11 |
| 27 | Speed faults in computation by chemical reaction networks. Distributed Computing, 2017, 30, 373-390. | 0.7 | 11 |
| 28 | Random Number Selection in Self-assembly. Lecture Notes in Computer Science, 2009, , 143-157. | 1.0 | 11 |
| 29 | Limitations of Self-assembly at Temperature One. Lecture Notes in Computer Science, 2009, , 35-44. | 1.0 | 11 |
| 30 | Feasible Depth. Lecture Notes in Computer Science, 2007, , 228-237. | 1.0 | 10 |
| 31 | Dimension Extractors and Optimal Decompression. Theory of Computing Systems, 2008, 43, 425-463. | 0.7 | 9 |
| 32 | Thermodynamic Binding Networks. Lecture Notes in Computer Science, 2017, , 249-266. | 1.0 | 8 |
| 33 | Design of Geometric Molecular Bonds. IEEE Transactions on Molecular, Biological, and Multi-Scale Communications, 2017, 3, 13-23. | 1.4 | 7 |
| 34 | Time-Optimal Self-Stabilizing Leader Election in Population Protocols. , 2021, , . | | 7 |
| 35 | Producibility in Hierarchical Self-assembly. Lecture Notes in Computer Science, 2014, , 142-154. | 1.0 | 7 |
| 36 | Constructive Dimension and Weak Truth-Table Degrees. Lecture Notes in Computer Science, 2007, , 63-72. | 1.0 | 7 |

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|----|---|-----|-----------|
| 37 | Producibility in hierarchical self-assembly. <i>Natural Computing</i> , 2016, 15, 41-49. | 1.8 | 6 |
| 38 | Probability 1 Computation with Chemical Reaction Networks. <i>Lecture Notes in Computer Science</i> , 2014, , 37-52. | 1.0 | 6 |
| 39 | Leaderless Deterministic Chemical Reaction Networks. <i>Lecture Notes in Computer Science</i> , 2013, , 46-60. | 1.0 | 5 |
| 40 | Parallelism and Time in Hierarchical Self-Assembly. <i>SIAM Journal on Computing</i> , 2017, 46, 661-709. | 0.8 | 4 |
| 41 | Finite-State Dimension and Real Arithmetic. <i>Lecture Notes in Computer Science</i> , 2006, , 537-547. | 1.0 | 4 |
| 42 | Composable Computation in Discrete Chemical Reaction Networks. , 2019, , . | | 4 |
| 43 | Computational complexity of atomic chemical reaction networks. <i>Natural Computing</i> , 2018, 17, 677-691. | 1.8 | 3 |
| 44 | A time and space optimal stable population protocol solving exact majority. , 2022, , . | | 3 |
| 45 | Designing ordered nucleic acid self-assembly processes. <i>Current Opinion in Structural Biology</i> , 2015, 31, 57-63. | 2.6 | 2 |
| 46 | Program Size and Temperature in Self-Assembly. <i>Algorithmica</i> , 2015, 72, 884-899. | 1.0 | 2 |
| 47 | Democratic, existential, and consensus-based output conventions in stable computation by chemical reaction networks. <i>Natural Computing</i> , 2018, 17, 97-108. | 1.8 | 2 |
| 48 | Composable computation in discrete chemical reaction networks. <i>Distributed Computing</i> , 2021, 34, 437-461. | 0.7 | 2 |
| 49 | Robustness of Expressivity in Chemical Reaction Networks. <i>Lecture Notes in Computer Science</i> , 2016, , 52-66. | 1.0 | 2 |
| 50 | Pushdown dimension. <i>Theoretical Computer Science</i> , 2007, 381, 105-123. | 0.5 | 1 |
| 51 | Design of geometric molecular bonds. , 2016, , . | | 1 |
| 52 | Programming Substrate-Independent Kinetic Barriers With Thermodynamic Binding Networks. <i>IEEE/ACM Transactions on Computational Biology and Bioinformatics</i> , 2021, 18, 283-295. | 1.9 | 1 |
| 53 | Ppsim: A Software Package for Efficiently Simulating and Visualizing Population Protocols. <i>Lecture Notes in Computer Science</i> , 2021, , 245-253. | 1.0 | 1 |
| 54 | A survey of size counting in population protocols. <i>Theoretical Computer Science</i> , 2021, 894, 91-91. | 0.5 | 1 |

| # | ARTICLE | IF | CITATIONS |
|----|--|-----|-----------|
| 55 | Computational Complexity of Atomic Chemical Reaction Networks. Lecture Notes in Computer Science, 2018, , 212-226. | 1.0 | 1 |
| 56 | Brief Announcement: A Time and Space Optimal Stable Population Protocol Solving Exact Majority. , 2021, , . | | 0 |
| 57 | Randomized Self-Assembly. , 2016, , 1759-1767. | | 0 |
| 58 | Hierarchical Self-Assembly. , 2016, , 903-909. | | 0 |