

Dieter Britz

List of Publications by Year in descending order

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Version: 2024-02-01

39
papers

655
citations

687363

13
h-index

580821

25
g-index

43
all docs

43
docs citations

43
times ranked

350
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 1 | Rectangular electrodes: Simulation of accurate steady state currents and the behaviour of square electrode arrays. <i>Electrochimica Acta</i> , 2022, 404, 139750. | 5.2 | 2 |
| 2 | Numerical Convergence Analysis of the Frank-Kamenetskii Equation. <i>Entropy</i> , 2020, 22, 84. | 2.2 | 0 |
| 3 | Revisiting rectangular electrodes; a simulation study. <i>Electrochimica Acta</i> , 2020, 338, 135728. | 5.2 | 7 |
| 4 | Comment on "Atmospheric chemistry of iodine anions: elementary reactions of $I^{\bullet-}$, $IO^{\bullet-}$ and $IO_2^{\bullet-}$ with ozone studied in the gas-phase at 300 K using an ion trap" Teiwes et al., <i>Phys. Chem. Chem. Phys.</i> , 2018, 20, 20608. <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 22654-22655. | 2.8 | 0 |
| 5 | Use of the Saul'yev method for the digital simulation of chronoamperometry at the disk electrode, in the presence of homogeneous chemical reactions. <i>Electrochimica Acta</i> , 2018, 283, 300-305. | 5.2 | 2 |
| 6 | Use of the Saul'yev method for the digital simulation of chronoamperometry and linear sweep voltammetry at the ultramicrodisk electrode. <i>Electrochimica Acta</i> , 2017, 258, 17-23. | 5.2 | 2 |
| 7 | Ketone Body Acetoacetate Buffers Methylglyoxal via a Non-enzymatic Conversion during Diabetic and Dietary Ketosis. <i>Cell Chemical Biology</i> , 2017, 24, 935-943.e7. | 5.2 | 32 |
| 8 | Two (and Three) Dimensions. <i>Monographs in Electrochemistry</i> , 2016, , 251-337. | 0.2 | 0 |
| 9 | Migrational Effects. <i>Monographs in Electrochemistry</i> , 2016, , 339-367. | 0.2 | 0 |
| 10 | Boundary Conditions. <i>Monographs in Electrochemistry</i> , 2016, , 101-121. | 0.2 | 0 |
| 11 | Unequal Intervals. <i>Monographs in Electrochemistry</i> , 2016, , 123-144. | 0.2 | 1 |
| 12 | The Commonly Used Implicit Methods. <i>Monographs in Electrochemistry</i> , 2016, , 145-176. | 0.2 | 0 |
| 13 | Other Methods. <i>Monographs in Electrochemistry</i> , 2016, , 177-234. | 0.2 | 0 |
| 14 | Basic Equations. <i>Monographs in Electrochemistry</i> , 2016, , 5-37. | 0.2 | 0 |
| 15 | Approximations to Derivatives. <i>Monographs in Electrochemistry</i> , 2016, , 39-59. | 0.2 | 0 |
| 16 | Digital Simulation in Electrochemistry. <i>Monographs in Electrochemistry</i> , 2016, , . | 0.2 | 47 |
| 17 | Surface concentration nonuniformities resulting from chronoamperometry of a reversible reaction at an ultramicrodisk electrode. <i>Journal of Electroanalytical Chemistry</i> , 2016, 776, 202-205. | 3.8 | 2 |
| 18 | Digital simulation of chronoamperometry at a disk electrode under a flat polymer film containing an enzyme. <i>Electrochimica Acta</i> , 2015, 152, 302-307. | 5.2 | 13 |

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|----|---|------|-----------|
| 19 | Several ways to simulate time dependent liquid junction potentials by finite differences. <i>Electrochimica Acta</i> , 2014, 137, 328-335. | 5.2 | 19 |
| 20 | Digital simulation of chronoamperometry at an electrode within a hemispherical polymer drop containing an enzyme: Comparison of a hemispherical with a flat disk electrode. <i>Biosensors and Bioelectronics</i> , 2013, 50, 269-277. | 10.1 | 5 |
| 21 | Minimum grid digital simulation of chronoamperometry at a disk electrode. <i>Electrochimica Acta</i> , 2012, 78, 365-376. | 5.2 | 21 |
| 22 | Digital simulation of thermal reactions. <i>Applied Mathematics and Computation</i> , 2011, 218, 1280-1290. | 2.2 | 14 |
| 23 | The true history of adaptive grids in electrochemical simulation. <i>Electrochimica Acta</i> , 2011, 56, 4420-4421. | 5.2 | 6 |
| 24 | Reference values of the chronoamperometric response at cylindrical and capped cylindrical electrodes. <i>Electrochimica Acta</i> , 2010, 55, 5629-5635. | 5.2 | 25 |
| 25 | Diffusion-limited chronoamperometry at conical-tip microelectrodes. <i>Electrochimica Acta</i> , 2010, 55, 1272-1277. | 5.2 | 11 |
| 26 | Strategies for damping the oscillations of the alternating direction implicit method of simulation of diffusion-limited chronoamperometry at disk electrodes. <i>Electrochimica Acta</i> , 2009, 54, 4822-4828. | 5.2 | 9 |
| 27 | The Higher Weight Enumerators of the Doubly-Even, Self-Dual $[48, 24, 12]$ Code. <i>IEEE Transactions on Information Theory</i> , 2007, 53, 2567-2571. | 2.4 | 8 |
| 28 | An interesting global stabilisation of a locally short-range unstable high-order scheme for the digital simulation of the diffusion equation. <i>Computers and Chemical Engineering</i> , 1999, 23, 297-300. | 3.8 | 5 |
| 29 | The effect of the discretization of the mixed boundary conditions on the numerical stability of the Crank-Nicolson algorithm of electrochemical kinetic simulations. <i>Computers & Chemistry</i> , 1997, 21, 391-401. | 1.2 | 14 |
| 30 | Brute force digital simulation. <i>Journal of Electroanalytical Chemistry</i> , 1996, 406, 15-21. | 3.8 | 27 |
| 31 | Numerical stability of finite difference algorithms for electrochemical kinetic simulations: Matrix stability analysis of the classic explicit, fully implicit and Crank-Nicolson methods and typical problems involving mixed boundary conditions. <i>Computers & Chemistry</i> , 1995, 19, 121-136. | 1.2 | 24 |
| 32 | Numerical stability of the Saul'yev finite difference algorithms for electrochemical kinetic simulations: Matrix stability analysis for an example problem involving mixed boundary conditions. <i>Computers & Chemistry</i> , 1995, 19, 357-370. | 1.2 | 14 |
| 33 | Numerical stability of finite difference algorithms for electrochemical kinetic simulations. Matrix stability analysis of the classic explicit, fully implicit and Crank-Nicolson methods, extended to the 3- and 4-point gradient approximation at the electrodes. <i>Computers & Chemistry</i> , 1995, 19, 351-355. | 1.2 | 12 |
| 34 | Some numerical investigations of the stability of electrochemical digital simulation, particularly as affected by first-order homogeneous reactions. <i>Journal of Electroanalytical Chemistry</i> , 1994, 368, 143-147. | 3.8 | 23 |
| 35 | A matter of degree. <i>Nature</i> , 1994, 372, 214-214. | 27.8 | 0 |
| 36 | Electrochemical kinetic simulations of mixed diffusion/homogeneous reaction problems by the Saul'yev finite difference algorithms. <i>Analytica Chimica Acta</i> , 1993, 278, 59-70. | 5.4 | 16 |

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|----|---|-----|-----------|
| 37 | Accuracy contours in (nT, \hat{I}) space in electrochemical digital simulations. Collection of Czechoslovak Chemical Communications, 1991, 56, 20-41. | 1.0 | 6 |
| 38 | Electrochemical digital simulation: re-evaluation of the crank-nicolson scheme. Analytica Chimica Acta, 1987, 194, 317-322. | 5.4 | 7 |
| 39 | Investigation of the relative merits of some n-point current approximations in digital simulation. Analytica Chimica Acta, 1987, 193, 277-285. | 5.4 | 55 |