

Michael F Doherty

List of Publications by Year in Descending Order

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The third column is the impact factor (IF) of the journal, and the fourth column is the number of citations of the article.

171
papers

7,528
citations

52
h-index

78
g-index

180
ext. papers

8,123
ext. citations

4.6
avg, IF

6.12
L-index

| # | Paper | IF | Citations |
|-----|--|------|-----------|
| 171 | Innovation in Chemical Reactor Engineering Practice and Science. <i>Computers and Chemical Engineering</i> , 2022 , 107699 | 4 | |
| 170 | Free energies of crystals computed using Einstein crystal with fixed center of mass and differing spring constants. <i>Journal of Chemical Physics</i> , 2021 , 154, 164509 | 3.9 | 2 |
| 169 | Crystal Morphology Modeling of Solvates and Hydrates of Organic Molecular Crystals: Olanzapine Solvate and Dihydrate. <i>Crystal Growth and Design</i> , 2021 , 21, 4871-4877 | 3.5 | |
| 168 | A New Software Framework for Implementing Crystal Growth Models to Materials of Any Crystallographic Complexity. <i>Crystal Growth and Design</i> , 2020 , 20, 2885-2892 | 3.5 | 4 |
| 167 | Performing solvation free energy calculations in LAMMPS using the decoupling approach. <i>Journal of Computer-Aided Molecular Design</i> , 2020 , 34, 641-646 | 4.2 | 5 |
| 166 | Absolute chemical potentials for complex molecules in fluid phases: A centroid reference for predicting phase equilibria. <i>Journal of Chemical Physics</i> , 2020 , 153, 214504 | 3.9 | 0 |
| 165 | Digital design of crystalline solids. <i>Computers and Chemical Engineering</i> , 2020 , 133, 106637 | 4 | 4 |
| 164 | A rapid screening methodology for chemical processes. <i>Computers and Chemical Engineering</i> , 2020 , 142, 107039 | 4 | 2 |
| 163 | Thermodynamic assessment of carbon dioxide emission reduction during fossil fuel derived energy production. <i>Energy</i> , 2019 , 177, 565-573 | 7.9 | 8 |
| 162 | Ultimate Reaction Selectivities and Screening-Level Metrics for Conceptual Design. <i>Computer Aided Chemical Engineering</i> , 2019 , 329-334 | 0.6 | |
| 161 | Ultimate Reaction Selectivity Limits for Intensified ReactorSeparators. <i>Industrial & Engineering Chemistry Research</i> , 2019 , 58, 6042-6048 | 3.9 | 4 |
| 160 | Ultimate bounds on reaction selectivity for batch reactors. <i>Chemical Engineering Science</i> , 2019 , 199, 652-660 | 4.4 | 6 |
| 159 | LithiumSulfur Batteries: State of the Art and Future Directions. <i>ACS Applied Energy Materials</i> , 2018 , 1, 1783-1814 | 6.1 | 74 |
| 158 | Nonequilibrium Kink Density from One-Dimensional Nucleation for Step Velocity Predictions. <i>Crystal Growth and Design</i> , 2018 , 18, 723-727 | 3.5 | 3 |
| 157 | Ion dissolution mechanism and kinetics at kink sites on NaCl surfaces. <i>Proceedings of the National Academy of Sciences of the United States of America</i> , 2018 , 115, 656-661 | 11.5 | 37 |
| 156 | Target bounds on reaction selectivity via Feinberg's CFSTR equivalence principle. <i>AICHE Journal</i> , 2018 , 64, 926-939 | 3.6 | 13 |
| 155 | Polymorph Selection by Continuous Precipitation. <i>Crystal Growth and Design</i> , 2018 , 18, 4306-4319 | 3.5 | 10 |

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| 154 | Modeling Olanzapine Solution Growth Morphologies. <i>Crystal Growth and Design</i> , 2018 , 18, 905-911 | 3.5 | 21 |
| 153 | In Silico Crystal Growth Rate Prediction for NaCl from Aqueous Solution. <i>Crystal Growth and Design</i> , 2018 , 18, 6302-6306 | 3.5 | 12 |
| 152 | New Tricks of the Trade for Crystal Structure Refinement. <i>ACS Central Science</i> , 2017 , 3, 726-733 | 16.8 | 8 |
| 151 | Modeling Step Velocities and Edge Surface Structures during Growth of Non-Centrosymmetric Crystals. <i>Crystal Growth and Design</i> , 2017 , 17, 2066-2080 | 3.5 | 11 |
| 150 | Modeling layered crystal growth at increasing supersaturation by connecting growth regimes. <i>AIChE Journal</i> , 2017 , 63, 1338-1352 | 3.6 | 24 |
| 149 | Steady State Morphologies of Paracetamol Crystal from Different Solvents. <i>Crystal Growth and Design</i> , 2017 , 17, 659-670 | 3.5 | 24 |
| 148 | A tribute to professor Roger Sargent: Intellectual leader of process systems engineering. <i>AIChE Journal</i> , 2016 , 62, 2951-2958 | 3.6 | 4 |
| 147 | Predicting the Effect of Solvent on the Crystal Habit of Small Organic Molecules. <i>Crystal Growth and Design</i> , 2016 , 16, 2590-2604 | 3.5 | 39 |
| 146 | Polymorph selection by continuous crystallization. <i>AIChE Journal</i> , 2016 , 62, 3505-3514 | 3.6 | 18 |
| 145 | Energetic and entropic components of the Tolman length for mW and TIP4P/2005 water nanodroplets. <i>Journal of Chemical Physics</i> , 2016 , 145, 204703 | 3.9 | 20 |
| 144 | Critical length of a one-dimensional nucleus. <i>Journal of Chemical Physics</i> , 2016 , 145, 211916 | 3.9 | 4 |
| 143 | A design aid for crystal growth engineering. <i>Progress in Materials Science</i> , 2016 , 82, 1-38 | 42.2 | 48 |
| 142 | Rate Expressions for Kink Attachment and Detachment During Crystal Growth. <i>Crystal Growth and Design</i> , 2016 , 16, 3313-3322 | 3.5 | 28 |
| 141 | Designing Robust Crystallization Processes in the Presence of Parameter Uncertainty Using Attainable Regions. <i>Industrial & Engineering Chemistry Research</i> , 2015 , 54, 10350-10363 | 3.9 | 22 |
| 140 | Separation of conglomerate forming enantiomers using a novel continuous preferential crystallization process. <i>AIChE Journal</i> , 2015 , 61, 2810-2823 | 3.6 | 27 |
| 139 | Crystallization of selective polymorph using relationship between supersaturation and solubility. <i>AIChE Journal</i> , 2015 , 61, 1372-1379 | 3.6 | 12 |
| 138 | Regions of attainable particle sizes in continuous and batch crystallization processes. <i>Chemical Engineering Science</i> , 2014 , 106, 167-180 | 4.4 | 91 |
| 137 | Materials science. Imaging crystallization. <i>Science</i> , 2014 , 344, 705-6 | 33.3 | 15 |

- 136 A mechanistic growth model for inorganic crystals: Solid-state interactions. *AIChE Journal*, **2014**, 60, 3703-3719. 3.6 12
- 135 Kink Rate Model for the General Case of Organic Molecular Crystals. *Crystal Growth and Design*, **2014**, 14, 2460-2467. 3.5 16
- 134 Attainable Regions in Crystallization Processes. *Computer Aided Chemical Engineering*, **2014**, 34, 465-470. 0.6 3
- 133 A mechanistic growth model for inorganic crystals: Growth mechanism. *AIChE Journal*, **2014**, 60, 3720-3736. 3.6 17
- 132 Size-Dependent Surface Free Energy and Tolman-Corrected Droplet Nucleation of TIP4P/2005 Water. *Journal of Physical Chemistry Letters*, **2013**, 4, 4267-72. 6.4 54
- 131 Reformulating multidimensional population balances for predicting crystal size and shape. *AIChE Journal*, **2013**, 59, 3468-3474. 3.6 8
- 130 Needle-Shaped Crystals: Causality and Solvent Selection Guidance Based on Periodic Bond Chains. *Crystal Growth and Design*, **2013**, 13, 3341-3352. 3.5 63
- 129 Engineering Crystal Morphology. *Annual Review of Materials Research*, **2013**, 43, 359-386. 12.8 91
- 128 Effect of Structurally Similar Additives on Crystal Habit of Organic Molecular Crystals at Low Supersaturation. *Crystal Growth and Design*, **2013**, 13, 1412-1428. 3.5 38
- 127 Homogeneous nucleation of methane hydrates: unrealistic under realistic conditions. *Journal of the American Chemical Society*, **2012**, 134, 19544-7. 16.4 188
- 126 Predictive Modeling of Supersaturation-Dependent Crystal Shapes. *Crystal Growth and Design*, **2012**, 12, 656-669. 3.5 63
- 125 Crystal shape modification through cycles of dissolution and growth: Attainable regions and experimental validation. *AIChE Journal*, **2012**, 58, 1465-1474. 3.6 39
- 124 Multisite models to determine the distribution of kink sites adjacent to low-energy edges. *Physical Review E*, **2012**, 85, 021604. 2.4 15
- 123 Spiral Growth Model for Faceted Crystals of Non-Centrosymmetric Organic Molecules Grown from Solution. *Crystal Growth and Design*, **2011**, 11, 2780-2802. 3.5 59
- 122 A new framework and a simpler method for the development of batch crystallization recipes. *AIChE Journal*, **2011**, 57, 606-617. 3.6 17
- 121 Analytical Design and Operation of Systems with Crystallization, Filtration, and Recycling. *Industrial & Engineering Chemistry Research*, **2011**, 50, 1196-1205. 3.9 4
- 120 Reinterpreting edge energies calculated from crystal growth experiments. *Journal of Crystal Growth*, **2011**, 327, 117-126. 1.6 13
- 119 Communication: Bubbles, crystals, and laser-induced nucleation. *Journal of Chemical Physics*, **2011**, 134, 171102. 3.9 45

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| 118 | A simulation test of the optical Kerr mechanism for laser-induced nucleation. <i>Journal of Chemical Physics</i> , 2011 , 134, 154501 | 3.9 | 29 |
| 117 | A stochastic model for the critical length of a spiral edge. <i>Journal of Crystal Growth</i> , 2010 , 312, 785-792 | 1.6 | 15 |
| 116 | Reducing the mean size of API crystals by continuous manufacturing with product classification and recycle. <i>Chemical Engineering Science</i> , 2010 , 65, 5770-5780 | 4.4 | 44 |
| 115 | Plantwide dynamics and control of processes with crystallization. <i>Computers and Chemical Engineering</i> , 2010 , 34, 112-121 | 4 | 14 |
| 114 | Predicting crystal growth by spiral motion. <i>Proceedings of the Royal Society A: Mathematical, Physical and Engineering Sciences</i> , 2009 , 465, 1145-1171 | 2.4 | 58 |
| 113 | A New Model for the Effect of Molecular Imposters on the Shape of Faceted Molecular Crystals. <i>Crystal Growth and Design</i> , 2009 , 9, 2637-2645 | 3.5 | 43 |
| 112 | Effect of Competing Reversible Reactions on Optimal Operating Policies for Plants with Recycle. <i>Industrial & Engineering Chemistry Research</i> , 2009 , 48, 8037-8047 | 3.9 | 12 |
| 111 | Estimating diffusivity along a reaction coordinate in the high friction limit: Insights on pulse times in laser-induced nucleation. <i>Journal of Chemical Physics</i> , 2009 , 131, 224112 | 3.9 | 26 |
| 110 | Crystal Shape Engineering. <i>Industrial & Engineering Chemistry Research</i> , 2008 , 47, 9812-9833 | 3.9 | 250 |
| 109 | The Evolution of Crystal Shape During Dissolution: Predictions and Experiments. <i>Crystal Growth and Design</i> , 2008 , 8, 1100-1101 | 3.5 | 39 |
| 108 | Nucleation and polymorph selection in a model colloidal fluid. <i>Physical Review E</i> , 2008 , 77, 041604 | 2.4 | 23 |
| 107 | From form to function: Crystallization of active pharmaceutical ingredients. <i>AIChE Journal</i> , 2008 , 54, 1682-1688 | 3.6 | 205 |
| 106 | Selectivity versus conversion and optimal operating policies for plants with recycle. <i>AIChE Journal</i> , 2008 , 54, 2597-2609 | 3.6 | 3 |
| 105 | Feasible Products for Double-Feed Reactive Distillation Columns. <i>Industrial & Engineering Chemistry Research</i> , 2007 , 46, 3255-3264 | 3.9 | 6 |
| 104 | Faceted crystal shape evolution during dissolution or growth. <i>AIChE Journal</i> , 2007 , 53, 1337-1348 | 3.6 | 65 |
| 103 | Manipulation of crystal shape by cycles of growth and dissolution. <i>AIChE Journal</i> , 2007 , 53, 1510-1517 | 3.6 | 30 |
| 102 | Plantwide operation of processes with crystallization. <i>AIChE Journal</i> , 2007 , 53, 2885-2896 | 3.6 | 13 |
| 101 | Shape evolution of 3-dimensional faceted crystals. <i>AIChE Journal</i> , 2006 , 52, 1906-1915 | 3.6 | 75 |

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| 100 | Choosing an operating policy for seeded batch crystallization. <i>AIChE Journal</i> , 2006 , 52, 2046-2054 | 3.6 | 56 |
| 99 | Steady-State Operating Policies for Plants with Multiple Reactions of Equal Overall Order. <i>Industrial & Engineering Chemistry Research</i> , 2006 , 45, 8056-8062 | 3.9 | 4 |
| 98 | Insight from Economically Optimal Steady-State Operating Policies for Dynamic Plantwide Control. <i>Industrial & Engineering Chemistry Research</i> , 2006 , 45, 1343-1353 | 3.9 | 15 |
| 97 | Novel Reactor Temperature and Recycle Flow Rate Policies for Optimal Process Operation in the Plantwide Context. <i>Industrial & Engineering Chemistry Research</i> , 2005 , 44, 6729-6740 | 3.9 | 10 |
| 96 | Natural gas conversion to liquid fuels in a zone reactor. <i>Catalysis Today</i> , 2005 , 106, 301-304 | 5.3 | 30 |
| 95 | Selection of reference components in reaction invariants. <i>Chemical Engineering Science</i> , 2005 , 60, 7168-7171 | 4.1 | 7 |
| 94 | Experimental study of feasibility in kinetically-controlled reactive distillation. <i>AIChE Journal</i> , 2005 , 51, 464-479 | 3.6 | 2 |
| 93 | Effects of vapor-liquid mass transfer on feasibility of reactive distillation. <i>AIChE Journal</i> , 2004 , 50, 1795-1803 | 3.3 | 2 |
| 92 | Simultaneous prediction of crystal shape and size for solution crystallization. <i>AIChE Journal</i> , 2004 , 50, 2101-2112 | 3.6 | 46 |
| 91 | A dynamic model for evolution of crystal shape. <i>Journal of Crystal Growth</i> , 2004 , 267, 239-250 | 1.6 | 35 |
| 90 | Importance of Process Chemistry in Selecting the Operating Policy for Plants with Recycle. <i>Industrial & Engineering Chemistry Research</i> , 2004 , 43, 3957-3971 | 3.9 | 20 |
| 89 | Evolution of Crystal Shape. <i>Crystal Growth and Design</i> , 2004 , 4, 109-112 | 3.5 | 36 |
| 88 | Feasible Regions for Step-Growth Melt Polycondensation Systems. <i>Industrial & Engineering Chemistry Research</i> , 2004 , 43, 428-440 | 3.9 | 8 |
| 87 | Distillation, Azeotropic, and Extractive 2004 , | | 1 |
| 86 | Simultaneous kinetic resolution of chiral propylene oxide and propylene glycol in a continuous reactive distillation column. <i>Chemical Engineering Science</i> , 2003 , 58, 1289-1300 | 4.4 | 13 |
| 85 | Reactive distillation for methyl acetate production. <i>Computers and Chemical Engineering</i> , 2003 , 27, 1855-1866 | 4.6 | 57 |
| 84 | Synthesis of Azeotropic Distillation Systems with Recycles. <i>Industrial & Engineering Chemistry Research</i> , 2003 , 42, 1783-1794 | 3.9 | 18 |
| 83 | Green chemical engineering aspects of reactive distillation. <i>Environmental Science & Technology</i> , 2003 , 37, 5325-9 | 10.3 | 48 |

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| 82 | Modeling Crystal Shape of Polar Organic Materials: Applications to Amino Acids. <i>Crystal Growth and Design</i> , 2003 , 3, 221-237 | 3.5 | 75 |
| 81 | Design of reactive extraction systems for bioproduct recovery. <i>AIChE Journal</i> , 2002 , 48, 514-526 | 3.6 | 47 |
| 80 | Feasible region for a countercurrent cascade of vapor-liquid CSTRs. <i>AIChE Journal</i> , 2002 , 48, 800-814 | 3.6 | 33 |
| 79 | Feasibility and synthesis of hybrid reactive distillation systems. <i>AIChE Journal</i> , 2002 , 48, 2754-2768 | 3.6 | 14 |
| 78 | Multiple steady states in reactive distillation: kinetic effects. <i>Computers and Chemical Engineering</i> , 2002 , 26, 81-93 | 4 | 72 |
| 77 | Predicting the shape of organic crystals grown from polar solvents. <i>Chemical Engineering Science</i> , 2002 , 57, 1805-1813 | 4.4 | 35 |
| 76 | Reaction Invariants and Mole Balances for Plant Complexes. <i>Industrial & Engineering Chemistry Research</i> , 2002 , 41, 3771-3783 | 3.9 | 9 |
| 75 | A systematic method for reaction invariants and mole balances for complex chemistries. <i>Computers and Chemical Engineering</i> , 2001 , 25, 1199-1217 | 4 | 23 |
| 74 | Effect of chemical kinetics on feasible splits for reactive distillation. <i>AIChE Journal</i> , 2001 , 47, 590-601 | 3.6 | 31 |
| 73 | Modeling the Crystal Shape of Polar Organic Materials: Prediction of Urea Crystals Grown from Polar and Nonpolar Solvents. <i>Crystal Growth and Design</i> , 2001 , 1, 455-461 | 3.5 | 46 |
| 72 | Feasible products for kinetically controlled reactive distillation of ternary mixtures. <i>AIChE Journal</i> , 2000 , 46, 923-936 | 3.6 | 22 |
| 71 | Modeling crystal shapes of organic materials grown from solution. <i>AIChE Journal</i> , 2000 , 46, 1348-1367 | 3.6 | 215 |
| 70 | Simulation of kinetic effects in reactive distillation. <i>Computers and Chemical Engineering</i> , 2000 , 24, 2457-2472 | 4 | 53 |
| 69 | Selectivity Targets for Batch Reactive Distillation. <i>Industrial & Engineering Chemistry Research</i> , 2000 , 39, 1565-1575 | 3.9 | 17 |
| 68 | Reactive Distillation. <i>Industrial & Engineering Chemistry Research</i> , 2000 , 39, 3953-3957 | 3.9 | 195 |
| 67 | Bifurcation study of kinetic effects in reactive distillation. <i>AIChE Journal</i> , 1999 , 45, 546-556 | 3.6 | 46 |
| 66 | A Novel Distillate Policy for Batch Reactive Distillation with Application to the Production of Butyl Acetate. <i>Industrial & Engineering Chemistry Research</i> , 1999 , 38, 714-722 | 3.9 | 66 |
| 65 | Computing All Homogeneous and Heterogeneous Azeotropes in Multicomponent Mixtures. <i>Industrial & Engineering Chemistry Research</i> , 1999 , 38, 4901-4912 | 3.9 | 20 |

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| 64 | Structure of Distillation Regions for Multicomponent Azeotropic Mixtures. <i>AIChE Journal</i> , 1998 , 44, 1382-1391 | 78 |
| 63 | A new technique for predicting the shape of solution-grown organic crystals. <i>AIChE Journal</i> , 1998 , 44, 2501-2514 | 84 |
| 62 | Design for Simultaneous Reaction and Liquid-Liquid Extraction. <i>Industrial & Engineering Chemistry Research</i> , 1998 , 37, 4748-4755 | 33 |
| 61 | Measurement of Residue Curve Maps and Heterogeneous Kinetics in Methyl Acetate Synthesis. <i>Industrial & Engineering Chemistry Research</i> , 1998 , 37, 1917-1928 | 115 |
| 60 | Design Method for Kinetically Controlled, Staged Reactive Distillation Columns. <i>Industrial & Engineering Chemistry Research</i> , 1998 , 37, 2821-2834 | 109 |
| 59 | Attainable regions for reaction with separation. <i>AIChE Journal</i> , 1997 , 43, 374-387 | 80 |
| 58 | Thermodynamic behavior of reactive azeotropes. <i>AIChE Journal</i> , 1997 , 43, 2227-2238 | 49 |
| 57 | Global Stability Analysis and Calculation of Liquid-Liquid Equilibrium in Multicomponent Mixtures. <i>Industrial & Engineering Chemistry Research</i> , 1996 , 35, 1395-1408 | 90 |
| 56 | A Geometric Method for the Design of Liquid Extractors. <i>Industrial & Engineering Chemistry Research</i> , 1996 , 35, 2672-2681 | 22 |
| 55 | A Geometric Design Method for Side-Stream Distillation Columns. <i>Industrial & Engineering Chemistry Research</i> , 1996 , 35, 3653-3664 | 26 |
| 54 | Calculation of residue curve maps for mixtures with multiple equilibrium chemical reactions. <i>Industrial & Engineering Chemistry Research</i> , 1995 , 34, 3195-3202 | 66 |
| 53 | Synthesis of Reactive Distillation Systems with Multiple Equilibrium Chemical Reactions. <i>Industrial & Engineering Chemistry Research</i> , 1995 , 34, 2555-2565 | 95 |
| 52 | Necessary and sufficient conditions for reactive azeotropes in multireaction mixtures. <i>AIChE Journal</i> , 1995 , 41, 2383-2392 | 42 |
| 51 | Vapor-liquid phase equilibrium in systems with multiple chemical reactions. <i>Chemical Engineering Science</i> , 1995 , 50, 23-48 | 165 |
| 50 | Calculation of solid-liquid equilibrium and crystallization paths for melt crystallization processes. <i>Chemical Engineering Science</i> , 1995 , 50, 1679-1694 | 41 |
| 49 | Theory of phase equilibria in multireaction systems. <i>Chemical Engineering Science</i> , 1995 , 50, 3201-3216 | 62 |
| 48 | Design of three-component kinetically controlled reactive distillation columns using fixed-points methods. <i>Chemical Engineering Science</i> , 1994 , 49, 1947-1963 | 48 |
| 47 | Minimum entrainer flows for extractive distillation: A bifurcation theoretic approach. <i>AIChE Journal</i> , 1994 , 40, 243-268 | 84 |

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| 46 | Effect of kinetics on residue curve maps for reactive distillation. <i>AICHE Journal</i> , 1994 , 40, 1814-1824 | 3.6 | 97 |
| 45 | New tools for the design of kinetically controlled reactive distillation columns. <i>Computers and Chemical Engineering</i> , 1994 , 18, S1-S13 | 4 | 10 |
| 44 | Design and operating targets for nonideal multicomponent batch distillation. <i>Industrial & Engineering Chemistry Research</i> , 1993 , 32, 293-301 | 3.9 | 14 |
| 43 | Geometric nonlinear analysis of multicomponent nonideal distillation: a simple computer-aided design procedure. <i>Chemical Engineering Science</i> , 1993 , 48, 1367-1391 | 4.4 | 33 |
| 42 | A new pressure-swing-distillation process for separating homogeneous azeotropic mixtures. <i>Industrial & Engineering Chemistry Research</i> , 1992 , 31, 346-357 | 3.9 | 127 |
| 41 | Feasibility and separation sequencing in multicomponent batch distillation. <i>Chemical Engineering Science</i> , 1991 , 46, 1311-1326 | 4.4 | 72 |
| 40 | Automatic screening of entrainers for homogeneous azeotropic distillation. <i>Industrial & Engineering Chemistry Research</i> , 1991 , 30, 760-772 | 3.9 | 82 |
| 39 | Dynamics of heterogeneous azeotropic distillation columns. <i>AICHE Journal</i> , 1990 , 36, 39-52 | 3.6 | 47 |
| 38 | Thermal integration of homogeneous azeotropic distillation sequences. <i>AICHE Journal</i> , 1990 , 36, 969-984 | 3.6 | 97 |
| 37 | Geometric behavior and minimum flows for nonideal multicomponent distillation. <i>Chemical Engineering Science</i> , 1990 , 45, 1801-1822 | 4.4 | 110 |
| 36 | Design and synthesis of heterogeneous azeotropic distillationsII. Heterogeneous phase diagrams. <i>Chemical Engineering Science</i> , 1990 , 45, 1823-1836 | 4.4 | 56 |
| 35 | Design and synthesis of heterogeneous azeotropic distillationsII. Residue curve maps. <i>Chemical Engineering Science</i> , 1990 , 45, 1837-1843 | 4.4 | 63 |
| 34 | Design and synthesis of heterogeneous azeotropic distillationsIII. Column sequences. <i>Chemical Engineering Science</i> , 1990 , 45, 1845-1854 | 4.4 | 65 |
| 33 | Patterns of composition change in multicomponent batch distillation. <i>Chemical Engineering Science</i> , 1990 , 45, 1207-1221 | 4.4 | 55 |
| 32 | Wiped film reactor model for nylon 6,6 polymerization. <i>Industrial & Engineering Chemistry Research</i> , 1990 , 29, 2012-2020 | 3.9 | 14 |
| 31 | Design and minimum reflux for heterogeneous azeotropic distillation columns. <i>AICHE Journal</i> , 1989 , 35, 1585-1591 | 3.6 | 32 |
| 30 | Design/optimization of ternary heterogeneous azeotropic distillation sequences. <i>AICHE Journal</i> , 1989 , 35, 1592-1601 | 3.6 | 45 |
| 29 | Optimal design and synthesis of homogeneous azeotropic distillation sequences. <i>Industrial & Engineering Chemistry Research</i> , 1989 , 28, 564-572 | 3.9 | 61 |

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| 28 | Approximate dynamic models for chemical process systems. <i>Industrial & Engineering Chemistry Research</i> , 1989 , 28, 546-552 | 3.9 | 19 |
| 27 | Design and minimum-reflux calculations for single-feed multicomponent reactive distillation columns. <i>Chemical Engineering Science</i> , 1988 , 43, 1523-1537 | 4.4 | 139 |
| 26 | The influence of equilibrium chemical reactions on vapor-liquid phase diagrams. <i>Chemical Engineering Science</i> , 1988 , 43, 529-540 | 4.4 | 100 |
| 25 | The simple distillation of homogeneous reactive mixtures. <i>Chemical Engineering Science</i> , 1988 , 43, 541-550 | 4.4 | 138 |
| 24 | Chaos in deterministic systems: Strange attractors, turbulence, and applications in chemical engineering. <i>Chemical Engineering Science</i> , 1988 , 43, 139-183 | 4.4 | 79 |
| 23 | Design and minimum-reflux calculations for double-feed multicomponent reactive distillation columns. <i>Chemical Engineering Science</i> , 1988 , 43, 2377-2389 | 4.4 | 68 |
| 22 | The interface between design and control. 3. Selecting a set of controlled variables. <i>Industrial & Engineering Chemistry Research</i> , 1988 , 27, 611-615 | 3.9 | 40 |
| 21 | The interface between design and control. 1. Process controllability. <i>Industrial & Engineering Chemistry Research</i> , 1988 , 27, 597-605 | 3.9 | 64 |
| 20 | The interface between design and control. 2. Process operability. <i>Industrial & Engineering Chemistry Research</i> , 1988 , 27, 606-611 | 3.9 | 31 |
| 19 | Relative gain array for units in plants with recycle. <i>Industrial & Engineering Chemistry Research</i> , 1987 , 26, 1259-1262 | 3.9 | 37 |
| 18 | Screening of process retrofit alternatives. <i>Industrial & Engineering Chemistry Research</i> , 1987 , 26, 2195-2204 | 3.9 | 21 |
| 17 | Modeling of reactive distillation systems. <i>Industrial & Engineering Chemistry Research</i> , 1987 , 26, 983-989 | 3.9 | 50 |
| 16 | A simple exact method for calculating tangent pinch points in multicomponent nonideal mixtures by bifurcation theory. <i>Chemical Engineering Science</i> , 1986 , 41, 3155-3160 | 4.4 | 23 |
| 15 | Design and synthesis of homogeneous azeotropic distillations. 5. Columns with nonnegligible heat effects. <i>Industrial & Engineering Chemistry Fundamentals</i> , 1986 , 25, 279-289 | | 22 |
| 14 | Design and synthesis of homogeneous azeotropic distillations. 4. Minimum reflux calculations for multiple-feed columns. <i>Industrial & Engineering Chemistry Fundamentals</i> , 1986 , 25, 269-279 | | 31 |
| 13 | The presynthesis problem for homogeneous azeotropic distillation has a unique explicit solution. <i>Chemical Engineering Science</i> , 1985 , 40, 1885-1889 | 4.4 | 22 |
| 12 | Properties of liquid-vapour composition surfaces for multicomponent mixtures with constant latent heat. <i>Chemical Engineering Science</i> , 1985 , 40, 1979-1980 | 4.4 | 11 |
| 11 | On the dynamics of distillation processes. I. batch distillation. <i>Chemical Engineering Science</i> , 1985 , 40, 2087-2093 | 4.4 | 34 |

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| 10 | Effect of overdesign on the operability of distillation columns. <i>Industrial & Engineering Chemistry Process Design and Development</i> , 1985 , 24, 593-598 | | 10 |
| 9 | Separation of closely boiling mixtures by reactive distillation. 2. Experiments. <i>Industrial & Engineering Chemistry Process Design and Development</i> , 1985 , 24, 1071-1073 | | 21 |
| 8 | Separation of closely boiling mixtures by reactive distillation. 1. Theory. <i>Industrial & Engineering Chemistry Process Design and Development</i> , 1985 , 24, 1062-1071 | | 36 |
| 7 | Design and synthesis of homogeneous azeotropic distillations. 3. The sequencing of columns for azeotropic and extractive distillations. <i>Industrial & Engineering Chemistry Fundamentals</i> , 1985 , 24, 474-485 | | 134 |
| 6 | Design and synthesis of homogeneous azeotropic distillations. 1. Problem formulation for a single column. <i>Industrial & Engineering Chemistry Fundamentals</i> , 1985 , 24, 454-463 | | 107 |
| 5 | Design and synthesis of homogeneous azeotropic distillations. 2. Minimum reflux calculations for nonideal and azeotropic columns. <i>Industrial & Engineering Chemistry Fundamentals</i> , 1985 , 24, 463-474 | | 117 |
| 4 | On the dynamics of distillation processes-V. <i>Chemical Engineering Science</i> , 1984 , 39, 883-892 | 4.4 | 26 |
| 3 | An approximate model for binary azeotropic distillation design. <i>Chemical Engineering Science</i> , 1984 , 39, 11-19 | 4.4 | 12 |
| 2 | Controlling thermally linked distillation columns. <i>Industrial & Engineering Chemistry Process Design and Development</i> , 1984 , 23, 483-490 | | 13 |
| 1 | Material stability of multicomponent mixtures and the multiplicity of solutions to phase-equilibrium equations. 1. Nonreacting mixtures. <i>Industrial & Engineering Chemistry Fundamentals</i> , 1983 , 22, 472-485 | | 24 |