

Stuart A Rice

List of Publications by Citations

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1,209
papers

48,920
citations

107
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g-index

1,585
ext. papers

50,674
ext. citations

4.5
avg, IF

7.62
L-index

| # | Paper | IF | Citations |
|------|--|-----|-----------|
| 1209 | Multimode Molecular Dynamics Beyond the Born-Oppenheimer Approximation. <i>Advances in Chemical Physics</i> , 2007 , 59-246 | | 886 |
| 1208 | Control of selectivity of chemical reaction via control of wave packet evolution. <i>Journal of Chemical Physics</i> , 1985 , 83, 5013-5018 | 3.9 | 823 |
| 1207 | Theoretical Aspects of Ionization Potentials and Photoelectron Spectroscopy: A Green's Function Approach. <i>Advances in Chemical Physics</i> , 2007 , 205-344 | | 717 |
| 1206 | Magnetic Relaxation in Fine-Particle Systems. <i>Advances in Chemical Physics</i> , 2007 , 283-494 | | 636 |
| 1205 | Molecular Fluorescence and Energy Transfer Near Interfaces. <i>Advances in Chemical Physics</i> , 2007 , 1-65 | | 613 |
| 1204 | Classical-Limit Quantum Mechanics and the Theory of Molecular Collisions. <i>Advances in Chemical Physics</i> , 2007 , 69-177 | | 610 |
| 1203 | Coherent pulse sequence induced control of selectivity of reactions: Exact quantum mechanical calculations. <i>Journal of Chemical Physics</i> , 1986 , 85, 5805-5820 | 3.9 | 590 |
| 1202 | Fluorescence-detected wave packet interferometry: Time resolved molecular spectroscopy with sequences of femtosecond phase-locked pulses. <i>Journal of Chemical Physics</i> , 1991 , 95, 1487-1511 | 3.9 | 483 |
| 1201 | Angle-Resolved Photoemission as a Tool for the Study of Surfaces. <i>Advances in Chemical Physics</i> , 2007 , 533-656 | | 450 |
| 1200 | Adiabatic Population Transfer with Control Fields. <i>Journal of Physical Chemistry A</i> , 2003 , 107, 9937-9945 | 2.8 | 440 |
| 1199 | Discrete-Variable Representations and their Utilization. <i>Advances in Chemical Physics</i> , 2007 , 263-310 | | 422 |
| 1198 | Semiclassical theory of Bound States. <i>Advances in Chemical Physics</i> , 2007 , 1-61 | | 389 |
| 1197 | Structure and Dynamics of Simple Microclusters. <i>Advances in Chemical Physics</i> , 2007 , 49-135 | | 386 |
| 1196 | Overtone Frequencies and Intensities in the Local Mode Picture. <i>Advances in Chemical Physics</i> , 2007 , 1-58 | | 376 |
| 1195 | Dynamics of Crystal Growth. <i>Advances in Chemical Physics</i> , 2007 , 157-228 | | 368 |
| 1194 | On the Calculation of Time Correlation Functions. <i>Advances in Chemical Physics</i> , 2007 , 63-227 | | 346 |
| 1193 | Dephasing of Molecular Vibrations in Liquids. <i>Advances in Chemical Physics</i> , 2007 , 1-48 | | 342 |

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| 1192 | Active control of the dynamics of atoms and molecules. <i>Annual Review of Physical Chemistry</i> , 1997 , 48, 601-41 | 15.7 | 314 |
| 1191 | Inhomogeneous RF Fields: A Versatile Tool for the Study of Processes with Slow Ions. <i>Advances in Chemical Physics</i> , 2007 , 1-176 | | 314 |
| 1190 | Scattering from a classically chaotic repeller. <i>Journal of Chemical Physics</i> , 1989 , 90, 2225-2241 | 3.9 | 309 |
| 1189 | Correlation Effects in the Ionization of Molecules: Breakdown of the Molecular Orbital Picture. <i>Advances in Chemical Physics</i> , 2007 , 115-159 | | 300 |
| 1188 | The Molecular Quasi-Species. <i>Advances in Chemical Physics</i> , 2007 , 149-263 | | 277 |
| 1187 | On the Calculation of Autocorrelation Functions of Dynamical Variables. <i>Journal of Chemical Physics</i> , 1966 , 45, 1086-1096 | 3.9 | 275 |
| 1186 | Ab Initio Quantum Molecular Dynamics. <i>Advances in Chemical Physics</i> , 2002 , 439-512 | | 260 |
| 1185 | A Critical Assessment of Coupled Cluster Method in Quantum Chemistry. <i>Advances in Chemical Physics</i> , 2007 , 1-175 | | 259 |
| 1184 | Transition Path Sampling. <i>Advances in Chemical Physics</i> , 2003 , 1-78 | | 259 |
| 1183 | Triplet Excitons in Crystals of Aromatic Molecules. <i>Journal of Chemical Physics</i> , 1965 , 42, 309-323 | 3.9 | 255 |
| 1182 | Theory of Ultrafast Nonadiabatic Excited-State Processes and their Spectroscopic Detection in Real Time. <i>Advances in Chemical Physics</i> , 2007 , 1-169 | | 252 |
| 1181 | Radiationless Transitions in Photochemistry. <i>Advances in Photochemistry</i> , 2007 , 149-309 | | 251 |
| 1180 | Electron Transfer from Isolated Molecules to Biomolecules. <i>Advances in Chemical Physics</i> , 2007 , 35-202 | | 247 |
| 1179 | Dielectric Constants of Fluid Models: Statistical Mechanical Theory and its Quantitative Implementation. <i>Advances in Chemical Physics</i> , 2007 , 183-328 | | 238 |
| 1178 | Study of the Lifetimes of Individual Vibronic States of the Isolated Benzene Molecule. <i>Journal of Chemical Physics</i> , 1971 , 55, 5561-5581 | 3.9 | 238 |
| 1177 | Multiconfigurational Perturbation Theory: Applications in Electronic Spectroscopy. <i>Advances in Chemical Physics</i> , 2007 , 219-331 | | 235 |
| 1176 | Assisted adiabatic passage revisited. <i>Journal of Physical Chemistry B</i> , 2005 , 109, 6838-44 | 3.4 | 230 |
| 1175 | Multiparticle Collision Dynamics: Simulation of Complex Systems on Mesoscales. <i>Advances in Chemical Physics</i> , 2008 , 89-146 | | 224 |

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| 1174 | Variational Approaches to Vibration-Rotation Spectroscopy for Polyatomic Molecules. <i>Advances in Chemical Physics</i> , 2007 , 305-379 | | 215 |
| 1173 | Theory of Polyelectrolyte Solutions. <i>Advances in Chemical Physics</i> , 2007 , 1-66 | | 214 |
| 1172 | A Consistent Molecular Treatment of Dielectric Phenomena. <i>Advances in Chemical Physics</i> , 2007 , 467-566 | | 212 |
| 1171 | Kinetic Theory of Dense Fluids. X. Measurement and Interpretation of Self-Diffusion in Liquid Ar, Kr, Xe, and CH ₄ . <i>Journal of Chemical Physics</i> , 1962 , 36, 2710-2720 | 3.9 | 212 |
| 1170 | Roles of Repulsive and Attractive Forces in Liquids : The Equilibrium Theory of Classical Fluids. <i>Advances in Chemical Physics</i> , 2007 , 105-156 | | 211 |
| 1169 | Direct measurements of constrained brownian motion of an isolated sphere between two walls. <i>Physical Review E</i> , 2000 , 62, 3909-19 | 2.4 | 209 |
| 1168 | Classical and Quantum Magnetization Reversal Studied in Nanometer-Sized Particles and Clusters. <i>Advances in Chemical Physics</i> , 2007 , 99-190 | | 206 |
| 1167 | Integral Equation Theories of the Structure, Thermodynamics, and Phase Transitions of Polymer Fluids. <i>Advances in Chemical Physics</i> , 2007 , 1-142 | | 200 |
| 1166 | The Statistical Mechanics of the Electrical Double Layer. <i>Advances in Chemical Physics</i> , 2007 , 141-253 | | 197 |
| 1165 | Functional Integrals and Polymer Statistics. <i>Advances in Chemical Physics</i> , 2007 , 1-128 | | 196 |
| 1164 | Colored Noise in Dynamical Systems. <i>Advances in Chemical Physics</i> , 2007 , 239-326 | | 196 |
| 1163 | Liquids, Glasses, and the Glass Transition: A Free-Volume Approach. <i>Advances in Chemical Physics</i> , 2007 , 455-525 | | 194 |
| 1162 | Study of the Properties of an Excess Electron in Liquid Helium. I. The Nature of the Electron-Helium Interactions. <i>Journal of Chemical Physics</i> , 1965 , 43, 2614-2625 | 3.9 | 186 |
| 1161 | On the Excess Electron and Hole Band Structures and Carrier Mobility in Naphthalene, Anthracene, and Several Polyphenyls. <i>Journal of Chemical Physics</i> , 1963 , 39, 1683-1697 | 3.9 | 185 |
| 1160 | Adiabatic and Quasidiabatic States in a Gauge Theoretical Framework. <i>Advances in Chemical Physics</i> , 2007 , 293-391 | | 180 |
| 1159 | Hard Convex Body Fluids. <i>Advances in Chemical Physics</i> , 2007 , 1-166 | | 179 |
| 1158 | Quantum ergodicity and vibrational relaxation in isolated molecules. <i>Journal of Chemical Physics</i> , 1974 , 61, 203-223 | 3.9 | 174 |
| 1157 | Photophysics of Internal Twisting. <i>Advances in Chemical Physics</i> , 2007 , 1-174 | | 173 |

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| 1156 | Analysis and Evaluation of Ionization Potentials, Electron Affinities, and Excitation Energies by the Equations of Motion Green's Function Method. <i>Advances in Chemical Physics</i> , 2007 , 1-69 | | 169 |
| 1155 | Self-Consistent-Field Methods for Vibrational Excitations in Polyatomic Systems. <i>Advances in Chemical Physics</i> , 2007 , 97-132 | | 168 |
| 1154 | Localized Excitations in Condensed Ne, Ar, Kr, and Xe. <i>Journal of Chemical Physics</i> , 1965 , 42, 4250-4253 | 3.9 | 168 |
| 1153 | On the consistency, extremal, and global properties of counterdiabatic fields. <i>Journal of Chemical Physics</i> , 2008 , 129, 154111 | 3.9 | 167 |
| 1152 | Experimental Study of Luminescence and Excitation Trapping in Vinyl Polymers, Paracyclophanes, and Related Compounds. <i>Journal of Chemical Physics</i> , 1965 , 43, 886-897 | 3.9 | 167 |
| 1151 | Semiclassical quantization of the scattering from a classically chaotic repellor. <i>Journal of Chemical Physics</i> , 1989 , 90, 2242-2254 | 3.9 | 164 |
| 1150 | The Nature and Structural Properties of Graphite Intercalation Compounds. <i>Advances in Chemical Physics</i> , 2007 , 455-532 | | 159 |
| 1149 | Collective Orientational Relaxation in Dense Dipolar Liquids. <i>Advances in Chemical Physics</i> , 2007 , 1-126 | | 157 |
| 1148 | The helix-coil transition in charged macromolecules. <i>Molecular Physics</i> , 1960 , 3, 391-407 | 1.7 | 152 |
| 1147 | Interaction Potentials and Glass Formation: A Survey of Computer Experiments. <i>Advances in Chemical Physics</i> , 2007 , 397-453 | | 151 |
| 1146 | On the Kinetic Theory of Dense Fluids. VI. Singlet Distribution Function for Rigid Spheres with an Attractive Potential. <i>Journal of Chemical Physics</i> , 1961 , 34, 2144-2155 | 3.9 | 149 |
| 1145 | Stress and Structure in Fluid Interfaces. <i>Advances in Chemical Physics</i> , 2007 , 357-454 | | 148 |
| 1144 | Exact quantization of the scattering from a classically chaotic repellor. <i>Journal of Chemical Physics</i> , 1989 , 90, 2255-2262 | 3.9 | 146 |
| 1143 | Theoretical Studies of Transannular Interactions. I. Benzene Excimer Fluorescence and the Singlet States of the Paracyclophanes. <i>Journal of Chemical Physics</i> , 1966 , 44, 23-35 | 3.9 | 146 |
| 1142 | Coherent Pulse Sequence Control of Product Formation in Chemical Reactions. <i>Advances in Chemical Physics</i> , 2007 , 441-523 | | 145 |
| 1141 | Spectral Line Shapes in Gases in the Binary-Collision Approximation. <i>Advances in Chemical Physics</i> , 2007 , 235-293 | | 141 |
| 1140 | Theory and Molecular Models for Water. <i>Advances in Chemical Physics</i> , 2007 , 1-101 | | 140 |
| 1139 | Magnetic Circular Dichroism. <i>Advances in Chemical Physics</i> , 2007 , 197-264 | | 140 |

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| 1138 | Exchange Effects on the Electron and Hole Mobility in Crystalline Anthracene and Naphthalene. <i>Journal of Chemical Physics</i> , 1965 , 42, 733-737 | 3.9 | 140 |
| 1137 | Stability of Complex Reaction Networks. <i>Advances in Chemical Physics</i> , 2007 , 1-215 | | 139 |
| 1136 | The Expansion of the Master Equation. <i>Advances in Chemical Physics</i> , 2007 , 245-309 | | 136 |
| 1135 | The Rotation of Molecules in Dense Phases. <i>Advances in Chemical Physics</i> , 2007 , 1-104 | | 134 |
| 1134 | Path-Integral Centroid Methods in Quantum Statistical Mechanics and Dynamics. <i>Advances in Chemical Physics</i> , 2007 , 135-218 | | 133 |
| 1133 | Irreversible Thermodynamics for Quantum Systems Weakly Coupled to Thermal Reservoirs. <i>Advances in Chemical Physics</i> , 2007 , 109-142 | | 131 |
| 1132 | Models, Interpretations, and Calculations Concerning Resonant Electron Scattering Processes in Atoms and Molecules. <i>Advances in Chemical Physics</i> , 2007 , 91-147 | | 131 |
| 1131 | The Transition from Analytic Dynamics to Statistical Mechanics. <i>Advances in Chemical Physics</i> , 2007 , 155-185 | | 131 |
| 1130 | The OH stretching region infrared spectra of low density amorphous solid water and polycrystalline ice Ih. <i>Journal of Chemical Physics</i> , 1978 , 69, 3477-3482 | 3.9 | 131 |
| 1129 | Spectra of the Alkali Halides. II. The Infrared Spectra of the Sodium and Potassium Halides, RbCl, and CsCl. <i>Journal of Chemical Physics</i> , 1957 , 27, 573-579 | 3.9 | 128 |
| 1128 | Monte Carlo Sampling for Classical Trajectory Simulations. <i>Advances in Chemical Physics</i> , 2007 , 171-201 | | 127 |
| 1127 | Anomalous hydrodynamic interaction in a quasi-two-dimensional suspension. <i>Physical Review Letters</i> , 2004 , 92, 258301 | 7.4 | 127 |
| 1126 | Nonlinear resonance and stochasticity in intramolecular energy exchange. <i>Journal of Chemical Physics</i> , 1976 , 65, 1676-1683 | 3.9 | 127 |
| 1125 | Tests of effective pair potentials for water: Predicted ice structures. <i>Journal of Chemical Physics</i> , 1982 , 76, 650-660 | 3.9 | 125 |
| 1124 | The Redfield Equation in Condensed-Phase Quantum Dynamics. <i>Advances in Chemical Physics</i> , 2007 , 77-134 | | 124 |
| 1123 | Energy Landscapes: From Clusters to Biomolecules. <i>Advances in Chemical Physics</i> , 2007 , 1-111 | | 124 |
| 1122 | Intermolecular and Intramolecular Potentials: Topographical Aspects, Calculation, and Functional Representation via A Double Many-Body Expansion Method. <i>Advances in Chemical Physics</i> , 2007 , 255-338 | | 123 |
| 1121 | A lattice model of a supported monolayer of amphiphile molecules: Monte Carlo simulations. <i>Journal of Chemical Physics</i> , 1988 , 88, 1298-1306 | 3.9 | 123 |

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| 1120 | Star Polymers: Experiment, Theory, and Simulation. <i>Advances in Chemical Physics</i> , 2007 , 67-163 | | 122 |
| 1119 | Bottlenecks to unimolecular reactions and an alternative form for classical RRKM theory. <i>The Journal of Physical Chemistry</i> , 1986 , 90, 3470-3482 | | 122 |
| 1118 | Spectroscopic properties of polyenes. III. 1,3,5,7-Octatetraene. <i>Journal of Chemical Physics</i> , 1978 , 68, 522-529 | 3.9 | 122 |
| 1117 | Internal Rotation and the Breakdown of the Adiabatic Approximation: Many-Phonon Radiationless Transitions. <i>Journal of Chemical Physics</i> , 1970 , 52, 2460-2473 | 3.9 | 119 |
| 1116 | Solid-Liquid Phase Behavior in Microclusters. <i>Advances in Chemical Physics</i> , 2007 , 75-138 | | 118 |
| 1115 | Generalized Langevin Equations and Many-Body Problems in Chemical Dynamics. <i>Advances in Chemical Physics</i> , 2007 , 143-253 | | 117 |
| 1114 | The distribution of rings of hydrogen-bonded molecules in a model of liquid water. <i>Journal of Chemical Physics</i> , 1987 , 86, 5676-5682 | 3.9 | 117 |
| 1113 | Phase transitions in a confined quasi-two-dimensional colloid suspension. <i>Physical Review E</i> , 1997 , 55, 637-656 | 2.4 | 116 |
| 1112 | Natural Chiroptical Spectroscopy: Theory and Computations. <i>Advances in Chemical Physics</i> , 2007 , 545-644 | | 116 |
| 1111 | Radiative and Nonradiative Processes in Benzene. <i>Advances in Chemical Physics</i> , 2007 , 365-421 | | 115 |
| 1110 | Statistical Mechanics of Static and Low-Velocity Kinetic Friction. <i>Advances in Chemical Physics</i> , 2003 , 187-272 | | 115 |
| 1109 | Molecular dynamics studies of the liquid-vapor interface of water. <i>Journal of Chemical Physics</i> , 1991 , 94, 2207-2218 | 3.9 | 114 |
| 1108 | Tethered Polymer Layers. <i>Advances in Chemical Physics</i> , 2007 , 165-260 | | 113 |
| 1107 | The water-water pair potential near the hydrogen bonded equilibrium configuration. <i>Journal of Chemical Physics</i> , 1980 , 72, 3236-3247 | 3.9 | 112 |
| 1106 | Fractal behavior in classical collisional energy transfer. <i>Journal of Chemical Physics</i> , 1986 , 84, 2649-2652 | 3.9 | 110 |
| 1105 | Theory of Inhomogeneous Electron Systems: Spin-Density-Functional Formalism. <i>Advances in Chemical Physics</i> , 2007 , 59-193 | | 108 |
| 1104 | Chemical Reaction Dynamics in Liquid Solution. <i>Advances in Chemical Physics</i> , 2007 , 61-223 | | 108 |
| 1103 | Charge-Transfer Exciton States in Aromatic Molecular Crystals. <i>Journal of Chemical Physics</i> , 1964 , 41, 3294-3306 | 3.9 | 108 |

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| 1102 | Relaxation dynamics of photoexcited benzene-gas van der Waals complexes. <i>Journal of Chemical Physics</i> , 1984 , 81, 1083-1101 | 3.9 | 107 |
| 1101 | Generalized Entropy Theory of Polymer Glass Formation. <i>Advances in Chemical Physics</i> , 2008 , 125-222 | | 105 |
| 1100 | Exciton-Exciton Interactions and Photoconductivity in Crystalline Anthracene. <i>Journal of Chemical Physics</i> , 1963 , 38, 366-373 | 3.9 | 105 |
| 1099 | 1B _{2u} -4A _{1g} spectroscopy of jet-cooled benzene: Single vibronic level fluorescence studies. <i>Journal of Chemical Physics</i> , 1984 , 81, 1060-1072 | 3.9 | 104 |
| 1098 | On the Use of Pseudopotentials in the Quantum Theory of Atoms and Molecules. <i>Advances in Chemical Physics</i> , 2007 , 283-342 | | 100 |
| 1097 | Electron Resonance of Gaseous Diatomic Molecules. <i>Advances in Chemical Physics</i> , 2007 , 149-248 | | 100 |
| 1096 | Molecular Vibration and Nonlinear Optics. <i>Advances in Chemical Physics</i> , 2007 , 1-40 | | 100 |
| 1095 | The mean spherical approximation and effective pair potentials in liquids. <i>Journal of Chemical Physics</i> , 1980 , 72, 4208-4215 | 3.9 | 100 |
| 1094 | On the Singlet-Exciton States of Crystalline Anthracene. <i>Journal of Chemical Physics</i> , 1965 , 42, 1515-1534 | 3.9 | 100 |
| 1093 | A synchrotron x-ray liquid surface spectrometer. <i>Review of Scientific Instruments</i> , 1997 , 68, 4372-4384 | 1.7 | 99 |
| 1092 | On using shaped light pulses to control the selectivity of product formation in a chemical reaction: An application to a multiple level system. <i>Journal of Chemical Physics</i> , 1990 , 93, 1670-1680 | 3.9 | 99 |
| 1091 | Quantum ergodicity and vibrational relaxation in isolated molecules. II. Independent effects and relaxation to the asymptotic limit. <i>Journal of Chemical Physics</i> , 1974 , 61, 768-779 | 3.9 | 99 |
| 1090 | Oscillations and Complex Dynamical Bifurcations in Electrochemical Systems. <i>Advances in Chemical Physics</i> , 2007 , 161-298 | | 98 |
| 1089 | Classical Description of Nonadiabatic Quantum Dynamics. <i>Advances in Chemical Physics</i> , 2005 , 243-375 | | 98 |
| 1088 | On the Statistical Theory of Unimolecular Processes. <i>Advances in Chemical Physics</i> , 2007 , 231-263 | | 97 |
| 1087 | A zeroth order random network model of liquid water. <i>Journal of Chemical Physics</i> , 1979 , 70, 3927-3938 | 3.9 | 97 |
| 1086 | Use of Pseudopotentials in Atomic-Structure Calculations. <i>Journal of Chemical Physics</i> , 1968 , 49, 2741-2755 | 3.9 | 97 |
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| 1084 | Lifetimes and Quantum Yields of Individual Vibronic States of C ₆ D ₆ and C ₆ H ₅ F. <i>Journal of Chemical Physics</i> , 1972 , 56, 2291-2308 | 3.9 | 95 |
| 1083 | Modeling Viral Capsid Assembly. <i>Advances in Chemical Physics</i> , 2014 , 155, 1-68 | | 93 |
| 1082 | Hydrogen Bonds with Large Proton Polarizability and Proton Transfer Processes in Electrochemistry and Biology. <i>Advances in Chemical Physics</i> , 2007 , 1-217 | | 93 |
| 1081 | A molecular dynamics study of the structure of a model Langmuir monolayer of amphiphile molecules. <i>Journal of Chemical Physics</i> , 1988 , 89, 5898-5908 | 3.9 | 93 |
| 1080 | A pseudoatom theory for the liquid-vapor interface of simple metals: Computer simulation studies of sodium and cesium. <i>Journal of Chemical Physics</i> , 1983 , 78, 5225-5249 | 3.9 | 93 |
| 1079 | Study of Impurity-Host Coupling in Shpol'skii Matrices. <i>Journal of Chemical Physics</i> , 1971 , 54, 2014-2023 | 3.9 | 93 |
| 1078 | Statistical Physics of Polymer Solutions: Conformation-Space Renormalization-Group Approach. <i>Advances in Chemical Physics</i> , 2007 , 301-437 | | 92 |
| 1077 | A study of the liquid-vapor interface of mercury: Computer simulation results. <i>Journal of Chemical Physics</i> , 1983 , 78, 5081-5095 | 3.9 | 92 |
| 1076 | Study of the Properties of an Excess Electron in Liquid Helium. II. A Refined Description of Configuration Changes in the Liquid. <i>Journal of Chemical Physics</i> , 1965 , 43, 2625-2632 | 3.9 | 92 |
| 1075 | Superparamagnetism and Spin Glass Dynamics of Interacting Magnetic Nanoparticle Systems. <i>Advances in Chemical Physics</i> , 2004 , 191-248 | | 91 |
| 1074 | Formation of an ordered Langmuir monolayer by a non-polar chain molecule. <i>Nature</i> , 1994 , 367, 151-153 | 5.0 | 91 |
| 1073 | Viscoelastic Subdiffusion: Generalized Langevin Equation Approach. <i>Advances in Chemical Physics</i> , 2012 , 187-253 | | 89 |
| 1072 | Time-Resolved Photoelectron Spectroscopy of Nonadiabatic Dynamics in Polyatomic Molecules. <i>Advances in Chemical Physics</i> , 2008 , 497-584 | | 89 |
| 1071 | Optimization and Characterization of a Multiconfigurational Self-Consistent Field (MCSCF) State. <i>Advances in Chemical Physics</i> , 2007 , 1-176 | | 89 |
| 1070 | Chemical Kinetics of Flue Gas Cleaning by Irradiation with Electrons. <i>Advances in Chemical Physics</i> , 2007 , 315-402 | | 89 |
| 1069 | High-precision molecular wave-packet interferometry with HgAr dimers. <i>Physical Review Letters</i> , 2003 , 91, 243003 | 7.4 | 88 |
| 1068 | Determination of the density profile in the liquid-vapor interface near the triple point. <i>Journal of Chemical Physics</i> , 1978 , 68, 5558-5567 | 3.9 | 88 |
| 1067 | Nonequilibrium Phase Transitions and Chemical Instabilities. <i>Advances in Chemical Physics</i> , 2007 , 311-355 | | 85 |

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| 1066 | Nonequilibrium Fluctuations in Small Systems: From Physics to Biology. <i>Advances in Chemical Physics</i> , 2008 , 31-123 | | 84 |
| 1065 | Theory of Collision-Induced Line Shapes—Absorption and Light Scattering at Low Density. <i>Advances in Chemical Physics</i> , 2007 , 49-112 | | 84 |
| 1064 | Control of Quantum Dynamics by Laser Pulses: Adiabatic Floquet Theory. <i>Advances in Chemical Physics</i> , 2003 , 147-267 | | 83 |
| 1063 | Spectroscopic properties of polyenes. II. The vacuum ultraviolet spectra of cis- and trans-1,3,5-hexatriene. <i>Journal of Chemical Physics</i> , 1974 , 60, 3231-3237 | 3.9 | 81 |
| 1062 | Unimolecular decomposition of the long-lived complex formed in the reaction F+C ₄ H ₈ . <i>Journal of Chemical Physics</i> , 1973 , 59, 1402-1415 | 3.9 | 81 |
| 1061 | Accurate Quantum Chemical Calculations. <i>Advances in Chemical Physics</i> , 2007 , 103-161 | | 80 |
| 1060 | Structure of the liquid-vapor interface of water. <i>Journal of Chemical Physics</i> , 1985 , 82, 4391-4392 | 3.9 | 80 |
| 1059 | The influence of quantization on the onset of chaos in Hamiltonian systems: The Kolmogorov entropy interpretation. <i>Journal of Chemical Physics</i> , 1981 , 74, 1340-1349 | 3.9 | 80 |
| 1058 | Do Exciton States Exist in the Liquid Phase?. <i>Journal of Chemical Physics</i> , 1966 , 44, 4470-4472 | 3.9 | 80 |
| 1057 | Ultrafast Dynamics and Spectroscopy of Bacterial Photosynthetic Reaction Centers. <i>Advances in Chemical Physics</i> , 2002 , 1-88 | | 79 |
| 1056 | Structure and Dynamics of Low-Temperature Water as Studied by Scattering Techniques. <i>Advances in Chemical Physics</i> , 2007 , 1-45 | | 76 |
| 1055 | Electrolytes and the Electric Double Layer. <i>Advances in Chemical Physics</i> , 2007 , 1-159 | | 76 |
| 1054 | Self-consistent Monte Carlo simulations of the electron and ion distributions of inhomogeneous liquid alkali metals. I. Longitudinal and transverse density distributions in the liquid-vapor interface of a one-component system. <i>Journal of Chemical Physics</i> , 1987 , 87, 3069-3081 | 3.9 | 76 |
| 1053 | On the Kinetic Theory of Simple Dense Fluids. XI. Experimental and Theoretical Studies of Positive Ion Mobility in Liquid Ar, Kr, and Xe. <i>Journal of Chemical Physics</i> , 1962 , 37, 947-956 | 3.9 | 76 |
| 1052 | The Theoretical Investigation of the Electron Affinity of Chemical Compounds. <i>Advances in Chemical Physics</i> , 2007 , 169-221 | | 75 |
| 1051 | Phase space bottlenecks and statistical theories of isomerization reactions. <i>Journal of Chemical Physics</i> , 1987 , 86, 2020-2035 | 3.9 | 75 |
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| 1047 | Theory of Radiationless Transitions in an Isolated Molecule. <i>Journal of Chemical Physics</i> , 1968 , 49, 610-623 | 3.9 | 74 |
| 1046 | Reduced Dimensionality Theories of Quantum Reactive Scattering. <i>Advances in Chemical Physics</i> , 2007 , 115-167 | | 73 |
| 1045 | Electron-Impact Spectrometry. <i>Advances in Chemical Physics</i> , 2007 , 15-90 | | 73 |
| 1044 | Polar and Nonpolar Solvation Dynamics, Ion Diffusion, and Vibrational Relaxation: Role of Biphasic Solvent Response in Chemical Dynamics. <i>Advances in Chemical Physics</i> , 2007 , 207-433 | | 73 |
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| 1041 | An improved analysis of the OH stretching region of the vibrational spectrum of ice Ih. <i>Journal of Chemical Physics</i> , 1982 , 77, 583-602 | 3.9 | 71 |
| 1040 | Low-Energy Electron Diffraction. <i>Advances in Chemical Physics</i> , 215-339 | | 71 |
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