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Curriculum Vita

Thomas Nicola Rescigno

PERSONAL

Home Address: 3661 Bernal Avenue, Pleasanton, California 94566
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Birthdate: September 10, 1947
Birthplace: New York City, New York

EDUCATION

1965-1969 Columbia College, New York, New York
B.A. June 1969 (Chemical Physics), graduated summa cum laude

1969-1973 Harvard University, Cambridge, Massachusetts
M.A. June 1971 (Chemistry)
Ph.D., November 1973 (Chemical Physics)
Research Director: William P. Reinhardt
Thesis subject: The Fredholm Method in Electron-Atom Scattering

PROFESSIONAL EXPERIENCE

June 2001 - Present Senior Scientist, Chemical Sciences
Lawrence Berkeley National Laboratory
Berkeley, California 94720

Jan 2000 - June 2001 Staff Scientist, Computing Sciences Directorate
Lawrence Berkeley National Laboratory
Berkeley, California 94720

May 1986 - Oct 2002 Senior Scientist, Physics Directorate
Lawrence Livermore National Laboratory, University of California
Livermore, California 94550

May 1996 - Dec 1999 Participating Guest and Co-Principal Investigator
Computing Sciences Directorate
Lawrence Berkeley National Laboratory
Berkeley, California 94720

Summer 1994 Lecturer, Summer School in Theoretical Atomic and Molecular

Physics

- Summer 1993 University of New Mexico
Los Alamos, New Mexico 87545
- Summer 1991 Visiting Scientist, Los Alamos National Laboratory
Summer 1987 Los Alamos, New Mexico 87545
Los Alamos, New Mexico 87545
- May 1985 - Leader, Theoretical Atomic and Molecular Physics Group
May 1986 Physics Department, Lawrence Livermore National Laboratory
University of California, Livermore, California 94550
- October 1981 - Leader, Atomic and Molecular Theory Group
May 1985 Chemistry and Materials Science Department
Lawrence Livermore National Laboratory, University of California
Livermore, California 94550
- June 1979 - Leader, Theoretical Atomic and Molecular Physics Group
October 1981 Laser Program, Lawrence Livermore National Laboratory
University of California, Livermore, California 94550
- October 1975 - Staff Physicist, Theoretical Atomic and Molecular Physics Group
June 1979 Lawrence Livermore National Laboratory, University of California
Livermore, California 94550
- October 1973 - Research Fellow in Chemistry, California Institute of Technology
September 1975 Pasadena, California 91125
Research Director: Professor Vincent McKoy
- June 1969 - National Science Foundation Trainee, Brookhaven National
September 1969 Laboratory, Upton, New York
Research Director: Dr. Norman Sutin
- May 1969 - National Science Foundation Trainee, Columbia University
September 1968 New York, New York
Research Director: Professor George Flynn

AWARDS AND HONORS

- March 2009 Outstanding Referee, American Physical Society
- October 1988 Fellow, American Physical Society
- April 1975 National Science Foundation Energy Fellowship
- June 1971 Harvard Teaching Fellow Award
- 1969 - 1972 National Science Foundation Predoctoral Fellow
- June 1969 American Institute of Chemists Medal

PROFESSIONAL ACTIVITIES

Research fields: low and intermediate energy electron-atom and electron-molecule collisions, photoionization of atoms and molecules, analyticity and resonance phenomena, atomic and molecular structure, many-body theory, numerical methods, kinetic modeling

Publications:
209 papers in refereed journals
5 review articles
16 chapters in books
1 book

Presentations:
49 invited talks at national and international meetings
26 invited lectures at universities and research institutions
134 technical presentations at scientific meetings

Leadership:**American Physical Society:**

Division of Atomic, Molecular and Optical Physics

Member, Executive Committee (2003-2004, 1996-2000)

Member, Nominating Committee (2003-2004)

Member, Fellowship Committee (2001-2002)

Local Chair, Annual Meeting (1990)

Topical Group on Few-Body Physics

Member, Nominating Committee (2006)

Chair (2003)

Vice-Chair(2002)

Vice-Chair Elect (2001)

Allis Prize Committee

Chair (2001)

Vice-Chair (1999)

Gaseous Electronics Conference

Member, Executive Committee, (2003-2004, 1998-2000)

Physical Review A

Member, Editorial Board, (1995-2001)

National Science Foundation:

AMO Theory Review Panel (2002)

ITR Review Panel (2000)

Review Panel, Institute for Theoretical Atomic and Molecular Physics, Harvard University (1993)

National Research Council Panelist (1989-92)

Department of Energy:

AMO Physics Review Panel, Oak Ridge National Laboratory, (2007)

International Conference on Photonic, Electronic and Atomic Colisions:

Satellite Symposium on Electron- Molecule Collisions
and Swarms (Organizing Committee (2003-2005,
1995-1999), Local Chair (1995))

Journal Review: Physical Review A, Physical Review Letters, Journal of
Chemical Physics, Journal of Physical Chemistry, Journal
of Physics B, Zeitschrift für Physik D

Proposal Review: National Science Foundation
Department of Energy
Petroleum Research Council
Engineering and Physical Sciences Research Council (UK)

PROFESSIONAL AND HONORARY SOCIETIES

American Physical Society
Phi Beta Kappa
Phi Lambda Upsilon

PUBLICATIONS

Thomas Nicola Rescigno

1. Laser Induced Vibrational Fluorescence in Sulphur Hexafluoride (with J. T. Knudton, R. D. Bates, Jr., G. W. Flynn, and M. Ronn), *Bull. Amer. Phys. Soc.* **14**, 65 (1969).
2. Application of the Fredholm Method to Elastic Scattering from Coulomb Admixtures (with W. P. Reinhardt), *Bull. Amer. Phys. Soc.* **16**, 1412 (1971).
3. Extraction of Accurate Scattering Information from Apparently Divergent Fredholm Series: The Centrifugal Barrier (with W. P. Reinhardt), *J. Phys. B* **4**, 1279 (1971).
4. Computation of Elastic Scattering Phase Shifts via Analytic Continuation of Fredholm Determinants Constructed Using an L2 Basis (with W. P. Reinhardt and D. W. Oxtoby), *Phys. Rev. Lett.* **28**, 401 (1972).
5. On Extraction of Scattering Information from Fredholm Determinants Calculated in an L2 Basis: A Chebyshev Discretization of the Continuum (with E. J. Heller and W. P. Reinhardt), *Phys. Rev. A* **8**, 2946 (1973).
6. Calculation of Substituted Fredholm Determinants Using Complex Basis Functions (with W. P. Reinhardt), *Phys. Rev. A* **8**, 2828 (1973).
7. Singlet e-H Scattering at Intermediate Energies (with W. P. Reinhardt), *Phys. Rev. A* **10**, 158 (1974).
8. Calculation of Helium Photoionization in the Random Phase Approximation Using Square-Integrable Basis Functions (with C. W. McCurdy and V. McKoy), *Phys. Rev. A* **9**, 2409 (1974).
9. A Discrete Basis Set Approach to Nonspherical Scattering (with C. W. McCurdy and V. McKoy), *Phys. Rev. A* **10**, 401 (1974).
10. A Discrete Basis Set Approach to Nonspherical Scattering II. (with C. W. McCurdy and V. McKoy), *Phys. Rev. A* **10**, 2240 (1974).
11. Photoionization Cross Sections for H₂ in the Random Phase Approximation with a Square-Integrable Basis (with P. H. S. Martin, V. McKoy and W. H. Henneker), *Chem. Phys. Lett.* **29**, 496 (1974).
12. A Relationship Between the Many-Body Theory of Inelastic Scattering and the Distorted Wave Approximation (with C. W. McCurdy and V. McKoy), *J. Phys. B*

- 7, 2396 (1974).
13. Low-Energy Elastic e⁻-H₂ Cross Sections Using Discrete Basis Functions (with C. W. McCurdy and V. McKoy), *Phys. Rev. A* **11**, 825 (1975).
 14. A Many-Body Treatment of Feshbach Theory Applied to Electron-Atom and Electron-Molecule Collisions (with C. W. McCurdy and V. McKoy), *Phys. Rev. A* **12**, 406 (1975).
 15. A Rigorous Method for Computing Atomic and Molecular Photoionization Cross Sections from a Basis Set Expansion (with V. McKoy), *Phys. Rev. A* **12**, 522 (1975).
 16. Photoabsorption Cross Sections of Two Electron Atoms by the Coordinate Rotation Method: Application to H- and Several States of He (with C. W. McCurdy and V. McKoy), *J. Chem. Phys.* **64**, 477 (1976).
 17. The Equations of Motion Method: An Approach to the Dynamical Properties of Atoms and Molecules (with C. W. McCurdy, V. McKoy, and D. L. Yeager) in "Modern Theoretical Chemistry," edited by H. F. Schaefer, III (Plenum Press, New York, 1977).
 18. Ab Initio Cross Sections for the Excitation of the b³ Σ_u State of H₂ by Electron Impact in the Distorted Wave Approximation (with C. W. McCurdy, V. McKoy and C. F. Bender), *Phys. Rev. A* **13**, 216 (1976).
 19. The Low-Energy Electron Impact Excitation of the Hydrogen Molecule (with C. W. McCurdy, V. McKoy, and C. F. Bender), *Phys. Rev. A* **13**, 216 (1976).
 20. A Simple Method for Evaluating Low-Energy Electron-Molecule Scattering Cross Sections Using Discrete Basis Functions (with C. W. McCurdy and V. McKoy), *J. Phys. B* **9**, 691 (1976).
 21. Cross Sections for the Elastic Scattering of Low-Energy Electrons by Molecular Fluorine: An Approximate Theoretical Treatment Using Discrete Basis Functions (with C. F. Bender, C. W. McCurdy, and V. McKoy), *J. Phys. B* **9**, 2141 (1976).
 22. The Stability of the F₂⁻ Ion: A Model for Dissociative Attachment (with C. F. Bender), *J. Phys. B* **9**, L329 (1976).
 23. A Discrete Basis Set Approach to the Minimum-Variance Method in Electron-Scattering (with D. A. Levin and V. McKoy), *Phys. Rev. A* **16**, 157 (1977).
 24. Theoretical Modeling of the KrF Fluorescence Spectrum (with N. W. Winter), Proceedings of the Third Summer Colloquium on Electronic Transition Lasers, edited by L. E. Wilson, S. N. Suchard, and J. I. Steinfeld, (MIT Press, Cambridge,

- Massachusetts, 1977).
25. The Equations of Motion Method for F₂: Transition Energies, Oscillator Strengths and Born Cross Sections (with C. F. Bender and B. V. McKoy), *Chem. Phys. Lett.* **45**, 307 (1977).
 26. The Photoionization of N₂: An Application of Moment Theory to the Results of a Discrete Basis Set Calculation (with C. F. Bender, V. McKoy, and P. W. Langhoff), *J. Chem. Phys.* **68**, 970 (1978).
 27. Orthogonality Constraints in Electron Scattering by Open-Shell Systems: Comments on a Paper by Riley and Truhlar, *J. Chem. Phys.* **66**, 5255 (1977).
 28. Potential Energy Curves and Predicted Fluorescence for NeF (with N. W. Winter and C. F. Bender), *J. Chem. Phys.* **67**, 3122 (1977).
 29. Comments on the Existence of Low Energy d-Wave Shape Resonances in Electron Fluorine Atom Scattering (with A. U. Hazi and N. W. Winter), *Phys. Rev. A* **16**, 2488 (1977).
 30. A Theoretical Study of the Photoionization of Metastable Neon (with A. U. Hazi), *Phys. Rev. A* **16**, 2376 (1977).
 31. Study of the Photodetachment Cross Section of F⁻ (with C. F. Bender and V. McKoy), *Phys. Rev. A* **17**, 645 (1978).
 32. K-Shell Photoionization in Molecular Nitrogen (with P. W. Langhoff), *Chem. Phys. Lett.* **51**, 65 (1977).
 33. Calculation of the Photoionization Cross Section of the ¹Σ_u Excimer State of Ar₂ (with A. U. Hazi and A. E. Orel), *J. Chem. Phys.* **68**, 5283 (1978).
 34. Extensions of the Complex Coordinate Method to the Study of Resonances in Many-Electron Systems (with C. W. McCurdy and A. E. Orel) *Phys. Rev. A* **17**, 1931 (1978).
 35. Theoretical Studies Involving Candidate Group VIA Laser Systems (with C. F. Bender, A. U. Hazi, A. E. Orel, and N. W. Winter), Proceedings of the Topical Meeting on Inertial Confinement Fusion, San Diego, California, Optical Society of America (1977).
 36. Photoabsorption in Formaldehyde: Intensities and Assignments in the Discrete and Continuous Spectral Intervals (with A. E. Orel, V. McKoy and P. W. Langhoff), *J. Chem. Phys.* **69**, 4689 (1978).
 37. Extension of the Method of Complex Coordinates to Molecular Resonances (with

- C. W. McCurdy), *Phys. Rev. Lett.* **41**, 1364 (1978).
38. Theoretical Study of the 400 eV Core-Valence Excited States of N₂ (with A. E. Orel), *J. Chem. Phys.* **70**, 3390 (1979).
 39. Cross Sections for Excitation of the B³Π_g, C³Π_u, and E³Σ_g States of N₂ by Low-Energy Electron Impact in the Distorted-Wave Approximation (with A. W. Fliflet and V. McKoy), *J. Phys. B* **12**, 3281 (1979).
 40. Electron and Photon-Molecule Collisions (co-edited with V. McKoy and B. Schneider), Plenum Press, New York, 1979.
 41. Potential Energy Curves for Diatomic Zinc and Cadmium (with C. F. Bender, H. F. Schaefer, and A. E. Orel), *J. Chem. Phys.* **71**, 1122 (1979).
 42. Basis Set Calculation of Siegert Eigenvalues: Partial Resonance Widths (with C. W. McCurdy), *Phys. Rev. A* **20**, 2346 (1979).
 43. Photoexcitation and Ionization in Molecular Fluorine: Stieltjes-Tchebycheff Calculations in the Static-Exchange Approximation (with A. E. Orel, B. V. McKoy and P. W. Langhoff), *J. Chem. Phys.* **72**, 1265 (1980).
 44. Theoretical Study of the De-excitation of KrF and XeF Excimers by Low-Energy Electrons (with A. U. Hazi and A. E. Orel), *Appl. Phys. Lett.* **35**, 477 (1979).
 45. On Π-Shell Photoionization in Molecular Nitrogen (with A. Gerwer, B. V. McKoy and P. W. Langhoff), *Chem. Phys. Lett.* **66**, 116 (1979).
 46. Dissociation of F₂ by Electron Impact Excitation of the Lowest 3_u Electronic State (with A. W. Fliflet and V. McKoy), *Phys. Rev. A* **21**, 788 (1980).
 47. The Schwinger Variational Principle for Electron-Molecule Scattering: Applications to Electron-H₂ Scattering (with D. K. Watson, R. R. Lucchese and V. McKoy), *Phys. Rev. A* **21**, 738 (1980).
 48. Complex Basis Function Calculations of Resolvent Matrix Elements: Molecular Photoionization (with C. W. McCurdy), *Phys. Rev. A* **21**, 1499 (1980).
 49. Theoretical Studies of Inner-Valence-Shell Photoionization in N₂ and CO (with P. W. Langhoff, S. R. Langhoff, J. Schirmer, L. S. Cederbaum, W. Domcke and W. Von Niessen), *Chem. Phys.* **58**, 71 (1981).
 50. Applicability of Self-Consistent Field Techniques Based on the Complex Coordinate Method to Metastable Electronic States (with C. W. McCurdy, E. R. Davidson and J. G. Lauderdale), *J. Chem. Phys.* **73**, 3268 (1980).

51. Theoretical Studies of Partial-Channel Photoionization Cross Sections in Diatomic and Polyatomic Molecules (with P. W. Langhoff, N. Padial, G. Csanak and V. McKoy), *Journal de Chimie Physique* **77**, 589 (1980).
52. Generalization of the Separable Approximation to Scattering in the Presence of Long-Range Forces, (with A. E. Orel), *Phys. Rev. A* **23**, 1134 (1981).
53. Application of the Complex Coordinate SCF Technique to a Molecular Shape Resonance: The $^2\Pi_g$ State of N₂, (with A. E. Orel and C. W. McCurdy), *J. of Chem. Phys.* **73**, 6347 (1980).
54. Cross Sections for Resonant Vibrational Excitation of N₂ by Electron Impact, (with A. U. Hazi and M. Kurila), *Phys. Rev. A.* **23**, 1089 (1981).
55. Recent Developments in Complex Scaling, Electron - Atom and Electron - Molecule Collisions, edited by J. Hinze (Plenum, New York, 1981).
56. Schwinger Variational Calculations for Electron Scattering by Polar Molecules, (with D. K. Watson, R. Lucchese and V. McKoy), *J. Phys. B.* **14**, 1875 (1981).
57. Ab Initio Study of Dissociative Attachment of Low-Energy Electrons to F₂, (with A. U. Hazi and A. E. Orel), *Phys. Rev. Letts.* **46**, 918 (1981).
58. A Separable Approximation for Exchange Interactions in Electron-Molecule Scattering, (with A. E. Orel), *Phys. Rev. A* **24**, 1267 (1981).
59. Theoretical Studies of Photoionization in Diatomic and Polyatomic Molecules, (with P. W. Langhoff, N. Padial, C. Csanak, and V. McKoy), *Int. J. Quant. Chem.* **14**, 285 (1980).
60. Theoretical Studies of Photoexcitation and Ionization in H₂O, (with G. Diercksen, W. Kraemer, V. McKoy, C. F. Bender, S. R. Langhoff and P. W. Langhoff), *J. Chem. Phys.* **76**, 1043 (1982).
61. Separable Approximation for Exchange Interactions in Electron-Molecule Scattering: Numerical Stabilization Procedures, (with A. E. Orel), *Phys. Rev. A* **25**, 2402 (1982).
62. Ab Initio Study of Vibrational Excitation of HF by Low Energy Electrons, (with A. E. Orel, A. U. Hazi and B. V. McKoy), *Phys. Rev. A* **26**, 690 (1982).
63. Hole Localization in Molecular Inner Shell Excitation Processes, (with A. E. Orel), in "Wavefunctions and Mechanisms from Electron Scattering Processes," edited by F. Gianturco and G. Stefani, (Springer Verlag, 1984).
64. Studies of Threshold Vibrational Excitation in Electron-Polar Molecule

- Collisions, in "Electron-Molecule Collisions and Photoionization Processes," edited by V. McKoy, H. Suzuki, K. Takayanagi and S. Trajmar, (Verlag Chemie: 1983).
65. Locally Complex Distortions of the Energy Spectrum and the Calculation of Scattering Amplitudes and Photoionization Cross Sections, (with C. W. McCurdy), *Phys. Rev. A* **31**, 624 (1984).
 66. Atomic Photoionization by the Complex Basis Expansion Method: Application to Ground State and Metastable Mg, *Phys. Rev. A* **31**, 607 (1984).
 67. Extension of Time-Independent Wave Operator Methods to the Calculation of the Two-Body Coulomb Amplitude (with B. Johnson, W. P. Reinhardt and C. W. McCurdy), *Phys. Rev. A* **32**, 1998 (1985).
 68. A Note on the Normalization of Resonance Wavefunctions and the Calculation of Resonance Widths (with C. W. McCurdy), *Phys. Rev. A* **34**, 1882 (1986).
 69. A Rigorous Finite Basis Set Approach to the Calculation of the Angular Dependence of Photoionization (with C. W. McCurdy) *Phys. Rev. A* **35**, 657 (1987).
 70. Complex Basis Function Treatment of Photoionization in the Random Phase Approximation (with S. Yabushita and C. W. McCurdy), *Phys. Rev. A* **36**, 3146 (1987).
 71. On the Interrelation Between Variational Methods for Scattering Amplitudes and Generalized R-Matrix Theory (with C. W. McCurdy and B. Schneider), *Phys. Rev. A* **36**, 2061 (1987).
 72. Multichannel Electron-Atom Scattering Calculations Using Local Complex Distortions of the Energy Spectrum (with C. W. McCurdy), *Chem. Phys. Letts.* **140**, 232 (1987).
 73. On the Disappearance of Continuum Exchange Integrals From Algebraic Variational Calculations of Electron Scattering (with B. Schneider) *Phys. Rev. A* **37**, 1044 (1988).
 74. The Complex Kohn Variational Method: Application to Low-Energy Electron Molecule Collisions (with B. Schneider) *Phys. Rev. A* **37**, 3749 (1988).
 75. An Algebraic Variational Approach to the Calculation of Total Cross Sections for One- and Two-Photon Ionization (with A. E. Orel), *Chem. Phys. Letts.* **146**, 434 (1988).
 76. Pseudospectral Techniques in Minimum-Variance Calculations of Electron-

- Scattering Cross Sections (with C. W. McCurdy and B. I. Schneider), *Phys. Rev A* **38**, 5921 (1988).
77. Electron-Impact Excitation of the $b^3\Sigma_u$ State of H₂ Using the Complex Kohn Method: R-Dependence of the Cross Section (with B. I. Schneider), *J. Phys B* **21**, L691 (1988).
 78. Collisions of Electrons with Polyatomic Molecules: Electron-Methane Scattering by the Complex Kohn Variational Method (with. C. W. McCurdy), *Phys. Rev. A* **39**, 4487 (1989).
 79. Accurate Ab Initio Treatment of Low-Energy Electron Collisions with Polyatomic Molecules: Resonant Electron-Formaldehyde Scattering (with. C. W. McCurdy and B. Schneider), *Phys. Rev. Lett.* **63**, 248 (1989).
 80. Minimum Variance Approach to Electron Scattering Calculations: Elimination of Anomalies and Monte Carlo Implementation (with. C. W. McCurdy), *Phys. Rev. A* **40**, 1297 (1989).
 81. New Theoretical Approaches for Studying Electron Collisions with Small Molecules, in "Non-Equilibrium Processes in Partially Ionized Gases", edited by M. Capitelli (Plenum, New York, 1990).
 82. New Developments in the *Ab Initio* Treatment of Low Energy Electron Collisions with Molecules (with B. I. Schneider, C. W. McCurdy and B. H. Lengsfeld III), Proceedings of the ICPEAC Satellite Meeting on Electron-Molecule Collisions, Yale University, July, 1989 (Plenum Press, 1990).
 83. Variational Expressions for First-Order Properties Involving Continuum Wave Functions (with A. E. Orel), *Phys. Rev. A* **41**, 1695 (1990).
 84. Electronic Excitation of Formaldehyde by Low Energy Electrons: A Theoretical Study Using the Complex Kohn Variational Method, T. N. Rescigno, B. H. Lengsfeld, III and C. W. McCurdy, *Phys. Rev. A* **41**, 2462(1990).
 85. Resonant Vibrational Excitation of H₂CO by Low Energy Electron Impact, (with B. I. Schneider and C. W. McCurdy) *Phys. Rev. A* **42**, 3132 (1990).
 86. Theoretical Study of Electron-Impact Excitation of N₂⁺, (with A. E. Orel and B. H. Lengsfeld) *Phys. Rev. A* **42**, 5292 (1990).
 87. Effective Potentials for Time-Dependent Calculations of Multiphoton Processes in Atoms, (with K. C. Kulander) *Comp. Phys. Comm.* **63**, 523 (1991).
 88. Electronic Excitation of H₂ by Electron Impact: Close-coupling calculations

- using the complex Kohn variational method, (with Steven D. Parker, C. William McCurdy and Byron H. Lengsfeld) *Phys. Rev. A*, **43**, 3514 (1991).
89. Continuum Basis Functions in the Complex Kohn Variational Method (with A. E. Orel) *Phys. Rev. A* **43**, 1625 (1991).
 90. Accurate *Ab Initio* Treatment of Low Energy Electron Collisions with Ethylene (with B. I. Schneider and B. H. Lengsfeld) *Phys. Rev. Letts.* **66**, 2728 (1991).
 91. Dissociative Excitation of HeH⁺ by Electron Impact (with A. E. Orel and B. H. Lengsfeld) *Phys. Rev. A* **44**, 4328 (1991).
 92. *Ab Initio* Study of Low Energy Electron-Methane Scattering (with B. H. Lengsfeld and C. W. McCurdy) *Phys. Rev. A* **44**, 4296 (1991).
 93. Electron-Molecule Close-Coupling with Correlated Target Wave Functions: Application to Impact Dissociation of F₂ (with B. H. Lengsfeld) *Phys. Rev. A* **44**, 2913 (1991).
 94. *Ab Initio* Calculations on Collisions of Low Energy Electrons with Polyatomic Molecules, in *Electronic and Atomic Collisions, Invited Papers, Proceedings of the XVII International Conference on the Physics of Electronic and Atomic Collisions*, Brisbane, Australia 1991 (Adam Hilger, Bristol, 1992).
 95. Electron-Impact Excitation of the T and V States of Ethylene: An *Ab Initio* Study (with B. I. Schneider) *Phys Rev A* **45**, 2894 (1992).
 96. Beyond the Primitive Separable Approximation in Electron-Molecule Scattering, (with C. W. McCurdy) *Phys. Rev. A*, **46**, 255 (1992).
 97. *Ab Initio* Description of Polarization in Low-Energy Electron Collisions with Polar Molecules: Application to Electron-NH₃ Scattering, (with B. H. Lengsfeld, C. W. McCurdy and S. D. Parker), *Phys. Rev. A* **45**, 7800 (1992).
 98. A Fixed-Nuclei, *Ab Initio* Treatment of Low-Energy Electron-H₂O Scattering, (with B. H. Lengsfeld), *Zeitschrift fur Physique D*, **24**, 117 (1992).
 99. Polarization and Correlation Effects in Elastic Electron-Li₂ Scattering, (with T. J. Gil, C. W. McCurdy and B. H. Lengsfeld), *Phys. Rev. A*, **47**, 255 (1993).
 100. On the Dependence of Electron Impact Excitation Cross Sections on Initial Vibrational Quantum Number in H₂ and D₂ Molecules, (with R. Celiberto), *Phys. Rev. A*, **47**, 1939 (1993).
 101. An *Ab Initio* Treatment of Near-Threshold Vibrational Excitation of H₂ by Electron Impact: New Perspectives on Discrepancies Between Crossed-Beam and

- Swarm Data, (with B. K. Elza and B. H. Lengsfeld), *J. Phys. B.* **26**, L567 (1993).
102. The Incorporation of Modern Electronic Structure Techniques in Electron-Molecule Collision Problems: Variational Calculations Using the Complex Kohn Method (with B. H. Lengsfeld and C. W. McCurdy), in "Modern Electronic Structure", edited by David Yarkony (World Scientific Publishing, 1995) p.501.
 103. Interchannel Coupling and Ground-State Correlation Effects in the Photoionization of CO (with B. H. Lengsfeld and A. E. Orel), *J. Chem Phys.* **99**, 5097 (1993).
 104. Non-Adiabatic Effects in Low-Energy Electron-Molecule Scattering, in *Electron Collisions with Molecules, Clusters and Surfaces, Proceedings of the XVIII ICPEAC Satellite Conference*, edited by H. Ehrhardt and L. Morgan (Plenum: London, 1993)
 105. *Ab Initio* Complex Kohn Calculations of Dissociative Excitation of Water, (with T. J. Gil, C. W. McCurdy and B. H. Lengsfeld), *Phys. Rev. A* **49**, 2642 (1994).
 106. *Ab Initio* Complex Kohn Calculations of Dissociative Excitation of Methane: Close-Coupling Convergence Studies, (with T. J. Gil, C. W. McCurdy and B. H. Lengsfeld), *Phys. Rev. A* **49**, 2551 (1994).
 107. Low-Energy Electron Collision Processes in Molecular Chlorine, *Phys. Rev. A* **50**, 1382 (1994).
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 110. The Effects of Open Inelastic Channels in the Resonant Dissociative Recombination of HeH⁺ (with A. E. Orel and K. C. Kulander) *Phys. Rev. Letts.* **74**, 4807 (1995).
 111. Low-Energy Electron Collision Processes in NF₃, *Phys. Rev. A* **52**, 329 (1995)
 112. Effective Potential Methods in Variational Treatments of Electron-Molecule Collisions I. Theoretical Formulation (with C. W. McCurdy), *J. Chem. Phys* **104**, 120 (1996).
 113. Effective Potential Methods in Variational Treatments of Electron-Molecule Collisions II. Application to HBr, *J. Chem. Phys* **104**, 125 (1996).

114. Algebraic Variational Approach to Atomic and Molecular Photoionization Cross Sections: Removing the Energy Dependence from the Basis (with A. E. Orel and C. W. McCurdy) *Phys. Rev. A* **55**, 342(1997).
115. Dissociative Excitation in Electron-Molecule Collisions, *Comments At. Mol. Phys.* **33**, 315 (1997).
116. Electron-Molecule Scattering (with C. W. McCurdy), in "Encyclopedia of Computational Chemistry" edited by Paul von Ragué Schleyer (Wiley, New York, 1997).
117. Making Complex Scaling Work for Long Range Potentials, (with M. Baertschy, D. Byrum and C. W. McCurdy) *Phys. Rev. A* **55**, 4253 (1997).
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119. Low-Energy Electron Scattering from CH₃Cl (with A. E. Orel, and C. W. McCurdy), *Phys. Rev. A* **56**, 2855 (1997).
120. An Approach to Electron Impact Ionization that Avoids the Three-Body Coulomb Asymptotic Form (with C. W. McCurdy and D. A. Byrum), *Phys. Rev. A* **56**, 1958 (1997).
121. Calculating Differential Cross Sections for Electron Impact Ionization Without Explicit Use of the Asymptotic Form (with C. W. McCurdy), *Phys. Rev. A* **56**, R4369 (1997).
122. The Calculation of Scattering Amplitudes as Continuous Functions of Energy: R-Matrix Theory without a Box (with C. W. McCurdy, W. A. Isaacs and D. E. Manolopoulos), *Phys. Rev. A* **57**, 3511 (1998).
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