Kenneth D Jordan

List of Publications by Year
in descending order

134 papers	7,748 citations	47006 47 h-index	51608 86 g-index
135 all docs	135 docs citations	135 times ranked	5423 citing authors

#	Article	IF	Citations
1	Progress toward a one-electron model for the non-valence correlation-bound anions of polycyclic aromatic hydrocarbons. Electronic Structure, 2022, 4, 014010.	2.8	1
2	Real-Time Modulation of Hydrogen Evolution Activity of Graphene Electrodes Using Mechanical Strain. ACS Applied Materials & Interfaces, 2022, 14, 10691-10700.	8.0	2
3	Vibrational Signatures of HNO₃ Acidity When Complexed with Microhydrated Alkali Metal Ions, M⁺Â·(HNO₃)(H₂O)_{n=5} (M = Li, K, Na, Rb, Cs), at 20 K. Journal of Physical Chemistry A, 2022, 126, 1640-1647.	2.5	4
4	The binding of atomic hydrogen on graphene from density functional theory and diffusion Monte Carlo calculations. Journal of Chemical Physics, 2022, 156, 144702.	3.0	3
5	Water Network Shape-Dependence of Local Interactions with the Microhydrated â^'NO₂^{â€“} and â^'CO₂^{â€“} Anionic Head Groups by Cold Ion Vibrational Spectroscopy. Journal of Physical Chemistry A, 2022, 126, 2471-2479.	2.5	2
6	A Fresh Look at the Role of the Coupling of a Discrete State with a Pseudocontinuum State in the Stabilization Method for Characterizing Metastable States. Journal of Physical Chemistry Letters, 2021, 12, 1202-1206.	4.6	8
7	Role of Overlap between the Discrete State and Pseudocontinuum States in Stabilization Calculations of Metastable States. Journal of Physical Chemistry A, 2021, 125, 4401-4408.	2.5	3
8	Frontiers of stochastic electronic structure calculations. Journal of Chemical Physics, 2021, 154, 170401.	3.0	5
9	Two-Dimensional Adiabatic Model for Calculating Progressions Resulting from Stretchâ€“Rock Coupling in Vibrational Spectra of Anionâ€“Water Complexes. Journal of Physical Chemistry Letters, 2021, 12, 6326-6329.	4.6	2
10	Temporary Anion Resonances of Pyrene: A 2D Photoelectron Imaging and Computational Study. Journal of Physical Chemistry A, 2021, 125, 7004-7013.	2.5	10
11	Mapping the temperature-dependent and network site-specific onset of spectral diffusion at the surface of a water cluster cage. Proceedings of the National Academy of Sciences of the United States of America, 2020, 117, 26047-26052.	7.1	15
12	The role of high-order electron correlation effects in a model system for non-valence correlation-bound anions. Journal of Chemical Physics, 2020, 153, 224118.	3.0	4
13	Analysis of the Contributions to the Kinetic and Potential Energies of an H Atom in the Presence of a Point Charge: The Molecular Virial Theorem Revisited. Journal of Physical Chemistry A, 2020, 124, 4534-4538.	2.5	0
14	QMCPACK: Advances in the development, efficiency, and application of auxiliary field and real-space variational and diffusion quantum Monte Carlo. Journal of Chemical Physics, 2020, 152, 174105.	3.0	80
15	Toward a systematic improvement of the fixed-node approximation in diffusion Monte Carlo for solidsâ€”A case study in diamond. Journal of Chemical Physics, 2020, 153, 184111.	3.0	16
16	One-Dimensional Adiabatic Model Approach for Calculating Progressions in Vibrational Spectra of Ionâ€“Water Complexes. Journal of Physical Chemistry A, 2019, 123, 7042-7050.	2.5	6
17	Going large(r): general discussion. Faraday Discussions, 2019, 217, 476-513.	3.2	1
18	Controlling internal degrees: general discussion. Faraday Discussions, 2019, 217, 138-171.	3.2	1

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19	Pushing resolution in frequency and time: general discussion. Faraday Discussions, 2019, 217, 290-321.	3.2	1
20	Exotic systems: general discussion. Faraday Discussions, 2019, 217, 601-622.	3.2	0
21	Molecular-level origin of the carboxylate head group response to divalent metal ion complexation at the airâ€“water interface. Proceedings of the National Academy of Sciences of the United States of America, 2019, 116, 14874-14880.	7.1	37
22	Prediction of a Non-Valence Temporary Anion State of (NaCl)₂. Journal of Physical Chemistry B, 2019, 123, 9198-9205.	2.6	3
23	Model potential study of non-valence correlation-bound anions of (C₆₀)_n clusters: the role of electric field-induced charge transfer. Faraday Discussions, 2019, 217, 547-560.	3.2	8
24	Comment on â€œPhysisorption of Water on Graphene: Subchemical Accuracy from Many-Body Electronic Structure Methodsâ€. Journal of Physical Chemistry C, 2019, 123, 10163-10165.	3.1	9
25	Prediction of a Nonvalence Temporary Anion Shape Resonance for a Model (H2O)4 System. Journal of Physical Chemistry A, 2019, 123, 2719-2726.	2.5	4
26	Smallest water clusters supporting the ice I structure. Proceedings of the National Academy of Sciences of the United States of America, 2019, 116, 24383-24385.	7.1	12
27	Tag-Free and Isotopomer-Selective Vibrational Spectroscopy of the Cryogenically Cooled H₉O₄⁺ Cation with Two-Color, IRâ€“IR Double-Resonance Photoexcitation: Isolating the Spectral Signature of a Single OH Group in the Hydronium Ion Core. Journal of Physical Chemistry A, 2018, 122, 9275-9284.	2.5	27
28	Accurate Predictions of Electron Binding Energies of Dipole-Bound Anions via Quantum Monte Carlo Methods. Journal of Physical Chemistry Letters, 2018, 9, 6185-6190.	4.6	24
29	Comment on a spurious prediction of a non-planar geometry for benzene at the MP2 level of theory. Chemical Physics Letters, 2017, 669, 230-232.	2.6	14
30	H4: A model system for assessing the performance of diffusion Monte Carlo calculations using a single Slater determinant trial function. Journal of Chemical Physics, 2017, 147, 074106.	3.0	8
31	Theoretical approaches for treating non-valence correlation-bound anions. Journal of Chemical Physics, 2017, 147, 214114.	3.0	34
32	Preface: Special Topic: From Quantum Mechanics to Force Fields. Journal of Chemical Physics, 2017, 147, 161401.	3.0	13
33	Stabilization calculations of the low-lying temporary anions states of Be, Mg, and Ca. Chemical Physics, 2017, 482, 239-243.	1.9	11
34	Implementation of analytical gradients and of a mixed real and momentum space DVR method for excess electron systems described by a self-consistent polarization model. Journal of Chemical Physics, 2017, 147, 161717.	3.0	2
35	Spectroscopic snapshots of the proton-transfer mechanism in water. Science, 2016, 354, 1131-1135.	12.6	213
36	Diffusion Monte Carlo Study of the Parallel Displaced Form of the Benzene Dimer. ACS Symposium Series, 2016, , 107-117.	0.5	3

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37	Characterization of the primary hydration shell of the hydroxide ion with H2 tagging vibrational spectroscopy of the OHâ˜ â‹¯ (H2O)n=2,3 and ODâ˜ â‹¯ (D2O)n=2,3 clusters. Journal of Chemical Physics, 2016, 145, 134304.	3.0	26
38	Ab initio calculation of the cross sections for electron impact vibrational excitation of CO via the 2Î shape resonance. Journal of Chemical Physics, 2016, 144, 104303.	3.0	7
39	Dispersion dipoles for coupled Drude oscillators. Journal of Chemical Physics, 2016, 144, 034111.	3.0	16
40	Proton-coupled electron transfer in [pyridineÂ·(H2O)]â˜, n= 3, 4, clusters. Chemical Physics Letters, 2016, 661, 196-199.	2.6	1
41	Symmetry-Adapted Perturbation Theory Energy Analysis of Alkyl Fluorineâ€“Aromatic Interactions in Torsion Balance Systems. Journal of Physical Chemistry A, 2016, 120, 9292-9298.	2.5	10
42	Theoretical Characterization of the Minimum-Energy Structure of (SF₆)₂. Journal of Physical Chemistry B, 2016, 120, 1788-1792.	2.6	3
43	Exploration of Brueckner orbital trial wave functions in diffusion Monte Carlo calculations. Chemical Physics Letters, 2016, 644, 117-120.	2.6	4
44	Water network-mediated, electron-induced proton transfer in [C5H5N â‹¯ (H2O)n]â˜ clusters. Journal of Chemical Physics, 2015, 143, 144305.	3.0	8
45	Quantum Monte Carlo calculation of the binding energy of the beryllium dimer. Journal of Chemical Physics, 2015, 143, 084116.	3.0	26
46	Oxygen Atom Exchange between Gaseous CO₂ and TiO₂ Nanoclusters. Journal of Physical Chemistry C, 2015, 119, 3605-3612.	3.1	18
47	Snapshots of Proton Accommodation at a Microscopic Water Surface: Understanding the Vibrational Spectral Signatures of the Charge Defect in Cryogenically Cooled H⁺(H₂O)_{n=2â€“28} Clusters. Journal of Physical Chemistry A, 2015, 119, 9425-9440.	2.5	111
48	Nonvalence Correlation-Bound Anion States of Polycyclic Aromatic Hydrocarbons. Journal of Physical Chemistry Letters, 2015, 6, 3994-3997.	4.6	21
49	Negative electron affinities from conventional electronic structure methods. Theoretical Chemistry Accounts, 2014, 133, 1.	1.4	50
50	Correcting density functionals for dispersion interactions using pseudopotentials. Chemical Physics Letters, 2014, 591, 133-136.	2.6	17
51	Nonvalence Correlation-Bound Anion State of C₆F₆: Doorway to Low-Energy Electron Capture. Journal of Physical Chemistry A, 2014, 118, 7201-7205.	2.5	51
52	Assessing the Performances of Dispersion-Corrected Density Functional Methods for Predicting the Crystallographic Properties of High Nitrogen Energetic Salts. Journal of Chemical Theory and Computation, 2014, 10, 4982-4994.	5.3	22
53	Assessment of Various Electronic Structure Methods for Characterizing Temporary Anion States: Application to the Ground State Anions of N2, C2H2, C2H4, and C6H6. Journal of Physical Chemistry A, 2014, 118, 7489-7497.	2.5	52
54	Nonvalence Correlation-Bound Anion States of Spherical Fullerenes. Nano Letters, 2014, 14, 4602-4606.	9.1	25

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55	Vibrational Spectroscopy of the Waterâ€“Nitrate Complex in the Oâ€“H Stretching Region. Journal of Physical Chemistry A, 2014, 118, 8188-8197.	2.5	39
56	Establishing the Ground State of the Disjoint Diradical Tetramethyleneethane with Quantum Monte Carlo. Journal of the American Chemical Society, 2013, 135, 13862-13869.	13.7	34
57	Correlation Consistent Gaussian Basis Sets for H, Bâ€“Ne with Diracâ€“Fock AREP Pseudopotentials: Applications in Quantum Monte Carlo Calculations. Journal of Chemical Theory and Computation, 2013, 9, 2170-2178.	5.3	27
58	Existence of a Correlation Bound s-Type Anion State of C₆₀. Journal of Physical Chemistry Letters, 2013, 4, 849-853.	4.6	71
59	A Self-Consistent Polarization Potential Model for Describing Excess Electrons Interacting with Water Clusters. Journal of Physical Chemistry B, 2013, 117, 4365-4370.	2.6	34
60	Benchmark Study of the Interaction Energy for an (H₂O)₁₆ Cluster: Quantum Monte Carlo and Complete Basis Set Limit MP2 Results. Journal of Physical Chemistry A, 2013, 117, 7606-7611.	2.5	20
61	An Assessment of the vdW-TS Method for Extended Systems. Journal of Chemical Theory and Computation, 2012, 8, 1503-1513.	5.3	112
62	Benchmark Calculations of the Energies for Binding Excess Electrons to Water Clusters. Journal of Chemical Theory and Computation, 2012, 8, 893-900.	5.3	39
63	Bottom-Up View of Water Network-Mediated CO₂ Reduction Using Cryogenic Cluster Ion Spectroscopy and Direct Dynamics Simulations. Journal of Physical Chemistry A, 2012, 116, 903-912.	2.5	19
64	Determination of conformational preferences in biomolecule mimics with localized orbital coupled cluster methods. , 2012, , .		2
65	A distributed point polarizable force field for carbon dioxide. Theoretical Chemistry Accounts, 2012, 131, 1.	1.4	10
66	From quantum mechanics to force fields: new methodologies for the classical simulation of complex systems. Theoretical Chemistry Accounts, 2012, 131, 1.	1.4	9
67	Unraveling the Anomalous Solvatochromic Response of the Formate Ion Vibrational Spectrum: An Infrared, Ar-Tagging Study of the HCO₂^{Â¯}, DCO₂^{Â¯}, and HCO₂^{Â¯}Â·H₂O Ions. Journal of Physical Chemistry Letters, 2011, 2, 2437-2441.	4.6	49
68	Evaluation of Theoretical Approaches for Describing the Interaction of Water with Linear Acenes. Journal of Physical Chemistry A, 2011, 115, 5955-5964.	2.5	24
69	Downsizing the Hydrated Electron's Lair. Science, 2010, 329, 42-43.	12.6	26
70	Benchmark calculations of waterâ€“acene interaction energies: Extrapolation to the waterâ€“graphene limit and assessment of dispersionâ€“corrected DFT methods. Physical Chemistry Chemical Physics, 2010, 12, 6375.	2.8	111
71	How the Shape of an H-Bonded Network Controls Proton-Coupled Water Activation in HONO Formation. Science, 2010, 327, 308-312.	12.6	99
72	Discrete Variable Representation Implementation of the One-Electron Polarization Model. Journal of Chemical Theory and Computation, 2010, 6, 2388-2394.	5.3	5

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73	A second generation distributed point polarizable water model. Journal of Chemical Physics, 2010, 132, 014309.	3.0	106
74	Isolating the Spectral Signatures of Individual Sites in Water Networks Using Vibrational Double-Resonance Spectroscopy of Cluster Isotopomers. Journal of Physical Chemistry Letters, 2010, 1, 2396-2401.	4.6	31
75	Potential energy landscape of the <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" altimg="si3.gif" display="inline" overflow="scroll"><mml:mrow><mml:mo stretchy="false">(</mml:mo><mml:msub><mml:mrow><mml:mtext>H</mml:mtext></mml:mrow><mml:mrow><mml:mn>2</mml:mn></mml:m	2.6	14
76	Vibrationally Induced Interconversion of H-Bonded NO₂^{â^'}Â·H₂O Isomers within NO₂^{â^'}Â·H₂OÂ·Ar_m Clusters Using IRâ^'IR Pumpâ^'Probe through the OH and NO Stretching Vibrations. Journal of Physical Chemistry A, 2009, 113, 975-981.	2.5	15
77	DF-DFT-SAPT Investigation of the Interaction of a Water Molecule to Coronene and Dodecabenzocoronene: Implications for the Waterâ^'Graphite Interaction. Journal of Physical Chemistry C, 2009, 113, 10242-10248.	3.1	79
78	Parallel tempering Monte Carlo simulations of the water heptamer anion. Chemical Physics Letters, 2008, 455, 135-138.	2.6	18
79	Analytical gradient for geometry optimizations of <mml:math xmlns:mml="http://www.w3.org/1998/Math/MathML" altimg="si28.gif" display="inline" overflow="scroll"><mml:mrow><mml:mo stretchy="false">(</mml:mo><mml:msub><mml:mrow><mml:mtext>H</mml:mtext></mml:mrow><mml:mrow><mml:mn>2</mml:mn></mml:m	2.6	9
80	Model Potential Approaches for Describing the Interaction of Excess Electrons with Water Clusters: Incorporation of Long-Range Correlation Effects. Journal of Physical Chemistry A, 2008, 112, 11021-11035.	2.5	68
81	Exploring the correlation between network structure and electron binding energy in the (H2O)7â^' cluster through isomer-photoselected vibrational predissociation spectroscopy and ab initio calculations: Addressing complexity beyond types I-III. Journal of Chemical Physics, 2008, 128, 104314.	3.0	32
82	Comparison of models with distributed polarizable sites for describing water clusters. Molecular Physics, 2007, 105, 2681-2696.	1.7	42
83	Computational Insight Concerning Catalytic Decision Points of the Transition Metal Catalyzed [2 + 2 + 1] Cyclocarbonylation Reaction of Allenes. Organometallics, 2006, 25, 5204-5206.	2.3	32
84	Interplay between hydrogen bonding and electron solvation on hydratedTiO2(110). Physical Review B, 2006, 73, .	3.2	50
85	Theoretical Characterization of the (H2O)21Cluster:Â Application of ann-body Decomposition Procedureâ€. Journal of Physical Chemistry B, 2006, 110, 18872-18878.	2.6	84
86	Electron Binding Motifs of (H2O)n-Clusters. Journal of the American Chemical Society, 2006, 128, 5828-5833.	13.7	94
87	Low-lying isomers and finite temperature behavior of (H2O)6â^'. Journal of Chemical Physics, 2006, 125, 174301.	3.0	43
88	Spectral Signatures of Hydrated Proton Vibrations in Water Clusters. Science, 2005, 308, 1765-1769.	12.6	712
89	Quantum Drude Oscillator Model for Describing the Interaction of Excess Electrons with Water Clusters:Â An Application to (H2O)13-Ââ€. Journal of Physical Chemistry A, 2005, 109, 11531-11538.	2.5	54
90	Molecular visualization in chemistry education: the role of multidisciplinary collaboration. Chemistry Education Research and Practice, 2005, 6, 136-149.	2.5	66

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91	Wet Electrons at the H2O/TiO2(110) Surface. Science, 2005, 308, 1154-1158.	12.6	239
92	Dipole-bound anions of highly polar molecules: Ethylene carbonate and vinylene carbonate. Journal of Chemical Physics, 2004, 120, 685-690.	3.0	63
93	CHEMISTRY: A Fresh Look at Electron Hydration. Science, 2004, 306, 618-619.	12.6	70
94	Parallel-Tempering Monte Carlo Study of (H2O)n=6-9. Journal of Physical Chemistry A, 2003, 107, 7380-7389.	2.5	65
95	THEORY OFDIPOLE-BOUNDANIONS. Annual Review of Physical Chemistry, 2003, 54, 367-396.	10.8	273
96	An ab initio study of the structure of two-, three- and five-dimersilicon clusters: An approach to the Si(100) surface. Theoretical Chemistry Accounts, 2003, 109, 268-273.	1.4	18
97	Dominant structural motifs of NOâˆ’â‹…(H2O)n complexes: Infrared spectroscopic and ab initio studies. Journal of Chemical Physics, 2003, 118, 4945-4953.	3.0	41
98	Large anharmonic effects in the infrared spectra of the symmetrical CH3NO2âˆ’â‹…(H2O) and CH3CO2âˆ’â‹…(H2O) complexes. Journal of Chemical Physics, 2003, 119, 10138-10145.	3.0	57
99	Dipole-bound anions of carbonyl, nitrile, and sulfoxide containing molecules. Journal of Chemical Physics, 2003, 119, 3650-3660.	3.0	79
100	Rearrangement pathways of the water trimer and tetramer anions. Journal of Chemical Physics, 2002, 116, 3612-3616.	3.0	18
101	Infrared predissociation spectroscopy of I[sup âˆ’]â‹…(CH[sub 3]OH)[sub n], n=1,2: Cooperativity in asymmetric solvation. Journal of Chemical Physics, 2002, 116, 4853.	3.0	41
102	Isolating the Charge-Transfer Component of the Anionic H Bond Via Spin Suppression of the Intracluster Proton Transfer Reaction in the NO-Â·H2O Entrance Channel Complexâ€. Journal of Physical Chemistry A, 2002, 106, 10010-10014.	2.5	24
103	Oxygen Atom Reactions with Circumtrindene and Related Molecules:â€‰ Analogues for the Oxidation of Nanotube Caps. Journal of Physical Chemistry A, 2002, 106, 2572-2579.	2.5	18
104	Theoretical Study of Oxygen Adsorption on Graphite and the (8,0) Single-walled Carbon Nanotube. Journal of Physical Chemistry B, 2001, 105, 11227-11232.	2.6	258
105	An interatomic potential for mercury dimer. Journal of Chemical Physics, 2001, 114, 5545-5551.	3.0	47
106	Resonant ion-dip infrared spectroscopy of benzeneâ€“(water)9: Expanding the cube. Journal of Chemical Physics, 2000, 113, 2290-2303.	3.0	81
107	Theoretical Study of the Low-Lying Electronically Excited States of Diacetylene. Journal of Physical Chemistry A, 2000, 104, 9009-9016.	2.5	15
108	Infrared spectroscopy of negatively charged water clusters: Evidence for a linear network. Journal of Chemical Physics, 1999, 110, 6268-6277.	3.0	116

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109	Near-Threshold Electron-Impact Excitation of the Low-Lying Rydberg States of Ethylene. Journal of Physical Chemistry A, 1999, 103, 5667-5670.	2.5	4
110	Electronic Structure of Dipole-Bound Anions. Journal of Physical Chemistry A, 1998, 102, 2624-2633.	2.5	148
111	CO2â^'Fluorocarbon and CO2â^'Hydrocarbon Interactions from First-Principles Calculations. Journal of Physical Chemistry A, 1998, 102, 2231-2236.	2.5	107
112	Resonant ion-dip infrared spectroscopy of the S4 and D2d water octamers in benzene-(water)8 and benzene2-(water)8. Journal of Chemical Physics, 1998, 109, 6601-6614.	3.0	123
113	LONG-RANGE INTRAMOLECULAR INTERACTIONS: IMPLICATIONS FOR ELECTRON TRANSFER. , 1997, , 257-287.		6
114	Infrared Spectrum of a Molecular Ice Cube: The S4 and D2d Water Octamers in Benzene-(Water)8. Science, 1997, 276, 1678-1681.	12.6	433
115	Energies of dipole-bound anionic states. International Journal of Quantum Chemistry, 1997, 64, 183-191.	2.0	86
116	Theoretical Characterization of the Structures and Vibrational Spectra of Benzeneâ^'(H2O)n(n= 1â^'3) Clusters. The Journal of Physical Chemistry, 1996, 100, 7810-7821.	2.9	187
117	Electronically Excited States of 1,4:5,8-Bismethano-1,4,4a,5,8,8a-hexahydronaphthalene, a Nonconjugated Diene:â€‰ Comparison of Theory and Experiment. Journal of the American Chemical Society, 1996, 118, 1235-1240.	13.7	7
118	Fluorescenceâ€dip infrared spectroscopy of the tropoloneâ€H2O complex. Journal of Chemical Physics, 1996, 105, 2605-2617.	3.0	47
119	Fluorescenceâ€dip infrared spectroscopy of tropolone and tropoloneâ€OD. Journal of Chemical Physics, 1996, 105, 2595-2604.	3.0	75
120	Contribution of electron correlation to the stability of dipole-bound anionic states. Physical Review A, 1996, 54, 1906-1909.	2.5	167
121	Low-Lying Electronically Excited States of CH3Cl:Â Comparison of Theory and Experiment. The Journal of Physical Chemistry, 1996, 100, 5642-5648.	2.9	11
122	Modeâ€selective photoisomerization in 5â€hydroxytropolone. II. Theory. Journal of Chemical Physics, 1995, 102, 5260-5270.	3.0	29
123	An Effective In-Situ O2 High Density Plasma Clean. Materials Research Society Symposia Proceedings, 1993, 315, 273.	0.1	1
124	Calculation of the Siâ€“H bond energies for the monohydride phase of Si(100). Journal of Chemical Physics, 1991, 95, 8652-8654.	3.0	116
125	Assignments of the temporary anion states of the chloromethanes. The Journal of Physical Chemistry, 1990, 94, 5666-5669.	2.9	80
126	Temporary anion states of polyatomic hydrocarbons. Chemical Reviews, 1987, 87, 557-588.	47.7	334

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127	Ab initio electronic structure of anions. Chemical Reviews, 1987, 87, 535-555.	47.7	336
128	Electron Transmission Study of the Splitting of the Ï€* MO's of Angleâ€strained Cyclic Acetylenes: Implications for the Electrophilicity of Alkynes. Bulletin Des SociÃ©tÃ©s Chimiques Belges, 1982, 91, 363-363.	0.0	3
129	Theoretical studies of positronâ€“molecule complexes. Journal of Chemical Physics, 1981, 75, 1876-1887.	3.0	48
130	Bonding in the Mg4 cluster. An example of chemical bonding originating from electron correlation effects. Journal of Chemical Physics, 1981, 75, 1044-1046.	3.0	37
131	Recent developments in electron-molecule scattering. International Journal of Quantum Chemistry, 1981, 20, 331-340.	2.0	1
132	Theoretical studies of positron complexes with atomic anions. Journal of Chemical Physics, 1980, 72, 493-503.	3.0	25
133	Comment on the structure and stability of (CO2)â¯2. Journal of Chemical Physics, 1979, 70, 4422-4424.	3.0	56
134	Studies of the temporary anion states of unsaturated hydrocarbons by electron transmission spectroscopy. Accounts of Chemical Research, 1978, 11, 341-348.	15.6	425