

Martin K Beyer

List of Publications by Year in descending order

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190
papers

7,859
citations

76031

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82
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205
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205
docs citations

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times ranked

6285
citing authors

| # | ARTICLE | IF | CITATIONS |
|----|---|-----|-----------|
| 1 | Formation of Temporary Negative Ions and Their Subsequent Fragmentation upon Electron Attachment to CoQ^0 and CoQ^0H^2 . <i>ChemPhysChem</i> , 2022, 23, e202100834. | 1.0 | 2 |
| 2 | Size-dependent H and H^2 formation by infrared multiple photon dissociation spectroscopy of hydrated vanadium cations, $\text{V}^+(\text{H}^2\text{O})_n$, $n = 3\text{--}51$. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 14699-14708. | 1.3 | 4 |
| 3 | Formation of Temporary Negative Ions and Their Subsequent Fragmentation upon Electron Attachment to CoQ^0 and CoQ^0H^2 . <i>ChemPhysChem</i> , 2022, 23, e202200094. | 1.0 | 0 |
| 4 | Infrared Multiple Photon Dissociation Spectroscopy Confirms Reversible Water Activation in $\text{Mn}^+(\text{H}^2\text{O})_n$, $n = 8$. <i>Journal of Physical Chemistry Letters</i> , 2022, 13, 3269-3275. | 2.1 | 6 |
| 5 | Rearrangement and decomposition pathways of bare and hydrogenated molybdenum oxysulfides in the gas phase. <i>Physical Chemistry Chemical Physics</i> , 2022, 24, 16576-16585. | 1.3 | 4 |
| 6 | Toward Detection of FeH^+ in the Interstellar Medium: Infrared Multiple Photon Dissociation Spectroscopy of Ar^2FeH^+ . <i>Journal of Physical Chemistry Letters</i> , 2022, 13, 5867-5872. | 2.1 | 2 |
| 7 | $[\text{Mo}_3\text{S}_{13}]^{2+}$ as a Model System for Hydrogen Evolution Catalysis by MoS_x : Probing Protonation Sites in the Gas Phase by Infrared Multiple Photon Dissociation Spectroscopy. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 5074-5077. | 7.2 | 18 |
| 8 | Microsolvation of Zn cations: infrared multiple photon dissociation spectroscopy of $\text{Zn}^+(\text{H}^2\text{O})_n$ ($n = 2\text{--}35$). <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 3627-3636. | 1.3 | 8 |
| 9 | Photochemistry and UV/vis spectroscopy of hydrated vanadium cations, $\text{V}^+(\text{H}^2\text{O})_n$, $n = 1\text{--}41$, a model system for photochemical hydrogen evolution. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 22251-22262. | 1.3 | 14 |
| 10 | Spectroscopy and photochemistry of copper nitrate clusters. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 9911-9920. | 1.3 | 3 |
| 11 | $[\text{Mo}_3\text{S}_{13}]^{2+}$ als Modellsystem für die katalytische Wasserstoffentwicklung durch MoS_x : Untersuchung der Protonierungsstellen in der Gasphase durch Infrarot-Mehrphotonendissoziationsspektroskopie. <i>Angewandte Chemie</i> , 2021, 133, 5133-5137. | 1.6 | 1 |
| 12 | Infrared spectroscopy of $\text{CO}_3^-(\text{H}_2\text{O})_{1,2}$ and $\text{CO}_4^-(\text{H}_2\text{O})_{1,2}$. <i>Journal of Chemical Physics</i> , 2021, 154, 084301. | 1.2 | 4 |
| 13 | Activation of a Copper Biscarbene Mechano-Catalyst Using Single-Molecule Force Spectroscopy Supported by Quantum Chemical Calculations. <i>Chemistry - A European Journal</i> , 2021, 27, 8723-8729. | 1.7 | 6 |
| 14 | Functionalization of Diamond-Like Carbon Surfaces to Access High Rupture Forces in Single-Molecule Force Spectroscopy of Covalent Bonds. <i>Chemistry Methods</i> , 2021, 1, 271-277. | 1.8 | 1 |
| 15 | Proton transfer reactivity of molybdenum oxysulfide dianions $[\text{Mo}_2\text{O}_2\text{S}_6]^{2-}$ and $[\text{Mo}_2\text{O}_2\text{S}_5]^{2-}$: The role of Coulomb barriers. <i>International Journal of Mass Spectrometry</i> , 2021, 464, 116558. | 0.7 | 4 |
| 16 | Getting Ready for the Hydrogen Evolution Reaction: The Infrared Spectrum of Hydrated Aluminum Hydride-Hydroxide $\text{AlOH}^+(\text{H}^2\text{O})_n$, $n = 1\text{--}9$. <i>Angewandte Chemie - International Edition</i> , 2021, 60, 16858-16863. | 7.2 | 14 |
| 17 | Auf zur Wasserstoffentwicklung: Das Infrarot-Spektrum von hydratisiertem Aluminiumhydrid-Hydroxid $\text{AlOH}^+(\text{H}_2\text{O})_n$, $n = 9\text{--}14$. <i>Angewandte Chemie</i> , 2021, 133, 16994-16999. | 1.6 | 2 |
| 18 | Asymmetric Solvation of the Zinc Dimer Cation Revealed by Infrared Multiple Photon Dissociation Spectroscopy of $\text{Zn}_2^+(\text{H}_2\text{O})_n$ ($n = 1\text{--}20$). <i>International Journal of Molecular Sciences</i> , 2021, 22, 6026. | 1.8 | 3 |

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|----|---|-----|-----------|
| 19 | Frontispiz: Auf zur Wasserstoffentwicklung: Das Infrarot-Spektrum von hydratisiertem Aluminiumhydrid-Hydroxid $\text{Al}(\text{OH})_2(\text{H}_2\text{O})_n$, $n=9$ 14. Angewandte Chemie, 2021, 133, . | 1.6 | 0 |
| 20 | Frontispiece: Getting Ready for the Hydrogen Evolution Reaction: The Infrared Spectrum of Hydrated Aluminum Hydride-Hydroxide $\text{Al}(\text{OH})_2(\text{H}_2\text{O})_n$, $n=9$ 14. Angewandte Chemie - International Edition, 2021, 60, . | 7.2 | 2 |
| 21 | Photochemical Hydrogen Evolution at Metal Centers Probed with Hydrated Aluminium Cations, $\text{Al}(\text{OH})_2(\text{H}_2\text{O})_n$, $n=1$ 10. Chemistry - A European Journal, 2021, 27, 16367-16376. | 1.7 | 6 |
| 22 | Infrared Multiple Photon Dissociation Spectroscopy of Hydrated Cobalt Anions Doped with Carbon Dioxide $\text{CoCO}_2(\text{H}_2\text{O})_n$, $n=1$ 10, in the $\text{C}=\text{O}$ Stretch Region. Chemistry - A European Journal, 2020, 26, 1074-1081. | 1.7 | 11 |
| 23 | Infrared multiple photon dissociation spectroscopy of anionic copper formate clusters. Journal of Chemical Physics, 2020, 153, 184301. | 1.2 | 5 |
| 24 | Photochemical activation of carbon dioxide in $\text{Mg}^+(\text{CO}_2)(\text{H}_2\text{O})_0,1$. Theoretical Chemistry Accounts, 2020, 139, 127. | 0.5 | 8 |
| 25 | Evidence for lactone formation during infrared multiple photon dissociation spectroscopy of bromoalkanoate doped salt clusters. Physical Chemistry Chemical Physics, 2020, 22, 12028-12038. | 1.3 | 5 |
| 26 | Aktivierung von Kohlenstoffdioxid an Metallzentren: Entwicklung des Ladungstransfers von Mg^+ auf CO_2 in $[\text{MgCO}_2(\text{H}_2\text{O})_n]^+$, $n=0$ 8. Angewandte Chemie, 2020, 132, 7537-7541. | 1.6 | 0 |
| 27 | UV/Vis Spectroscopy of Copper Formate Clusters: Insight into Metal-Ligand Photochemistry. Chemistry - A European Journal, 2020, 26, 8286-8295. | 1.7 | 10 |
| 28 | Carbon Dioxide Activation at Metal Centers: Evolution of Charge Transfer from Mg^+ to CO_2 in $[\text{MgCO}_2(\text{H}_2\text{O})_n]^+$, $n=0$ 8. Angewandte Chemie - International Edition, 2020, 59, 7467-7471. | 7.2 | 24 |
| 29 | Chemistry of NO_x and HNO_3 Molecules with Gas-Phase Hydrated $\text{O}^{\cdot-}$ and $\text{OH}^{\cdot-}$ Ions. Chemistry - A European Journal, 2020, 26, 7861-7868. | 1.7 | 5 |
| 30 | Controlling internal degrees: general discussion. Faraday Discussions, 2019, 217, 138-171. | 1.6 | 1 |
| 31 | Pushing resolution in frequency and time: general discussion. Faraday Discussions, 2019, 217, 290-321. | 1.6 | 1 |
| 32 | Exotic systems: general discussion. Faraday Discussions, 2019, 217, 601-622. | 1.6 | 0 |
| 33 | Probing the Structural Evolution of the Hydrated Electron in Water Cluster Anions $(\text{H}_2\text{O})_n^{\cdot-}$, $n=200$, by Electronic Absorption Spectroscopy. Journal of the American Chemical Society, 2019, 141, 18000-18003. | 6.6 | 18 |
| 34 | Structural Properties of Gas-Phase Molybdenum Oxide Clusters $[\text{Mo}_4\text{O}_{13}]^{2+}$, $[\text{HMo}_4\text{O}_{13}]^+$, and $[\text{CH}_3\text{Mo}_4\text{O}_{13}]^+$ Studied by Collision-Induced Dissociation. Journal of the American Society for Mass Spectrometry, 2019, 30, 1946-1955. | 1.2 | 7 |
| 35 | Electronic spectroscopy and nanocalorimetry of hydrated magnesium ions $[\text{Mg}(\text{H}_2\text{O})_n]^+$, $n=20$ 70: spontaneous formation of a hydrated electron?. Faraday Discussions, 2019, 217, 584-600. | 1.6 | 20 |
| 36 | Carbon-carbon bond formation in the reaction of hydrated carbon dioxide radical anions with 3-butyne-1-ol. International Journal of Mass Spectrometry, 2019, 435, 101-106. | 0.7 | 6 |

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|----|---|-----|-----------|
| 37 | Infrared Spectroscopy of Size-Selected Hydrated Carbon Dioxide Radical Anions $\text{CO}_2^{\cdot-}(\text{H}_2\text{O})_n$ ($n=61$) in the $\text{C}\ddot{\text{O}}$ Stretch Region. <i>Chemistry - A European Journal</i> , 2019, 25, 10165-10171. | 1.7 | 27 |
| 38 | Mechanical Activation Drastically Accelerates Amide Bond Hydrolysis, Matching Enzyme Activity. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 9787-9790. | 7.2 | 37 |
| 39 | Mechanische Aktivierung beschleunigt die Hydrolyse der Amidbindung drastisch, vergleichbar der Aktivität von Enzymen. <i>Angewandte Chemie</i> , 2019, 131, 9890-9894. | 1.6 | 6 |
| 40 | Release of Formic Acid from Copper Formate: Hydride, Proton-Coupled Electron and Hydrogen Atom Transfer All Play their Role. <i>ChemPhysChem</i> , 2019, 20, 1420-1424. | 1.0 | 17 |
| 41 | Decomposition of Copper Formate Clusters: Insight into Elementary Steps of Calcination and Carbon Dioxide Activation. <i>ChemistryOpen</i> , 2019, 8, 1453-1459. | 0.9 | 11 |
| 42 | Considerable matrix shift in the electronic transitions of helium-solvated cesium dimer cation Cs_2He_n . <i>Physical Chemistry Chemical Physics</i> , 2019, 21, 25362-25368. | 1.3 | 7 |
| 43 | Personal foreword on Helmut Schwarz Honor Issue. <i>International Journal of Mass Spectrometry</i> , 2019, 436, 23-24. | 0.7 | 1 |
| 44 | Structural Properties of Gas Phase Molybdenum Sulfide Clusters $[\text{Mo}_3\text{S}_{13}]^{2+}$, $[\text{HMo}_3\text{S}_{13}]^+$, and $[\text{H}_3\text{Mo}_3\text{S}_{13}]^{+}$ as Model Systems of a Promising Hydrogen Evolution Catalyst. <i>Journal of Physical Chemistry C</i> , 2019, 123, 8177-8186. | 1.5 | 24 |
| 45 | CO_2/O_2 Exchange in Magnesium-Water Clusters $\text{Mg}^+(\text{H}_2\text{O})_n$. <i>Journal of Physical Chemistry A</i> , 2019, 123, 73-81. | 1.1 | 11 |
| 46 | Infrared multiple photon dissociation of cesium iodide clusters doped with mono-, di- and triglycine. <i>European Journal of Mass Spectrometry</i> , 2019, 25, 122-132. | 0.5 | 9 |
| 47 | Kinetics of the reaction of $\text{CO}_3\ddot{\text{E}}^{\text{TM}}(\text{H}_2\text{O})_n$, $n = 0, 1, 2$, with nitric acid, a key reaction in tropospheric negative ion chemistry. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 10838-10845. | 1.3 | 6 |
| 48 | Isomeric Broadening of C_{60}^+ Electronic Excitation in Helium Droplets: Experiments Meet Theory. <i>Journal of Physical Chemistry Letters</i> , 2018, 9, 1237-1242. | 2.1 | 26 |
| 49 | Gas-Phase Reactivity Studies of Small Molybdenum Cluster Ions with Dimethyl Disulfide. <i>Topics in Catalysis</i> , 2018, 61, 20-27. | 1.3 | 10 |
| 50 | Photochemistry of glyoxylate embedded in sodium chloride clusters, a laboratory model for tropospheric sea-salt aerosols. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 8143-8151. | 1.3 | 22 |
| 51 | Photochemistry and spectroscopy of small hydrated magnesium clusters $\text{Mg}+(\text{H}_2\text{O})_n$, $n = 1-5$. <i>Journal of Chemical Physics</i> , 2018, 149, 044309. | 1.2 | 28 |
| 52 | Photodissociation of Sodium Iodide Clusters Doped with Small Hydrocarbons. <i>Chemistry - A European Journal</i> , 2018, 24, 12433-12443. | 1.7 | 16 |
| 53 | Photodissociation spectroscopy of protonated leucine enkephalin. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 10786-10795. | 1.3 | 29 |
| 54 | Electron-triggered chemistry in $\text{HNO}_3/\text{H}_2\text{O}$ complexes. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 11753-11758. | 1.3 | 28 |

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|----|---|-----|-----------|
| 55 | Hydration Leads to Efficient Reactions of the Carbonate Radical Anion with Hydrogen Chloride in the Gas Phase. <i>Journal of Physical Chemistry A</i> , 2017, 121, 192-197. | 1.1 | 5 |
| 56 | Communication: Charge transfer dominates over proton transfer in the reaction of nitric acid with gas-phase hydrated electrons. <i>Journal of Chemical Physics</i> , 2017, 147, 101101. | 1.2 | 11 |
| 57 | Infrared spectroscopy of $\text{O}^{\ominus}\text{H}^{\ominus}$ and OH^{\ominus} in water clusters: evidence for fast interconversion between $\text{O}^{\ominus}\text{H}^{\ominus}$ and OH^{\ominus} . <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 25346-25351. | 1.3 | 21 |
| 58 | Evidence for Electron Transfer in the Reactions of Hydrated Monovalent First-Row Transition-Metal Ions $\text{M}(\text{H}_2\text{O})_n^+$, $\text{M} = \text{V}, \text{Cr}, \text{Mn}, \text{Fe}, \text{Co}, \text{Ni}, \text{Cu}$, and Zn , $n < 40$, toward 1-Iodopropane. <i>Journal of Physical Chemistry A</i> , 2017, 121, 9557-9566. | 1.1 | 10 |
| 59 | Do protons recombine with O_2^{\ominus} and CO_2^{\ominus} in water clusters?. <i>International Journal of Mass Spectrometry</i> , 2017, 418, 101-106. | 0.7 | 12 |
| 60 | Theoretical simulation of the infrared signature of mechanically stressed polymer solids. <i>Beilstein Journal of Organic Chemistry</i> , 2017, 13, 1710-1716. | 1.3 | 6 |
| 61 | Charge transfer reactions between gas-phase hydrated electrons, molecular oxygen and carbon dioxide at temperatures of 80–300 K. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 23528-23537. | 1.3 | 24 |
| 62 | Electrons Mediate the Gas-Phase Oxidation of Formic Acid with Ozone. <i>Chemistry - A European Journal</i> , 2016, 22, 12684-12687. | 1.7 | 14 |
| 63 | Atomically resolved phase transition of fullerene cations solvated in helium droplets. <i>Nature Communications</i> , 2016, 7, 13550. | 5.8 | 84 |
| 64 | Mechanochemical Cycloreversion of Cyclobutane Observed at the Single Molecule Level. <i>Chemistry - A European Journal</i> , 2016, 22, 12034-12039. | 1.7 | 34 |
| 65 | The reaction of CF_2Cl_2 with gas-phase hydrated electrons. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 23910-23915. | 1.3 | 9 |
| 66 | Force dependence of the infrared spectra of polypropylene calculated with density functional theory. <i>Polymer Degradation and Stability</i> , 2016, 128, 294-299. | 2.7 | 9 |
| 67 | Chemical Reactivity on Gas-Phase Metal Clusters Driven by Blackbody Infrared Radiation. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 1357-1360. | 7.2 | 19 |
| 68 | Lokalisierung eines mechanochemischen Bindungsbruchs durch Einbettung des Mechanophors in einen Makrocyclus. <i>Angewandte Chemie</i> , 2015, 127, 2587-2590. | 1.6 | 14 |
| 69 | Reactivity of Hydrated Monovalent First Row Transition Metal Ions $[\text{M}(\text{H}_2\text{O})_n]^+$, $\text{M} = \text{Cr}, \text{Mn}, \text{Fe}, \text{Co}, \text{Ni}, \text{Cu}$, and Zn , $n < 50$, Toward Acetonitrile. <i>Journal of Physical Chemistry A</i> , 2015, 119, 5566-5578. | 1.1 | 13 |
| 70 | Radiation Effects in Solid Nitrogen and Nitrogen-Containing Matrices: Fingerprints of N_4^+ Species. <i>Journal of Physical Chemistry A</i> , 2015, 119, 2475-2482. | 1.1 | 24 |
| 71 | Pinpointing Mechanochemical Bond Rupture by Embedding the Mechanophore into a Macrocycle. <i>Angewandte Chemie - International Edition</i> , 2015, 54, 2556-2559. | 7.2 | 47 |
| 72 | Thermochemistry of the Reaction of SF_6 with Gas-Phase Hydrated Electrons: A Benchmark for Nanocalorimetry. <i>Journal of Physical Chemistry A</i> , 2015, 119, 9978-9985. | 1.1 | 37 |

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|----|---|-----|-----------|
| 73 | Mechanically induced silyl ester cleavage under acidic conditions investigated by AFM-based single-molecule force spectroscopy in the force-ramp mode. <i>Faraday Discussions</i> , 2014, 170, 357-367. | 1.6 | 18 |
| 74 | Electron transfer reactions of nitromethane, acetaldehyde and benzaldehyde with $(\text{H}_2\text{O})_n^-$, CO_2^- , and O_2^- in the gas phase. <i>International Journal of Mass Spectrometry</i> , 2014, 365-366, 295-300. | 0.7 | 11 |
| 75 | A density functional theory model of mechanically activated silyl ester hydrolysis. <i>Journal of Chemical Physics</i> , 2014, 140, 044321. | 1.2 | 18 |
| 76 | Carboxylation of Methyl Acrylate by Carbon Dioxide Radical Anions in Gas-Phase Water Clusters. <i>Angewandte Chemie - International Edition</i> , 2013, 52, 9327-9330. | 7.2 | 28 |
| 77 | CC bond formation between CO_2^- and allyl alcohol: A mechanistic study. <i>International Journal of Mass Spectrometry</i> , 2013, 354-355, 175-180. | 0.7 | 16 |
| 78 | Collisional Activation of N_2O Decomposition and CO Oxidation Reactions on Isolated Rhodium Clusters. <i>Journal of Physical Chemistry A</i> , 2013, 117, 8855-8863. | 1.1 | 28 |
| 79 | Reduction of Acetonitrile by Hydrated Magnesium Cations $\text{Mg}^+(\text{H}_2\text{O})_2$ ($n=2$) in the Gas Phase. <i>ChemPlusChem</i> , 2013, 78, 1040-1048. | 1.3 | 22 |
| 80 | Reactivity of Hydrated Monovalent First Row Transition Metal Ions $\text{M}^+(\text{H}_2\text{O})_2$ ($n=2$), $\text{M} = \text{V}, \text{Cr}, \text{Mn}, \text{Fe}, \text{Co}, \text{Ni}, \text{Cu}, \text{Zn}$, toward Molecular Oxygen, Nitrous Oxide, and Carbon Dioxide. <i>Journal of Physical Chemistry A</i> , 2013, 117, 1011-1020. | 1.1 | 42 |
| 81 | Gas-Phase Reactions of Cationic Vanadium-Phosphorus Oxide Clusters with C_2H_4 ($n=4, 6$): A DFT-Based Analysis of Reactivity Patterns. <i>Chemistry - A European Journal</i> , 2013, 19, 3017-3028. | 1.7 | 24 |
| 82 | Reactions of Hydrated Singly Charged First-Row Transition-Metal Ions $\text{M}^+(\text{H}_2\text{O})_2$ ($\text{M}=\text{V}, \text{Cr}, \text{Mn}, \text{Fe}, \text{Co}, \text{Ni}, \text{Cu}, \text{and Zn}$) toward Nitric Oxide in the Gas Phase. <i>Chemistry - A European Journal</i> , 2013, 19, 3741-3750. | 1.7 | 21 |
| 83 | Photodissociation and photochemistry of $\text{V}^+(\text{H}_2\text{O})_4$ in the 360-680 nm region. <i>Low Temperature Physics</i> , 2012, 38, 717-722. | 0.2 | 17 |
| 84 | Electron induced reactions of unsaturated hydrocarbons in water clusters. <i>International Journal of Mass Spectrometry</i> , 2012, 330-332, 246-253. | 0.7 | 6 |
| 85 | Reactions of $\text{M}^+(\text{H}_2\text{O})_2$, $n=2$, $n=4$, $n=6$, $n=8$, $n=10$, $n=12$, $n=14$, $n=16$, $n=18$, $n=20$, $n=22$, $n=24$, $n=26$, $n=28$, $n=30$, $n=32$, $n=34$, $n=36$, $n=38$, $n=40$, $\text{M} = \text{V}, \text{Cr}, \text{Mn}, \text{Fe}, \text{Co}, \text{Ni}, \text{Cu}, \text{and Zn}$, with D_2O Reveal Water Activation in $\text{M}^+(\text{H}_2\text{O})_2$. <i>Journal of Physical Chemistry A</i> , 2012, 116, 10676-10682. | 1.1 | 31 |
| 86 | Reactions of CH_3SH and CH_3SCH_3 with Gas-Phase Hydrated Radical Anions $(\text{H}_2\text{O})_2^-$, CO_2^- , and O_2^- . <i>Journal of Physical Chemistry A</i> , 2012, 116, 3824-3835. | 1.1 | 28 |
| 87 | Single-Molecule Force-Clamp Experiments Reveal Kinetics of Mechanically Activated Silyl Ester Hydrolysis. <i>ACS Nano</i> , 2012, 6, 1314-1321. | 7.3 | 33 |
| 88 | The structure of gas-phase $[\text{Al}(\text{H}_2\text{O})_n]^+$: hydrated monovalent aluminium $\text{Al}^+(\text{H}_2\text{O})_n$ or hydride-hydroxide $\text{HAlOH}^+(\text{H}_2\text{O})_n$? <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 6776. | 1.3 | 21 |
| 89 | Mechanically activated rupture of single covalent bonds: evidence of force induced bond hydrolysis. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 5994. | 1.3 | 48 |
| 90 | Competition between Birch reduction and fluorine abstraction in reactions of hydrated electrons $(\text{H}_2\text{O})_n^-$ with the isomers of di- and trifluorobenzene. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 8924. | 1.3 | 10 |

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|-----|---|-----|-----------|
| 91 | Hydrated Magnesium Cations $Mg^{+}(H_2O)_n$, Exhibit Chemistry of the Hydrated Electron in Reactions with O_2 and CO_2 . Journal of Physical Chemistry A, 2011, 115, 10174-10180. | 1.1 | 37 |
| 92 | Ion-molecule reactions of $CoAr_6^+$ with di- and trifluorobenzenes probe absolute pressure in FT-ICR MS. International Journal of Mass Spectrometry, 2011, 300, 44-49. | 0.7 | 8 |
| 93 | Charging Effects in an Electron Bombarded Ar Matrix and the Role of Chemiluminescence-Driven Relaxation. Journal of Physical Chemistry A, 2011, 115, 7258-7266. | 1.1 | 10 |
| 94 | Darstellung, Reaktivität gegenüber Kohlenwasserstoffen und elektronische Struktur von heteronuclearen Vanadium-Phosphor-Sauerstoff-Clusterionen. Angewandte Chemie, 2011, 123, 1466-1470. | 1.6 | 42 |
| 95 | Generation, Reactivity Towards Hydrocarbons, and Electronic Structure of Heteronuclear Vanadium Phosphorous Oxygen Cluster Ions. Angewandte Chemie - International Edition, 2011, 50, 1430-1434. | 7.2 | 73 |
| 96 | Diatomic $[CuO]^+$ and Its Role in the Spin-Selective Hydrogen- and Oxygen-Atom Transfers in the Thermal Activation of Methane. Angewandte Chemie - International Edition, 2011, 50, 4966-4969. | 7.2 | 156 |
| 97 | Selective Formic Acid Synthesis from Nanoscale Electrochemistry. Angewandte Chemie - International Edition, 2010, 49, 8257-8259. | 7.2 | 25 |
| 98 | Ion-molecule reactions of $CoAr_6^+$ with nitrogen oxides N_2O , NO , and NO_2 : measuring absolute pressure by shock-freezing of the collision complex. Low Temperature Physics, 2010, 36, 411-416. | 0.2 | 7 |
| 99 | Formation of $(Xe_2H)^+$ centers in solid Xe via recombination: nonstationary luminescence and internal electron emission. Low Temperature Physics, 2010, 36, 407-410. | 0.2 | 6 |
| 100 | The $[Au_n, Si]^+$, $n=1-4$, potential energy surface: Competition between $Au-Si$ and $Au-Au$ bonding. Journal of Chemical Physics, 2010, 132, 224307. | 1.2 | 16 |
| 101 | Simple Coupling Chemistry Linking Carboxyl-Containing Organic Molecules to Silicon Oxide Surfaces under Acidic Conditions. Langmuir, 2010, 26, 15333-15338. | 1.6 | 26 |
| 102 | Gas-Phase Ion Chemistry of Small Gold Cluster Anions. Organometallics, 2010, 29, 3001-3006. | 1.1 | 19 |
| 103 | Thermochemistry from ion-molecule reactions of hydrated ions in the gas phase: a new variant of nanocalorimetry reveals product energy partitioning. Physical Chemistry Chemical Physics, 2010, 12, 3772. | 1.3 | 49 |
| 104 | Thermal N_2H Bond Activation on Anionic and Cationic Platinum Clusters: Non-Predetermined Reaction Pathways Indicate Transitions to a Bulk Surface Reactivity. Chemistry - A European Journal, 2009, 15, 8465-8474. | 1.7 | 15 |
| 105 | A novel design of a temperature-controlled FT-ICR cell for low-temperature black-body infrared radiative dissociation (BIRD) studies of hydrated ions. International Journal of Mass Spectrometry, 2009, 279, 5-9. | 0.7 | 55 |
| 106 | Gas Phase Ion Chemistry of Gold-Silicon Clusters. ChemPhysChem, 2008, 9, 1383-1386. | 1.0 | 8 |
| 107 | Activation of Methane by Oligomeric $(Al_2O_3)_x^+$ ($x=3,4,5$): The Role of Oxygen-Centered Radicals in Thermal Hydrogen-Atom Abstraction. Angewandte Chemie - International Edition, 2008, 47, 1946-1950. | 7.2 | 183 |
| 108 | Gas-phase reactivities of charged platinum dimers with ammonia: A combined experimental/theoretical study. Chemical Physics Letters, 2008, 450, 268-273. | 1.2 | 16 |

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|-----|---|-----|-----------|
| 109 | Birch reduction of chlorobenzene in gas-phase hydrated electrons. <i>International Journal of Mass Spectrometry</i> , 2008, 277, 206-209. | 0.7 | 8 |
| 110 | Dynamic Strength of the Silicon-Carbon Bond Observed over Three Decades of Force-Loading Rates. <i>Journal of the American Chemical Society</i> , 2008, 130, 3664-3668. | 6.6 | 70 |
| 111 | Electric deflection studies of rhodium clusters. <i>Journal of Chemical Physics</i> , 2007, 126, 104301. | 1.2 | 31 |
| 112 | Reactions of Large Water Cluster Anions with Hydrogen Chloride: Formation of Atomic Hydrogen and Phase Separation in the Gas Phase. <i>Journal of the American Chemical Society</i> , 2007, 129, 3238-3246. | 6.6 | 23 |
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