

# Carmine D'Agostino

## List of Publications by Year in descending order

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Version: 2024-02-01

74  
papers

3,639  
citations

136885

32  
h-index

133188

59  
g-index

78  
all docs

78  
docs citations

78  
times ranked

4133  
citing authors

| #  | ARTICLE  | IF  | CITATIONS |
|----|--|-----|-----------|
| 1  | Correlating the strength of reducing agent adsorption with Ag/Al <sub>2</sub> O <sub>3</sub> catalyst performances in selective catalytic reduction (SCR) of NO <sub>x</sub> . <i>Catalysis Today</i> , 2022, 384-386, 274-278.  | 2.2 | 13        |
| 2  | Alkaline pretreatment of walnut shells increases pore surface hydrophilicity of derived biochars. <i>Applied Surface Science</i> , 2022, 571, 151253.  | 3.1 | 17        |
| 3  | NMR relaxation time measurements of solvent effects in an organocatalysed asymmetric aldol reaction over silica SBA-15 supported proline. <i>Reaction Chemistry and Engineering</i> , 2022, 7, 269-274.  | 1.9 | 14        |
| 4  | Insights into Substituent Effects of Benzaldehyde Derivatives in a Heterogeneous Organocatalyzed Aldol Reaction. <i>ChemCatChem</i> , 2022, 14, .  | 1.8 | 6         |
| 5  | Deep eutectic solvents: alternative reaction media for organic oxidation reactions. <i>Reaction Chemistry and Engineering</i> , 2021, 6, 582-598.  | 1.9 | 57        |
| 6  | Nuclear spin relaxation as a probe of zeolite acidity: a combined NMR and TPD investigation of pyridine in HZSM-5. <i>Physical Chemistry Chemical Physics</i> , 2021, 23, 17752-17760.   | 1.3 | 19        |
| 7  | <sup>23</sup> Na NMR $T_1$ relaxation measurements as a probe for diffusion and dynamics of sodium ions in salt-glycerol mixtures. <i>Journal of Chemical Physics</i> , 2021, 154, 224501.   | 1.2 | 6         |
| 8  | Tailoring pore structure and surface chemistry of microporous Alumina-Carbon Molecular Sieve Membranes (Al-CMSMs) by altering carbonization temperature for optimal gas separation performance: An investigation using low-field NMR relaxation measurements. <i>Chemical Engineering Journal</i> , 2021, 424, 129313. | 6.6 | 21        |
| 9  | NMR Investigation into the Influence of Surface Interactions on Liquid Diffusion in a Mesoporous Catalyst Support. <i>Topics in Catalysis</i> , 2020, 63, 319-327.   | 1.3 | 10        |
| 10 | Light-driven, heterogeneous organocatalysts for C-C bond formation toward valuable perfluoroalkylated intermediates. <i>Science Advances</i> , 2020, 6, .  | 4.7 | 75        |
| 11 | Oxidative Coupling of Aldehydes with Alcohol for the Synthesis of Esters Promoted by Polystyrene-Supported N-Heterocyclic Carbene: Unraveling the Solvent Effect on the Catalyst Behavior Using NMR Relaxation. <i>Organic Letters</i> , 2020, 22, 4927-4931.  | 2.4 | 22        |
| 12 | Creation of Al-Enriched Mesoporous ZSM-5 Nanoboxes with High Catalytic Activity: Converting Tetrahedral Extra-framework Al into Framework Sites by Post Treatment. <i>Angewandte Chemie - International Edition</i> , 2020, 59, 19478-19486.   | 7.2 | 69        |
| 13 | Creation of Al-Enriched Mesoporous ZSM-5 Nanoboxes with High Catalytic Activity: Converting Tetrahedral Extra-framework Al into Framework Sites by Post Treatment. <i>Angewandte Chemie</i> , 2020, 132, 19646-19654.  | 1.6 | 12        |
| 14 | Tailoring morphology of hierarchical catalysts for tuning pore diffusion behaviour: a rational guideline exploiting bench-top pulsed-field gradient (PFG) nuclear magnetic resonance (NMR). <i>Molecular Systems Design and Engineering</i> , 2020, 5, 1193-1204.  | 1.7 | 12        |
| 15 | Self-diffusion of glycerol in $\gamma$ -alumina nanopores. The neglected role of pore saturation in the dynamics of confined polyalcohols. <i>Applied Surface Science</i> , 2020, 516, 146089.   | 3.1 | 6         |
| 16 | Solvent Effects in the Homogeneous Catalytic Reduction of Propionaldehyde with Aluminium Isopropoxide Catalyst: New Insights from PFG NMR and NMR Relaxation Studies. <i>ChemPhysChem</i> , 2020, 21, 1101-1106.   | 1.0 | 9         |
| 17 | In situ high-pressure <sup>13</sup> C/ <sup>1</sup> H NMR reaction studies of benzyl alcohol oxidation over a Pd/Al <sub>2</sub> O <sub>3</sub> catalyst. <i>Reaction Chemistry and Engineering</i> , 2020, 5, 1053-1057.  | 1.9 | 7         |
| 18 | Nanoscale Clustering of Alcoholic Solutes in Deep Eutectic Solvents Studied by Nuclear Magnetic Resonance and Dynamic Light Scattering. <i>ACS Sustainable Chemistry and Engineering</i> , 2019, 7, 15086-15092.   | 3.2 | 26        |

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|----|---|-----|-----------|
| 19 | HfN Nanoparticles: An Unexplored Catalyst for the Electrocatalytic Oxygen Evolution Reaction. <i>Angewandte Chemie</i> , 2019, 131, 15610-15616.  | 1.6 | 9         |
| 20 | HfN Nanoparticles: An Unexplored Catalyst for the Electrocatalytic Oxygen Evolution Reaction. <i>Angewandte Chemie - International Edition</i> , 2019, 58, 15464-15470.   | 7.2 | 31        |
| 21 | Globular and Fibrous Proteins Modified with Deep Eutectic Solvents: Materials for Drug Delivery. <i>Molecules</i> , 2019, 24, 3583.   | 1.7 | 18        |
| 22 | Exploiting enhanced paramagnetic NMR relaxation for monitoring catalyst preparation using <sup>1</sup> T <sub>1</sub> and <sup>19</sup> F PFG NMR: NMR correlation maps. <i>Reaction Chemistry and Engineering</i> , 2019, 4, 268-272.            | 1.9 | 6         |
| 23 | Microstructure evolution during nano-emulsification by NMR and microscopy. <i>Journal of Colloid and Interface Science</i> , 2019, 551, 138-146.  | 5.0 | 4         |
| 24 | Unravelling mass transport in hierarchically porous catalysts. <i>Journal of Materials Chemistry A</i> , 2019, 7, 11814-11825.  | 5.2 | 57        |
| 25 | Diffusion, Ion Pairing and Aggregation in 1-Ethyl-3-Methylimidazolium-Based Ionic Liquids Studied by <sup>1</sup> H and <sup>19</sup> F PFG NMR: Effect of Temperature, Anion and Glucose Dissolution. <i>ChemPhysChem</i> , 2018, 19, 1081-1088. | 1.0 | 50        |
| 26 | Effect of Al content on the strength of terminal silanol species in ZSM-5 zeolite catalysts: a quantitative DRIFTS study without the use of molar extinction coefficients. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 4250-4262.      | 1.3 | 28        |
| 27 | Inhibitory effect of oxygenated heterocyclic compounds in mesoporous catalytic materials: A pulsed-field gradient NMR diffusion study. <i>Microporous and Mesoporous Materials</i> , 2018, 269, 88-92.  | 2.2 | 7         |
| 28 | Direct Correlation between Adsorption Energetics and Nuclear Spin Relaxation in a Liquid-saturated Catalyst Material. <i>ChemPhysChem</i> , 2018, 19, 2448-2448.  | 1.0 | 1         |
| 29 | Base adsorption mechanism over zeolite catalysts at different Al contents probed by the tapered element oscillating microbalance (TEOM). <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 25357-25364.                                      | 1.3 | 2         |
| 30 | Direct Correlation between Adsorption Energetics and Nuclear Spin Relaxation in a Liquid-saturated Catalyst Material. <i>ChemPhysChem</i> , 2018, 19, 2472-2479.  | 1.0 | 40        |
| 31 | A predictive model for the diffusion of a highly non-ideal ternary system. <i>Physical Chemistry Chemical Physics</i> , 2018, 20, 18436-18446.  | 1.3 | 15        |
| 32 | Product Inhibition in Glycerol Oxidation over Au/TiO <sub>2</sub> Catalysts Quantified by NMR Relaxation. <i>ACS Catalysis</i> , 2018, 8, 7334-7339.  | 5.5 | 20        |
| 33 | Mutual diffusion in concentrated liquid solutions: A new model based on cluster theory. <i>Journal of Molecular Liquids</i> , 2017, 232, 516-521.   | 2.3 | 15        |
| 34 | Maxwell-Stefan diffusion coefficient estimation for ternary systems: an ideal ternary alcohol system. <i>Physical Chemistry Chemical Physics</i> , 2017, 19, 16071-16077.   | 1.3 | 14        |
| 35 | Increased Affinity of Small Gold Particles for Glycerol Oxidation over Au/TiO <sub>2</sub> Probed by NMR Relaxation Methods. <i>ACS Catalysis</i> , 2017, 7, 4235-4241.   | 5.5 | 43        |
| 36 | Exploring catalyst passivation with NMR relaxation. <i>Faraday Discussions</i> , 2017, 204, 439-452.  | 1.6 | 12        |

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|----|---|-----|-----------|
| 37 | Swelling-induced structural changes and microparticle uptake of gelatin gels probed by NMR and CLSM. <i>Soft Matter</i> , 2017, 13, 2952-2961.  | 1.2 | 12        |
| 38 | Liquid-liquid equilibrium for the ternary system ethanol/toluene/n-decane: a correction to the existing coexistence curve and NRTL parameters. <i>Physics and Chemistry of Liquids</i> , 2017, 55, 669-673.                                   | 0.4 | 2         |
| 39 | Effect of paramagnetic species on $T_{1\rho}$ , $T_{2\rho}$ and $T_{1\rho}/T_{2\rho}$ NMR relaxation times of liquids in porous $\text{CuSO}_4/\text{Al}_2\text{O}_3$ . <i>RSC Advances</i> , 2017, 7, 36163-36167.                           | 1.7 | 16        |
| 40 | Effect of Al content on number and location of hydroxyl acid species in zeolites: a DRIFTS quantitative protocol without the need for molar extinction coefficients. <i>RSC Advances</i> , 2017, 7, 52604-52613.                              | 1.7 | 37        |
| 41 | Hole theory as a prediction tool for Brownian diffusive motion in binary mixtures of liquids. <i>RSC Advances</i> , 2017, 7, 51864-51869.   | 1.7 | 10        |
| 42 | Do group 1 metal salts form deep eutectic solvents?. <i>Physical Chemistry Chemical Physics</i> , 2016, 18, 25528-25537.  | 1.3 | 43        |
| 43 | Adsorption of pyridine from aqueous solutions by polymeric adsorbents MN 200 and MN 500. Part 2: Kinetics and diffusion analysis. <i>Chemical Engineering Journal</i> , 2016, 306, 1223-1233.   | 6.6 | 85        |
| 44 | Adsorption of pyridine from aqueous solutions by polymeric adsorbents MN 200 and MN 500. Part 1: Adsorption performance and PFG-NMR studies. <i>Chemical Engineering Journal</i> , 2016, 306, 67-76.  | 6.6 | 52        |
| 45 | Solvent inhibition in the liquid-phase catalytic oxidation of 1,4-butanediol: understanding the catalyst behaviour from NMR relaxation time measurements. <i>Catalysis Science and Technology</i> , 2016, 6, 7896-7901.                       | 2.1 | 46        |
| 46 | Prediction of mutual diffusion coefficients in binary liquid systems with one self-associating component from viscosity data and intra-diffusion coefficients at infinite dilution. <i>Chemical Engineering Science</i> , 2016, 147, 118-127. | 1.9 | 27        |
| 47 | Assessing the effect of reducing agents on the selective catalytic reduction of $\text{NO}_x$ over $\text{Ag}/\text{Al}_2\text{O}_3$ catalysts. <i>Catalysis Science and Technology</i> , 2016, 6, 1661-1666.                                 | 2.1 | 32        |
| 48 | Molecular and ionic diffusion in aqueous "deep eutectic solvent mixtures: probing inter-molecular interactions using PFG NMR. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 15297-15304.   | 1.3 | 204       |
| 49 | Probing hydrogen-bonding in binary liquid mixtures with terahertz time-domain spectroscopy: a comparison of Debye and absorption analysis. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 5999-6008.                                  | 1.3 | 36        |
| 50 | A local composition model for the prediction of mutual diffusion coefficients in binary liquid mixtures from tracer diffusion coefficients. <i>Chemical Engineering Science</i> , 2015, 132, 250-258.   | 1.9 | 33        |
| 51 | Structure and dynamics of aqueous 2-propanol: a THz-TDS, NMR and neutron diffraction study. <i>Physical Chemistry Chemical Physics</i> , 2015, 17, 30481-30491.   | 1.3 | 29        |
| 52 | A kinetic analysis methodology to elucidate the roles of metal, support and solvent for the hydrogenation of 4-phenyl-2-butanone over Pt/TiO <sub>2</sub> . <i>Journal of Catalysis</i> , 2015, 330, 362-373.                                 | 3.1 | 12        |
| 53 | Effect of solvent on the hydrogenation of 4-phenyl-2-butanone over Pt based catalysts. <i>Journal of Catalysis</i> , 2015, 330, 344-353.  | 3.1 | 49        |
| 54 | Assessing the use of NMR chemical shifts for prediction of VLE in non-ideal binary liquid mixtures. <i>Chemical Engineering Science</i> , 2014, 119, 331-333.   | 1.9 | 1         |

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|----|--|-----|-----------|
| 55 | The Effect of Grafting Zirconia and Ceria onto Alumina as a Support for Silicotungstic Acid for the Catalytic Dehydration of Glycerol to Acrolein. <i>Chemistry - A European Journal</i> , 2014, 20, 1743-1752.                    | 1.7 | 36        |
| 56 | Assessing the surface modifications following the mechanochemical preparation of a Ag/Al <sub>2</sub> O <sub>3</sub> selective catalytic reduction catalyst. <i>Catalysis Science and Technology</i> , 2014, 4, 531-539.           | 2.1 | 46        |
| 57 | Deactivation studies of a carbon supported AuPt nanoparticulate catalyst in the liquid-phase aerobic oxidation of 1,2-propanediol. <i>Catalysis Science and Technology</i> , 2014, 4, 1313-1322.                                   | 2.1 | 34        |
| 58 | The effect of coke deposition on the activity and selectivity of the HZSM-5 zeolite during ethylbenzene alkylation reaction in the presence of ethanol. <i>Catalysis Science and Technology</i> , 2014, 4, 1017.                   | 2.1 | 31        |
| 59 | Interpretation of NMR Relaxation as a Tool for Characterising the Adsorption Strength of Liquids inside Porous Materials. <i>Chemistry - A European Journal</i> , 2014, 20, 13009-13015.   | 1.7 | 98        |
| 60 | <i>In situ</i> study of reaction kinetics using compressed sensing NMR. <i>Chemical Communications</i> , 2014, 50, 14137-14140.  | 2.2 | 35        |
| 61 | Mesoscopic Structuring and Dynamics of Alcohol/Water Solutions Probed by Terahertz Time-Domain Spectroscopy and Pulsed Field Gradient Nuclear Magnetic Resonance. <i>Journal of Physical Chemistry B</i> , 2014, 118, 10156-10166. | 1.2 | 109       |
| 62 | Prediction of the mutual diffusivity in acetone-chloroform liquid mixtures from the tracer diffusion coefficients. <i>Chemical Engineering Science</i> , 2013, 95, 43-47.  | 1.9 | 30        |
| 63 | Solvent Effect and Reactivity Trend in the Aerobic Oxidation of 1,3-Propanediols over Gold Supported on Titania: NMR Diffusion and Relaxation Studies. <i>Chemistry - A European Journal</i> , 2013, 19, 11725-11732.              | 1.7 | 46        |
| 64 | Hydrogen Bonding Network Disruption in Mesoporous Catalyst Supports Probed by PFG-NMR Diffusometry and NMR Relaxometry. <i>Journal of Physical Chemistry C</i> , 2012, 116, 8975-8982.   | 1.5 | 65        |
| 65 | Understanding the Solvent Effect on the Catalytic Oxidation of 1,4-Butanediol in Methanol over Au/TiO <sub>2</sub> Catalyst: NMR Diffusion and Relaxation Studies. <i>Chemistry - A European Journal</i> , 2012, 18, 14426-14433.  | 1.7 | 50        |
| 66 | Prediction of mutual diffusion coefficients in non-ideal mixtures from pulsed field gradient NMR data: Triethylamine-water near its consolute point. <i>Chemical Engineering Science</i> , 2012, 74, 105-113.                      | 1.9 | 32        |
| 67 | Solvent effects in the hydrogenation of 2-butanone. <i>Journal of Catalysis</i> , 2012, 289, 30-41.  | 3.1 | 140       |
| 68 | Neutron diffraction, NMR and molecular dynamics study of glucose dissolved in the ionic liquid 1-ethyl-3-methylimidazolium acetate. <i>Chemical Science</i> , 2011, 2, 1594.   | 3.7 | 121       |
| 69 | Pulsed-Field Gradient NMR Spectroscopic Studies of Alcohols in Supported Gold Catalysts. <i>Journal of Physical Chemistry C</i> , 2011, 115, 1073-1079.  | 1.5 | 35        |
| 70 | Glycerol eutectics as sustainable solvent systems. <i>Green Chemistry</i> , 2011, 13, 82-90.   | 4.6 | 666       |
| 71 | Molecular motion and ion diffusion in choline chloride based deep eutectic solvents studied by <sup>1</sup> H pulsed field gradient NMR spectroscopy. <i>Physical Chemistry Chemical Physics</i> , 2011, 13, 21383.                | 1.3 | 397       |
| 72 | Prediction of binary diffusion coefficients in non-ideal mixtures from NMR data: Hexane-nitrobenzene near its consolute point. <i>Chemical Engineering Science</i> , 2011, 66, 3898-3906.  | 1.9 | 48        |

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|----|--|-----|-----------|
| 73 | A Neutron Diffraction and Molecular Dynamics Investigation of Acetate-Based Ionic Liquids as Solvents for Glucose. ECS Transactions, 2010, 33, 611-620.              | 0.3 | 0         |
| 74 | Structure and Dynamics of 1-Ethyl-3-methylimidazolium Acetate via Molecular Dynamics and Neutron Diffraction. Journal of Physical Chemistry B, 2010, 114, 7760-7768. | 1.2 | 117       |