

Fall

Chemistry 501

2020

Seminar Chair:
Dr. Johnathan
Brantley

**All seminars will
be via zoom**



SEPTEMBER 3 • link: tba • Host: Johnathan Brantley <i>Frieder Jäkle · Distinguished Professor, Chair · Rutgers University</i> “Putting Boron into Polymers: From Catalysis to Supramolecular Material”
SEPTEMBER 10 • link: tba • Host: Johnathan Brantley <i>Konstantinos Vogiatzis · Assistant Professor · University of Tennessee</i> “Coupling Quantum Chemistry with Machine Learning”
SEPTEMBER 17 • link: tba • Host: John Larese <u>Lind Lecture</u> <i>Thomas Mallouk · Professor Emeritus · Penn State University</i> TBA
SEPTEMBER 24 • link: tba • Host: Michael Kilbey <i>Barry Thompson · Professor of Chemistry · University of Southern California</i> “Reimagining Semiconducting Polymers for Alternative Energy Applications”
OCTOBER 1 • link: tba • Host: Ben Xue <i>Danna Freedman · Professor · Northwestern University</i> “Chemistry for the Second Quantum Revolution (and some bismuth)”
OCTOBER 8 • link: tba • Host: Ampofo Darko <i>Michael Doyle · Professor, Rita and John Feik Distinguished University Chair in Medicinal Chemistry · University of Texas, San Antonio</i> “Enantioselective Cycloaddition Reactions and Their Applications”
OCTOBER 22 • link: tba • Host: Constance Bailey <i>William Wuest · GRA Distinguished Investigator & Assoc Professor · Emory University</i> “Slaying superbugs one natural product at a time”
OCTOBER 29 • link: tba • Host: Bin Zhao <i>Yue Zhao · Professor of Chemistry · Université de Sherbrooke - Canada</i> “Polymer Motions Driven by Stimuli: towards Soft Robots”
NOVEMBER 5 • link: tba • Host: Thanh Do <i>Etienne Garand · Associate Professor · University of Wisconsin-Madison</i> “Probing molecular interactions in reactive complexes with cryogenic ion chemistry and spectroscopy”
November 12 • link: tba • Host: Brian Long <i>Geoff Coates · Tisch University Professor · Cornell University</i> “New Polymers from Old Monomers: Advances Enabled through Catalyst Design and Discovery”
NOVEMBER 19 • link: tba • Host: Konstantinos Vogiatzis <i>Francesco Paesani · Professor of Chemistry & Biochem · Univ of California-San Diego</i> “Data-Driven Many-Body Models for Predictive Molecular Simulations from the Gas to the Condensed Phase”