## Microsymposium

## Designing Proton Conducting Metal Organic Frameworks

## G. Shimizu<sup>1</sup>, <u>B. Gelfand<sup>1</sup></u>

## <sup>1</sup>University of Calgary, Department of Chemistry, Calgary, Canada

Metal organic frameworks (MOFs) or porous coordination polymers (PCPs) represent a tunable molecular scaffolding that can be adjusted for a breadth of applications. This presentation will concern our efforts towards tailoring MOFs towards making new proton conductors ultimately for fuel cells. A major hurdle in these technologies is an electrolyte capable of conducting protons above 100°C. Higher operating temperatures will enhance electrode kinetics and decrease electrode poisoning among several critical operational benefits. In contrast to the macromolecular approaches typically employed towards these electrolytes, we have used a MOF strategy to generate crystalline networks with acidic pores. These MOFs present options to address higher temperature conduction,1 conduction over 10-2 Scm-1,2 and water stability.3 The emphasis in the talk will concern routes to designing these systems and subsequent challenges in their characterization.

[1] J. A. Hurd et al. Nature Chem. 2009, 1, 705., [2] S. Kim et al., J. Am. Chem. Soc. 2013, 135, 963., [3] J. M. Taylor et al., J. Am. Chem. Soc. 2013, 135, 1193.

Keywords: MOF, Proton Conductor