

## Design of new ionic liquids with increased O<sub>2</sub> solubility for Li-O<sub>2</sub> battery

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In the past decade, Li-O<sub>2</sub> battery received a lot of attention since it gives the promise of large energy density which would potentially increase the range of Electric Vehicles (EV) above 400kms, reducing the customer range anxiety and therefore supporting a better penetration of EV in the market. Despite numerous studies, Li-O<sub>2</sub> battery remain an immature technology facing several challenges [1]. An important one is the electrolyte selection which is until now not resolved since the requirements for Li-O<sub>2</sub> battery electrolyte are even stronger than for Li-ion battery. Especially high O<sub>2</sub> solubility and O<sub>2</sub> diffusivity and low vapor pressure are important extra parameters to consider.

Ionic liquids (ILs), solvent consisting entirely of ions, present unique properties such as low volatility, high (electro) chemical stability, intrinsic electrical conductivity, thermal stability and are highly tuneable. Previous studies reveal that N,N-diethyl-N-(2-methoxyethyl)-N-methylammonium bis(trifluoromethylsulfonyl)imide (DEME-TFSI) was stable against O<sub>2</sub> radicals which is not the case of the common organic solvents such as carbonates or glymes used in Li-ion battery [2-3]. On the downside, one of the disadvantage of DEME-TFSI lies in its low O<sub>2</sub> solubility combined with a high viscosity, penalizing in the same time the O<sub>2</sub> diffusivity.

In order to increase the O<sub>2</sub> solubility, our approach was therefore to design new ionic liquids functionalized with fluorine, making them an intrinsic part of the electrolyte [4]. We will report the synthesis and characterization of those new ionic liquids, before to discuss their O<sub>2</sub> properties determined by electrochemical measurements combining the Cottrell equation on a macro-electrode and the steady-state current on an ultra-microelectrode.

### References:

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