## Tailored MnO<sub>2</sub> Nanoparticles as Electrocatalysts for Metal-Air Batteries

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In next-generation electronics, electrified transportation and energy storage, metal-air batteries represent one class of promising power sources thanks to their remarkably high theoretical energy and light weight [1]. The main feature of metal-air batteries is the combination of a metal anode with high energy density and an air electrode with open structure that facilitates the drawing of cathode active materials (*i.e.* oxygen) from air [1]. In these types of devices, Gas Diffusion Electrodes (GDEs) are widely used as cathodes [2]. However, one of the main drawbacks related to the cathodic reaction (ORR) is the overpotential loss (about 0.3-0.4 V) under operative conditions. Thus, lots of efforts were spent to inhibit the voltage loss requiring an effective ORR catalyst [1,3]. One of the most promising materials, in terms of both performances and costs, seems to be MnOx. According to the recent literature, MnO<sub>2</sub> would ensure capacities comparable to those of platinum, letting higher capacity retention to be reached in non-aqueous electrolytes to prevent Li decomposition [1].

In the present work, the electrochemical performances of either bare or Fe/Co-doped MnO<sub>2</sub> nanoelectrocatalysts are evaluated by Linear Sweep Voltammetries (LSVs). The crystal structure and the surface properties are examined by means of XRPD, BET-BJH, TEM, SEM/EDX and XPS analyses. Correlations between their physico-chemical features and the final electrocatalytic performances are drawn. Experimental results reveal that the as-synthesized powders have excellent electrochemical properties in organic electrolytes (0.15 M LiNO<sub>3</sub> in propylene carbonate, PC) showing a shift of the onset potential of about 150 mV with 2% Co-doped MnO<sub>2</sub>, thus resulting very promising candidates to be used in lithium-air batteries [4].

## References

- [1] F. Cheng et al., Soc. Rev. 2012, 41, 2172–2192.
- [2] S. Orsini et al., J. of Electroan. Chem. 2018, 808, 439–445.
- [3] E. Pargoletti et al., J. of Power Sources 2016, 325, 116-128.
- [4] E. Pargoletti et al., *Nanomaterials* **2020**, 10, 1735.