

SUPPORTING INFORMATION

**The critical importance of the ionomer on the electrochemical activity of platinum and
platinum-free catalysts for anion-exchange membrane fuel cells**

Pietro G. Santori,^a Abhishek N. Mondal,^b Dario R. Dekel,^{b,c *} and Frédéric Jaouen^{a *}

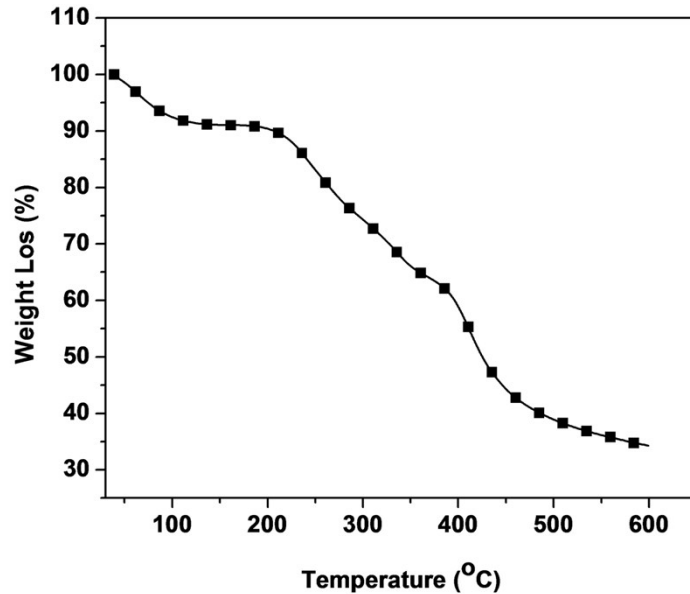


Figure S1. TGA curve of PPO-TMA in the temperature range of 30 to 600 °C. The measurement was performed under N₂ flow. The weight loss between 30 and 100°C is assigned to residual water. The PPO-TMA starts decomposing at *ca* 210°C, much earlier than pure PPO (*ca* 400°C). This is in line with TGA results reported for other AEIs based on a PPO polymer backbone, and in particular for another PPO functionalized with TMA cationic group.^{S1-S2}

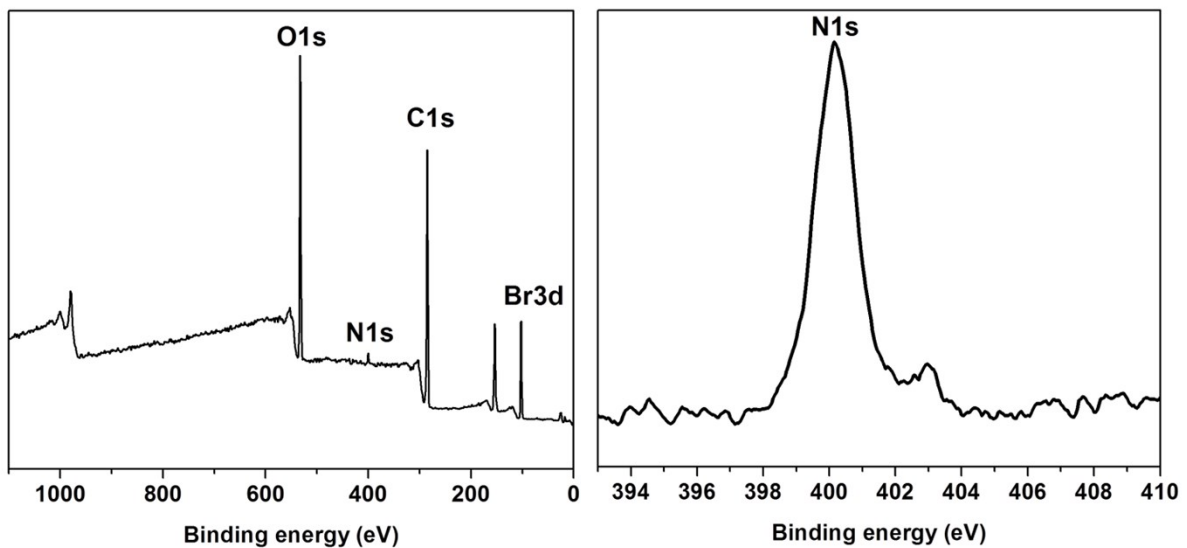


Figure S2. XPS survey spectrum and high resolution spectrum of N1s region of PPO-TMA in bromide form.

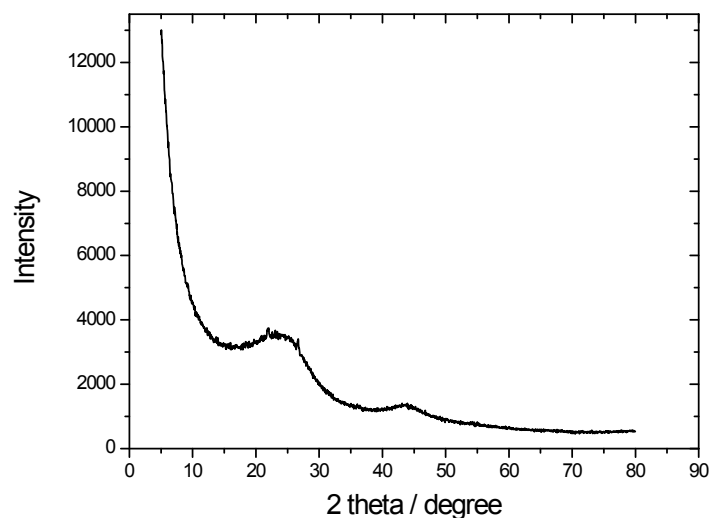


Figure S3. XRD pattern of Fe_{0.5}-950. The pattern shows only two broad peaks assigned to the amorphous carbon matrix with nanosized graphite crystallites.

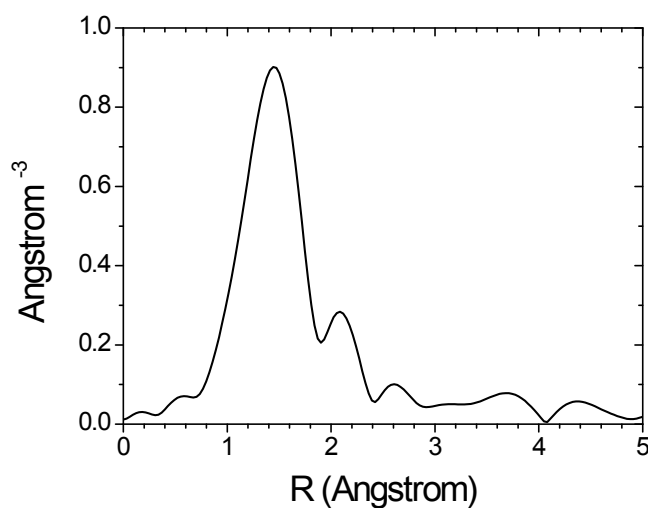


Figure S4. *Ex situ* Fe K-edge EXAFS spectrum of Fe_{0.5}-950. The spectrum was recorded in air at room temperature. The distance is not corrected for phase-shift. The signal at 1.5 Å is assigned to Fe-N or Fe-O interaction from FeN_x sites and O₂ adsorbed on Fe from such sites, at 2.2 Å is assigned to Fe-Fe interaction from a minor fraction of Fe having formed particles during the pyrolysis in NH₃, and at 2.6 Å is assigned to Fe-C interactions between Fe from FeN_x sites and C atoms from the second coordination sphere. This result is similar to our previous publication (Ref. 45, Fe-N-C sample prepared similarly to the present Fe_{0.5}-950 catalyst), where EXAFS assignments have been discussed also on the basis of a rigorous fitting procedure.

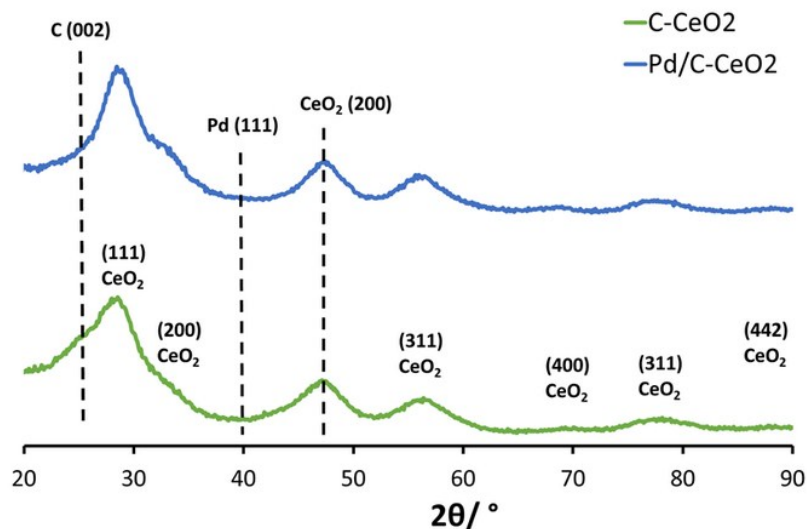


Figure S5. XRD patterns for CeO₂/C and Pd-CeO₂/C. Both patterns show only the reflections from the cubic structure of CeO₂ with the large FWHM consistent with nanosized CeO₂ particles. No signal for metallic Pd is visible in the pattern of Pd-CeO₂/C, indicating non-crystalline and/or high dispersion of palladium in this catalyst.

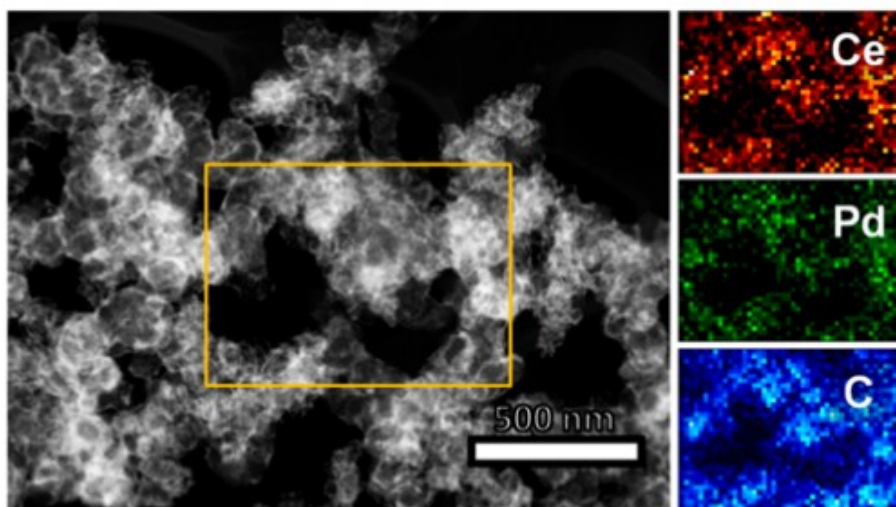


Figure S6. STEM images of Pd-CeO₂/C. HAADF image and STEM-EDS elemental mapping of the selected area (highlighted by orange rectangle on the left) showing Ce (red), Pd (green) and carbon (blue) elemental distribution. Reprinted with permission from *ACS Appl. Energy Mater.*, 2019, **2**, 4999–5008. Copyright (2019) American Chemical Society.

Supporting references

- S1. H.-S. Dang and P. Jannasch, *J. Mater. Chem. A*, 2016, **4**, 11924-11938.
- S2. H.-S. Dang and P. Jannasch, *Macromolecules*, 2015, **48**, 5742-5751.