

Jueves, 18 de enero 18:35 – 18:48

Study of mixed molybdates of lanthanum and praseodymium, $\text{La}_{5.4-x}\text{Pr}_x\text{Mo}_{1-y}\text{Nb}_y\text{O}_{12-\delta}$, for H_2 separation membranes.

Abraham Sánchez-Caballero*, José M. Porrás-Vázquez, Lucía dos Santos- Gómez, Javier Zamudio-García, Antonia Infantes-Molina, David Marrero- López, Enrique R. Losilla

a. Dpto. de Química Inorgánica, Cristalografía y Mineralogía: Universidad de Málaga, 29071-Málaga, España,

b. Dpto. de Física Aplicada I, Laboratorio de Materiales y Superficies (Unidad Asociada al C.S.I.C.): Universidad de Málaga, 29071-Málaga, España.

abraham11sc@uma.es

This research explores the impact of praseodymium and niobium doping on the $\text{La}_{5.4-x}\text{Pr}_x\text{Mo}_{1-y}\text{Nb}_y\text{O}_{12-\delta}$ series ($x = 1.35, 2.7, 4.05, 5.4$; $y = 0, 0.1$) as materials with mixed proton-ionic conductivity suitable for hydrogen separation membranes with two main objectives: i) enhancing the electronic conductivity through praseodymium doping, and ii) increasing oxide ion vacancies to improve the ionic conductivity via niobium doping [1].

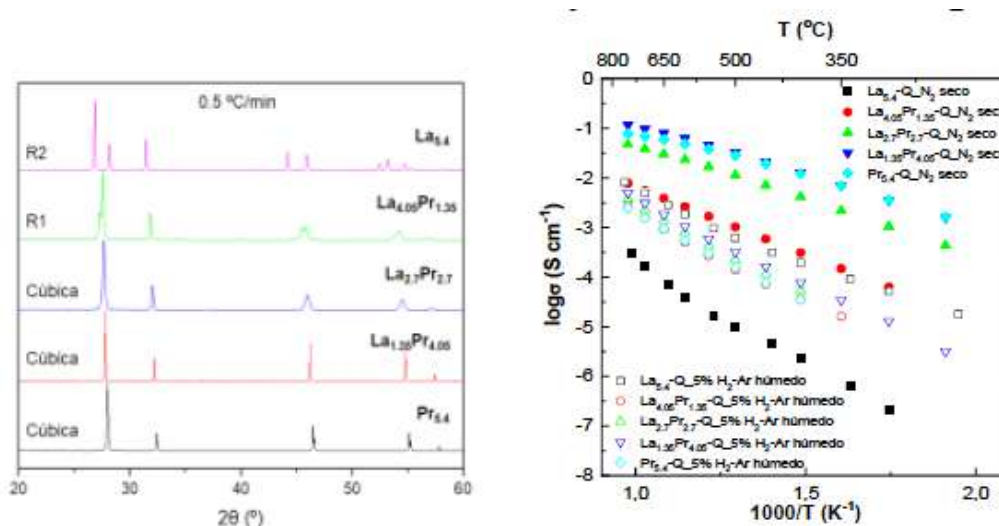


Figure 1. X-ray diffractograms of $\text{La}_{5.4-x}\text{Pr}_x\text{Mo}_{1-y}\text{Nb}_y\text{O}_{12-\delta}$ are presented on the left, accompanied by the Arrhenius diagram illustrating the conductivity on the right.

Material synthesis involves the freeze-drying method, and characterization includes X-ray diffraction (Rietveld method), X-ray photoelectron spectroscopy, scanning and transmission electron microscopy and complex impedance spectroscopy. The symmetry of the materials (cubic or rhombohedral) is influenced by synthetic conditions and composition, compositions with higher praseodymium content exhibit a consistent cubic fluorite-type structure. Optimal conductivity is observed in a nitrogen atmosphere for compositions with high praseodymium content due to the presence of the Pr⁴⁺/Pr³⁺ pair. In a reducing atmosphere (wet 5% H₂-Ar), conductivity significantly decreases due to the reduction of Pr⁴⁺ to Pr³⁺.

References

1. López-Vergara *et al.*, *Chemistry of Materials*, 2017, 29, 6966-6975.
<https://doi.org/10.1021/acs.chemmater.7b02481>

Jueves, 18 de enero 18:48 – 19:00

Transmission electron microscopy study of porous metal-organic frameworks (MOFs)

J. M. Vila Fungueiriño^{a*}, M. Ceballos^{b†}, S. Funes-Hernando^b, G. Zampini^c, M. Cedrún-Morales^b, M. Rodríguez-Pérez^c, M. F. Navarro Poupard^c, E. Polo^d, P. del Pino^b, B. Pelaz^e

a, Centro Singular de Investigación en Química Biolóxica e Materiais Moleculares (CiQUS), Departamento de Química Física, Universidade de Santiago de Compostela (USC), 15782 Santiago de Compostela, Galicia, Spain.

b, CiQUS, Departamento de Física de Partículas, USC, 15782 Santiago de Compostela, Galicia, Spain.

c, CiQUS, USC, 15782 Santiago de Compostela, Galicia, Spain.

d, CiQUS, Departamento de Bioquímica, USC, 15782 Santiago de Compostela, Galicia, Spain.

e, CiQUS, Departamento de Química Inorgánica, USC, 15782 Santiago de Compostela, Galicia, Spain.

* Presenting or corresponding author: josem.vila@usc.es

Metal-Organic Frameworks (MOFs) are an emerging family of crystalline and porous micro/nanomaterials composed of a extended network comprising