

STRUCTURAL AND MORPHOLOGICAL CHARACTERIZATION OF THE PEROVSKITE $\text{LaFe}_{0.2}\text{Cr}_{0.8-x}\text{Co}_x\text{O}_3$ ($x = 0.0, 0.2, 0.4, 0.6, 0.8$) FOR CO SELECTIVE OXIDATION

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ABSTRACT

In order to obtain perovskites with specific structural, and morphological properties for potential applications as catalytic materials, five compositions of the system $\text{LaFe}_{0.2}\text{Cr}_{0.8-x}\text{Co}_x\text{O}_3$ ($x = 0.0, 0.2, 0.4, 0.6, 0.8$) were synthesized by means of a combustion technique and evaluated in methane reforming of CH_4 .

The structural and morphological characterization of the calcined oxides, carried out by X-ray diffraction (XRD) (Figure 1) and electron microscopy (SEM-TEM) (Figures 2 and 3), confirmed the obtainment of a single perovskite type structure in all materials. It was noticed that the incorporation of cobalt in the structure is associated with an evolution from a rhombohedral to an orthorhombic structure with nanometric particle sizes and relevant surface properties [1, 2, 3]. From the micrographs, it is clear that the solids are made up of irregular multiparticulate aggregates, heterogeneously distributed with sizes ranging between 0.5 and 200 μm . The study showed average methane conversion values of $64.78 \pm 1.08\%$, corresponding to obtaining $63.43 \pm$

1.12% for H₂ and $61.73 \pm 0.99\%$ for CO. Similarly, the expressions of catalytic reactivity, in terms of the methane conversion, the selectivities (S) towards H₂, CO and the synthesis gas ratio obtained were estimated, showing that the material has a high degree of selectivity towards the production of hydrogen and carbon monoxide, always maintaining an H₂/CO ratio > 3.0, a beneficial situation in any reforming process [4]. The conversion results as a function of temperature, shown in Figure 4, confirm that reforming is a thermally activated process, whose maximum conversion value is reached at 700 °C.

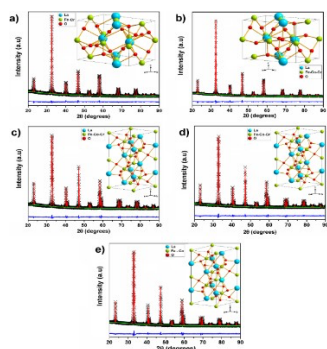


Figure 1. Rietveld refinement of the **a) LFM1, b) LFM2, c) LFM3, d) LFM4, and e) LFM5** samples.

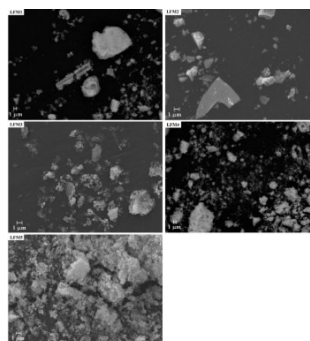


Figure 2. Scanning electron micrographs for synthesized samples

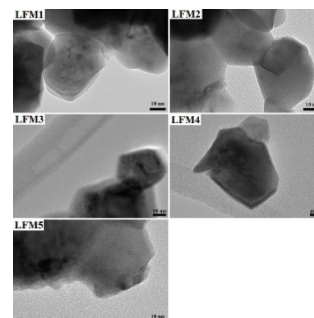


Figure 3. Transmission electron micrographs for synthesized samples.

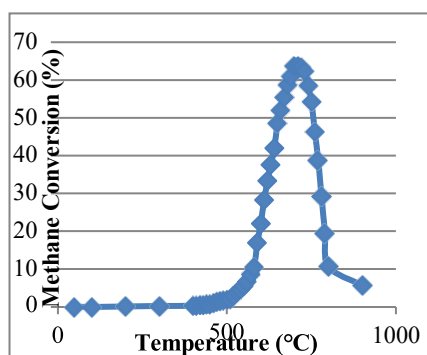


Figure 4. Methane conversion curve as a function of reaction temperature obtained by steam reforming using the LFM5 material.

CONCLUSIONS

Six perovskite-type materials based on the system $\text{LaFe}_{0.2}\text{Cr}_{0.8-x}\text{Co}_x\text{O}_3$ ($x = 0.0, 0.2, 0.4, 0.6, 0.8$) were synthesized employing a combustion method. Structural analysis using XRD, and electron microscopy showed that the oxides were single perovskite type phase. It was also noted the formation of particles with a high homogeneity and distribution, and outstanding interconnected porosity, generated in the transformation of the citrate species and the decomposition of volatile compounds. The study showed average methane conversion values of $64.78 \pm 1.08\%$, corresponding to obtaining $63.43 \pm 1.12\%$ for H_2 and $61.73 \pm 0.99\%$ for CO , which are within the normally accepted range for studies of this type, showing that under the proposed analysis conditions, the material has a high degree of selectivity towards the production of hydrogen and carbon monoxide, always maintaining an H_2/CO ratio > 3.0 .

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