



P-type transition metal carbodiimides MNCN ($M = \text{Ni, Mn}$)

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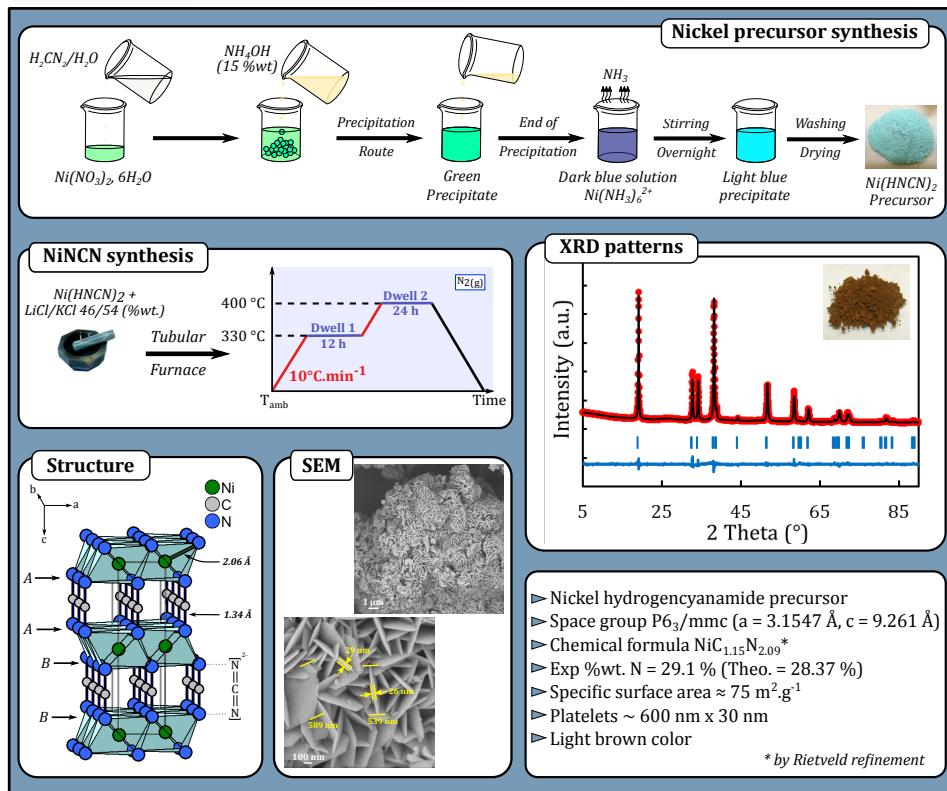
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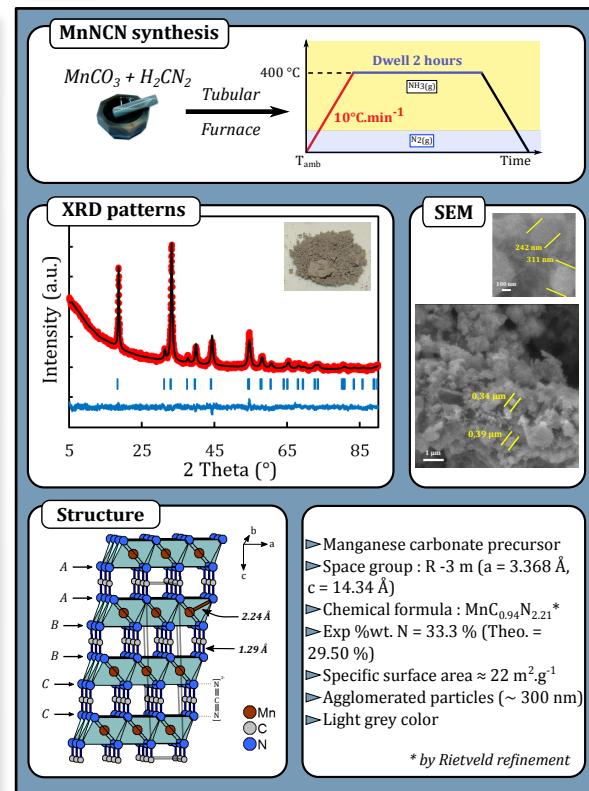
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Over the last decade, p-type semiconductors (SC) have known a renewed interest. Indeed these materials may have potential applications for light-emitting diodes, transistors, solar cells, etc. Since the stabilization of the first p-type semi-conductor CuAlO₂ with a delafossite structure type (A⁺B³⁺O₂) by H. Kawazoe in 1997 [1], many p-type semiconductors with the same structure have been reported such as CuGaO₂, CuCrO₂, CuScO₂... Recently, X. Liu *et. al* have reported the synthesis and the structure determination of a new family of materials: the transition metal carbodiimides MNCN (with M = Mn → Cu) [2,3]. These materials adopt a structure type related to the delafossite, formed by a divalent transition metal and NCN²⁻ anion. We report here the synthesis and the microstructural characterization of two transition metal carbodiimides MnNCN and NiNCN. Impedance measurements demonstrate the p-type conductivity of these materials.

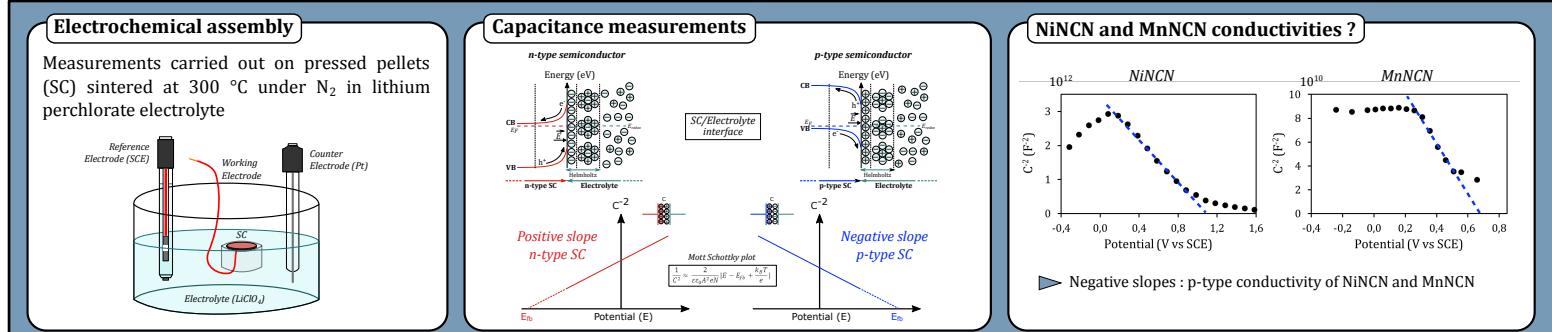
Nickel carbodiimide (NiNCN)



Manganese carbodiimide (MnNCN)



p-type characterization



References

- [1] Kawazoe, H.; Yasukawa, M.; Hyodo, H.; Kurita, M.; Yanagi, H.; Hosono, H. *Nature* **1997**, 389, 939. [2] Liu, X.; Krott, M.; Muller, P.; Hu, C.; Lueken, H.; Dronskowski, R. *Inorg. Chem.* **2005**, 44, 3001. [3] Krott, M.; Liu, X.; Fokwa, B. P.; Speldrich, M.; Lueken, H.; Dronskowski, R. *Inorg. Chem.* **2007**, 46, 2204.