Influence of the Co/Mo Ratio on the SWNT Synthesis from Carbon Monoxide

Toshiaki Nishii¹, Suguru Noda², Hisashi Sugime³, Naoto Masuyama⁴, Shigeo Maruyama⁵

¹ J-Power / The Univ. of Tokyo

² The Univ. of Tokyo

³ The Univ.of Tokyo

⁴ J-Power

⁵ The Univ. of Toyko

Contact e-mail: toshiaki_nishii@jpower.co.jp

By using carbon monoxide as carbon source, we have synthesized random, vertical and parallel aligned SW-NTs on quartz substrates dip-coated with Co and Mo (COCCVD). In order to study the influence of the Co/Mo ratio in details, combinatorial method [1] with wide nominal thickness profiles of metal catalysts was employed. By the uniquely masked sputter-deposition technique, we prepared a 2-dimensional catalyst library with Mo (0.2 - 4 nm) and Co (0.2 - 8 nm) thickness profiles on a SiO₂/Si wafer [2]. After the CVD using hydrogen / carbon monoxide (500 / 500 sccm) (800 degC, 1 atom), SWNT formation region was confirmed by HRSEM observation and micro Raman analysis. The optimum Co/Mo ratio for the SWNT formation was larger in the order of COCCVD, ACCVD [2] and CoMoCAT [3]. The difference of the optimum Co/Mo ratio between COCCVD and ACCVD is suspected to originate in the relation between the activity of Co and the energy requirement for graphite precipitation in a thermodynamic equilibrium process. Furthermore, the effect of support materials, heating or reducing processes, catalyst preparations will be discussed.

References:

[1] S. Noda, Y. Tsuji, Y. Murakami, S. Maruyama, Appl. Phys. Lett., 86 (2005) 173106.

[2] S. Noda, H. Sugime, T. Osawa, Y. Tsuji, S. Chiashi, Y. Murakami, and S. Maruyama, Carbon, in press.

[3] J.E.Herrera, L. Balzano, A. Borgna, W.E. Alvarez, D. E. Resasco, J. Catalysis, 204 (2001) 129.

[4] M. Hu, Y. Murakami, M. Ogura, S. Maruyama , T. Okubo, J. Catalysis, 225 (2004) 230.