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Single-wall carbon nanotubes (SWNTs) have opened up a new possibility for novel components in miniaturized electronic devices. SWNTs are one-dimensional structure that exhibit metallic or semiconducting behavior. Semiconducting nanotubes have been shown to function as field-effect active channels¹, nanorectifiers². Due to the doping effect of adsorbed oxygen molecules and the charge transfer with defined electrodes, pristine SWNTs show p-type semiconducting behavior. For complementary logic circuitry, various chemical modification achieved by doping semiconducting nanotubes have been demonstrated^{3,4}. However, the stability of alkali-metal doped nanotubes^{3,4} is sensitive to the ambient environment. Therefore, ambient stable chemical doping by metallofullerene will be a good starting point for the band engineering of carbon-based nanoelectronics.

The mono-metallofullerene encapsulating a single lanthanide element, Ln@C₈₂ (Ln = Ce, Nd, Gd, Dy, ...), has been confirmed by electron spin resonance and UV-Vis-NIR adsorption spectra to show charge transfer from the encaged atom to the C₈₂ cage,⁵ resulting in the charge state of Ln³⁺@C₈₂³⁻. The insertion of metallofullerenes into the nanotube is therefore expected to lead to further charge transfer from the C₈₂ cage to the tube and the electrical properties of the carbon nanotubes would be substantially modified.

The insertion of Dy@C₈₂ into the inner hollow space of the SWNTs host was carried out by gas phase diffusion of Dy@C₈₂.⁶ The individual metallofullerene peapods were deposited on a degenerately doped silicon substrate with SiO₂ insulating layer on top and predefined electrodes. Fig. 1 shows the transport measurement in one single tube. At room temperature, the doped nanotube shows p-type behavior as seen from the response to a back gate. As the temperature decreases, the conductance becomes n-type. It reveals that the Dy@C₈₂ molecules function as electron donors and transfer charge to the carbon nanotube host. The amount of charge transferred varies with temperature. At T < 215 K metallic behavior is observed, indicating the degenerate state by doping. Below about 75 K, single-electron charging phenomena dominate the transport and show irregular Coulomb blockade oscillation, implying that the insertion of Dy@C₈₂ splits the tube into a series of several quantum dots.

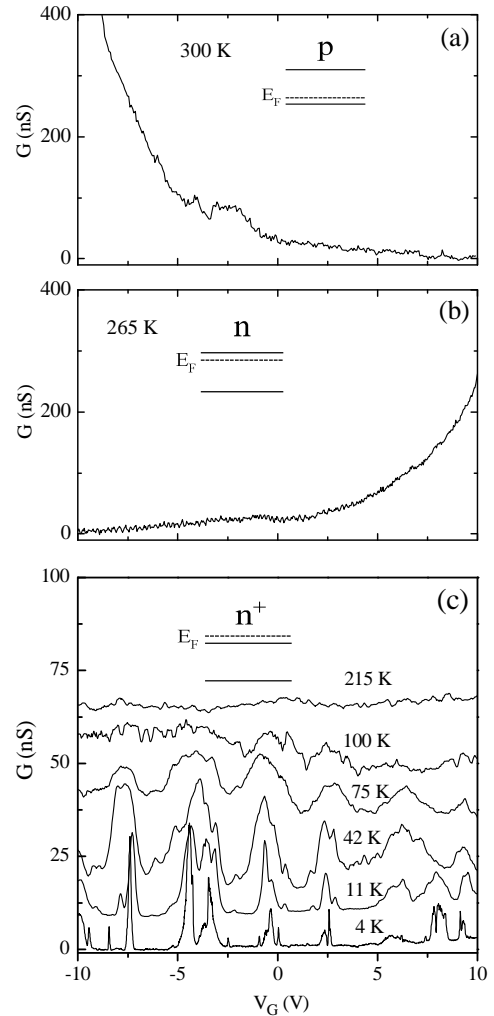


Fig. 1 (a)-(c) Gate voltage dependence of conductance measured at $V_{DS} = 4$ mV at various temperatures. The curves in (c) are displaced for clarity. The insets are schematic plots of the band diagram.

¹ S. J. Tans, A. R. M. Verschueren, and C. Dekker, *Nature* **393**, 49 (1998).

² R. D. Antonov and A. T. Johnson, *Phys. Rev. Lett.* **83**, 3274 (1999).

³ L. Grigorian, G. U. Sumanasekera, A. L. Loper, S. Fang, J. L. Allen, and P. C. Eklund, *Phys. Rev. B* **58**, R4195 (1998).

⁴ J. Kong, C. Zhou, E. Yenilmez, and H. Dai, *Appl. Phys. Lett.* **77**, 3977 (2000).

⁵ For a review, see H. Shinohara, *Rep. Prog. Phys.* **63**, 843 (2000).

⁶ K. Suenaga, M. Tence, C. Mory, C. Colliex, H. Kato, T. Okazaki, K. Hirahara, S. Bandow, and S. Iijima, *Science* **290**, 2280 (2000).