

Chemical p- and n-doping for MoS₂ transistor and its application

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Abstract

The discovery of isolation of the monolayer graphene using mechanically exfoliation has triggered the intense studies on the other two-dimensional (2D) materials, e.g., semiconducting molybdenum disulphide (MoS₂) and insulating hexagonal boron nitride (h-BN) [1]-[3]. MoS₂, one of the transition metal dichalcogenides (TMDCs), has a great potential for replacing the Si-based electronics due to its comparable band gap as large as 1.8 eV for monolayer and 1.2 eV for multilayers. For more versatile transistor applications, however, the doping process is inevitably necessary since various practical devices such as tunneling, transistors, logic circuits, and memory devices are composed of p-n junctions. Although there have been some reports about surface charge transfer doping effects of MoS₂ using potassium [4], MoO₃ [5], and molecules [6] recently, it is still quite ambiguous to apply for practical junction devices since they just demonstrated slight changes of electrical properties or improvement of contact resistance.

In this study, we demonstrate the chemical p- and n-doping for MoS₂ using AuCl₃ and benzyl viologen (BV), respectively. **Figure 1** shows the electrical properties (I_D - V_G) of MoS₂ transistors after (a) BV and (b) AuCl₃ doping. Before these doping, the MoS₂ transistor shows n-type semiconducting property as previously reported (see **Figure 1(a)**) [2]. However, it becomes to degenerate n-type and p-type semiconductor after BV and AuCl₃ chemical doping, respectively. Using potassium, which is a strong electron donor, there has been similar behavior compared with BV doping [4]. While the clear p-type semiconducting property using AuCl₃ has not been reported from our literature survey so far. Furthermore, it seems feasible to apply for the p-n diodes toward various optoelectronics. For achieving this, we stacked h-BN on a partial area of MoS₂ and dope the other area using AuCl₃ as shown in **Figure 2(a)**. When we measured the photo-response of this prototype diode, it showed the unique properties of the conventional diode such as open circuit voltage and short circuit current as shown in **Figure 2(b)**. Consequently, we demonstrated the proper chemical doping method for MoS₂ toward various optoelectronics.

References

- [1] Novoselov, K.S., et al. *Science*, **306**, (2004) 666-669.
- [2] Radisavljevic, B., et al. *Nat. Nanotech.*, **6**, (2011) 147-150.
- [3] Dean, C.R., et al. *Nat. Nanotech.*, **5**, (2010) 722-726.
- [4] Fang, H., et al., *Nano Lett.*, **13**, (2013) 1991-1995.
- [5] Lin, J., et al. *Appl. Phys. Lett.*, **103**, (2013) 063109.
- [6] Du, Y., et al. *IEEE Elec. Dev. Lett.*, **34**, (2013) 1328-1330.

Figures

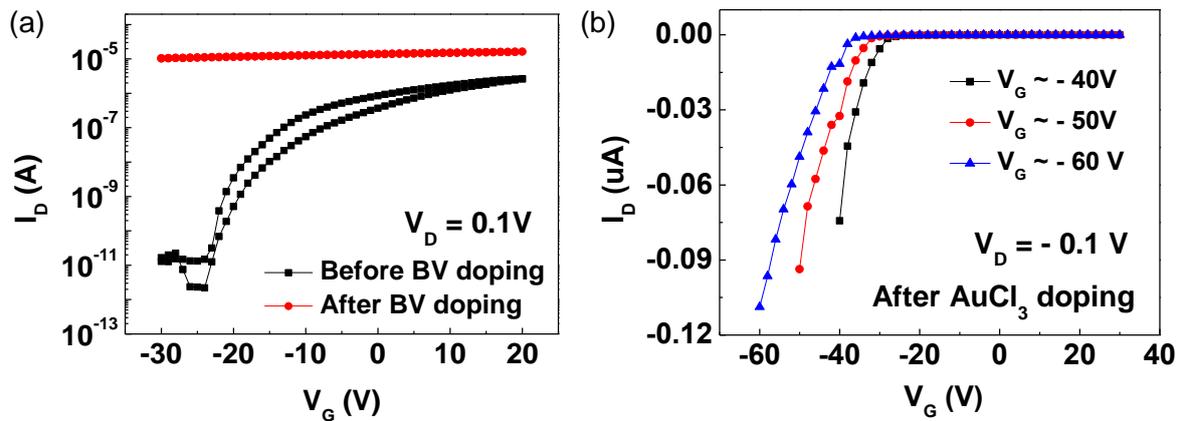


Figure 1: The electrical properties (I_D - V_G) of MoS₂ transistors (a) before and after BV doping, and (b) after AuCl₃ doping.

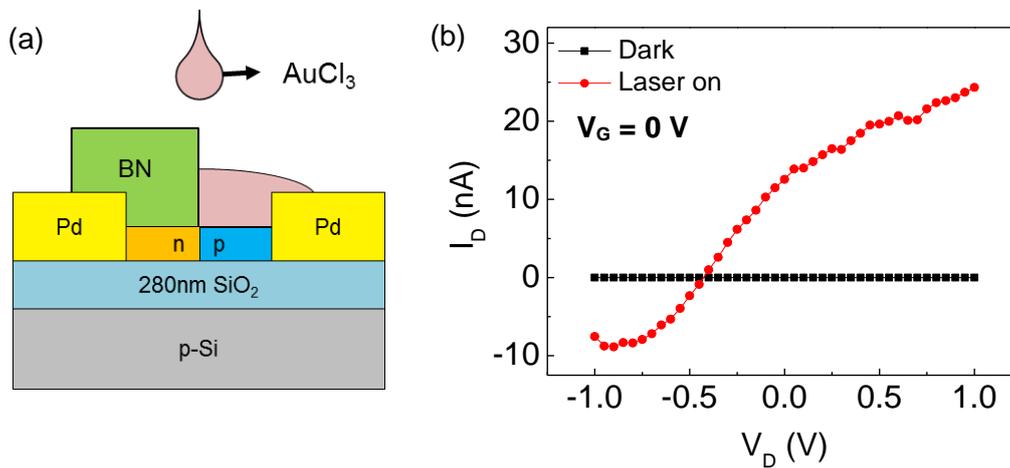


Figure 2: (a) Cross-sectional diagram of MoS₂ p-n diode. (b) The photo-response of a fabricated diode.