Realization of N-doping in Graphene via 4-Aminothiophenol Functionalization

Ashwini S. Gajarushi¹

C. Subramaniam¹, V. Ramgopal Rao^{1,2}

¹IITB, Powai, Mumbai, India ²IITD,Hauz Khas, New Delhi, Delhi, India

Contact as.gajarushi@gmail.com

Abstract

Graphene, possess exceptional electrical and optical properties due to its unique sp² hybridized hexagonal structure and linear energy-momentum relationship near Dirac point. Very high room temperature mobility and highest specific surface area makes graphene very promising candidate for electronics and sensing applications. However, graphene gets easily p-doped when exposed to ambient conditions, restricting its applicability in practice. Herein, explored have chemical we functionalization of graphene using 4-Aminothiophenol (4-ATP) for realization of ndoping in graphene. It has been observed that non-covalent functionalization via 4-ATP dopes the GFETs n-type and shifts the Dirac point of GFETs in negative direction by -40 V to -50 V gate voltage. Further it enhances the electron mobility and symmetry of current-voltage characteristics. The X-ray photoelectron Spectroscopy (XPS) of 4-ATP-doped graphene (4ATP Gr) shows occurrence of N1s peak confirming n-type doping of graphene. Moreover existence of S2s and S2p peaks elucidates thiol functionalization of graphene opening new route towards sensing application. Stability of such n-type doping (for >40 days), ensures its practical applicability towards development graphene of based electronics and sensing technology.

References

- [1] A. K. Geim et.al., Nat. Mater., 6 (2007) 183–191
- [2] W. Xu et. al., Small, 10 (2014), 1999-2005
- [3] V. Georgakilas et. al., Chem Rev, 112 (2012), 6156-6214

Figures

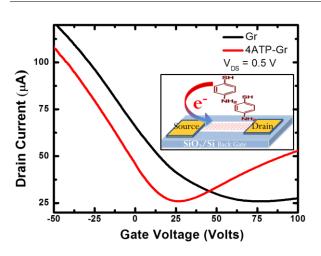


Figure 1: Gate voltage vs. Drain current characteristics of GFET (black) and 4ATP-doped GFET (red); Inset-Charge transfer interaction between 4-ATP and graphene