

Metal Contacts to MoS₂: a Two-Dimensional Semiconductor

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With increasing demands for electrostatic control as scaling continues in today's transistors, low dimensional structures continue to gain attention as a pathway for future device scaling because they offer excellent electrostatic control while remaining compatible with straightforward lithography techniques. In particular, MoS₂ has attracted interest for transistor applications because its large band gap allows for field effect devices with low off-current, unlike graphene [1]. One key bottleneck, however, is the realization of ohmic contacts on MoS₂ to improve FET device on-state performance. With this in mind, we evaluate Ni and Pd contacts on MoS₂ as potential alternatives to the already realized Au-MoS₂ and Ti-MoS₂ contacts [1]. Back-gated transfer length method (TLM) structures with Au, Ni, and Pd contact metals were fabricated on exfoliated MoS₂ flakes, with 300nm SiO₂ on degenerately doped Si as the substrate. The data indicate that Ni, like Au, makes an ohmic contact to the n-doped MoS₂ while the Pd metal contact shows Schottky behavior.

Figure 1 shows the alignment of the metal work-functions to the MoS₂ bands for the high work function metals on which we focus in this work. Figure 2 shows representative I_d - V_{ds} characteristics at $V_{bg}=50V$ for the three contact metals on MoS₂. The MoS₂ flake thicknesses are 5nm for Ni and 6nm for Au, measured by AFM, and ~13nm for Pd, estimated by optical contrast. One finds that the Ni and Au contacts are ohmic while the Pd contact shows Schottky behavior. Figure 3 shows the contact resistance (R_c) as a function of back-gate voltage (V_{bg}), extracted using the TLM. R_c for Ni and Au are found to be comparable, with a minimum R_c of about 4.5Ω.mm at $V_{bg}=50V$. The gate-dependent R_c reflects the gate modulation of the potential barrier at the metal-MoS₂ interface. Considering charge-transfer at the metal-MoS₂ interface alone for these high work function metals, the different contact nature of Pd from Ni and Au is somewhat unexpected. Indeed, recent *ab initio* calculations suggest that the modification of the electronic states at the interface by the metal is the key to understanding the contact to MoS₂, going beyond the simple charge-transfer considerations of a metal-semiconductor junction [2]. To investigate the Ni contact further, R_c was measured at low temperature via TLM, as shown in Figure 4, for a ~6nm thick flake, as estimated by optical contrast. The R_c varies weakly with temperature, increasing by less than a factor of two from 238K to 78K for any V_{bg} . Enabled by the low temperature Ni ohmic contact, the sheet resistance and mobility of MoS₂ could be extracted as shown in Figure 5. The mobility, extracted from fits to the sheet resistance as a function of V_{bg} , is determined to be 48 cm²/Vs at 289K, increasing to 237 cm²/Vs at 78K due to decreased phonon scattering. The mobility of MoS₂ could be significantly improved to several hundred cm²/Vs if its surface is passivated by atomic-layer-deposited dielectrics. [1, 3] The temperature dependence of the mobility is power law, $\mu \sim T^\gamma$, where $\gamma=1.36$ for $T=130K$ and above.

In conclusion, studies of Ni, Au, and Pd metal contacts on MoS₂ reveal that Ni, like Au, is ohmic while Pd shows Schottky contact behavior. Contact resistance for Ni and Au are comparable, with a minimum R_c for both metals of about 4.5Ω.mm at $V_{bg}=50V$.

References

- [1] B. Radisavljevic et al, *Nature Nanotechnology*, **6**, 147 (2011).
- [2] I. Popov et al, *Physical Review Letters*, **108** (2012).
- [3] H. Liu and P.D. Ye, *IEEE Electron Device Letters* (in press).

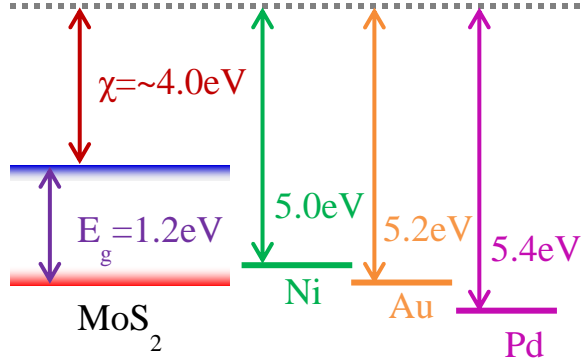


Figure 1: Band diagram showing the work function alignment of the contact metals to MoS₂

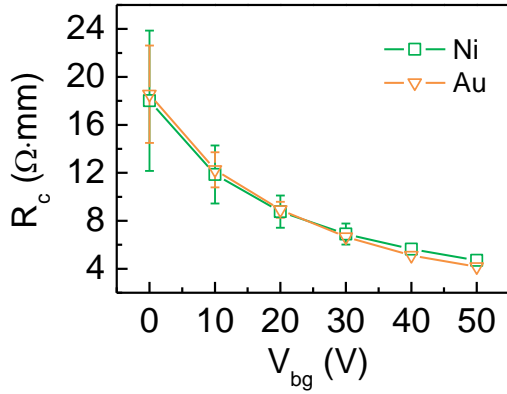


Figure 3: R_c extracted from TLM as a function of gate voltage for Ni and Au contact metals at $V_{ds}=50\text{mV}$. Error bars are determined from the liner fit used to extract R_c . Error bars are omitted for those smaller than the size of the symbol.

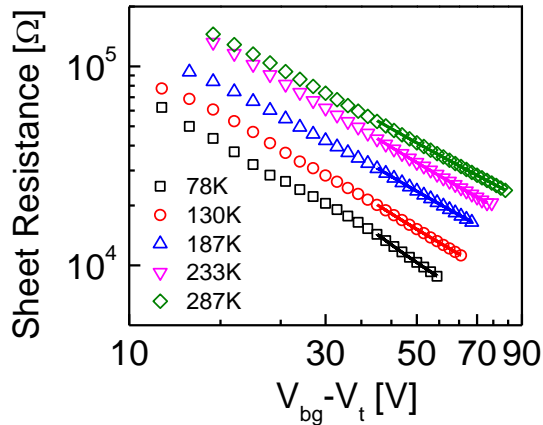


Figure 5: Symbols: Experimental sheet resistance of MoS₂ extracted at $V_{ds}=50\text{mV}$ as a function of $V_{bg}-V_t$ for different temperatures for an Ni TLM device. V_t is taken as the V_{bg} where I_d is 10nA. Solid lines: $R_{sheet}=1/C_{ox}(V_{bg}-V_t)\mu$ fits to the data used to extract mobility.

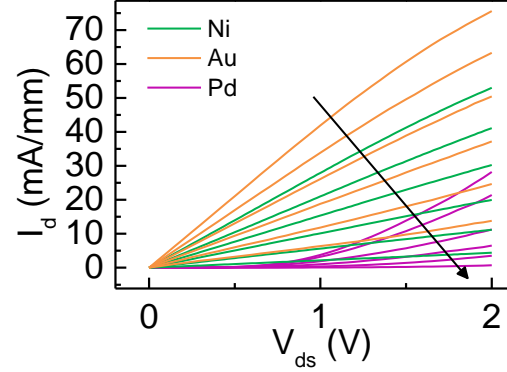


Figure 2: I_d - V_{ds} output characteristics for the three contact metals at $V_{bg}=0\text{V}$ to 50V with 10V step. Arrow indicates decreasing V_{bg} . The channel length for all three devices is 500nm .

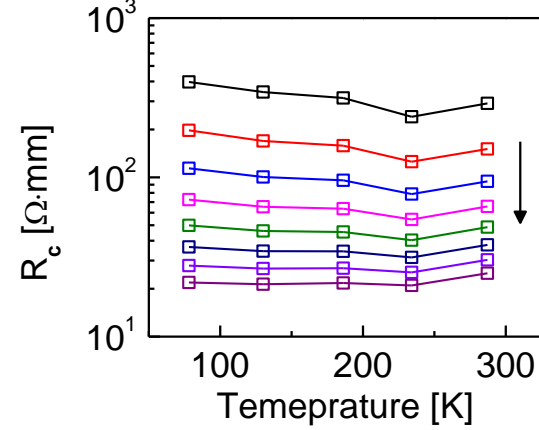


Figure 4: R_c extracted at $V_{ds}=50\text{mV}$ as a function of temperature for $V_{bg}-V_t$ from 14V to 56V with 6V step. The sample is different from the one measured in Figure 3. V_t is taken as the V_{bg} where I_d is 10nA . The arrow indicates increasing gate voltage.

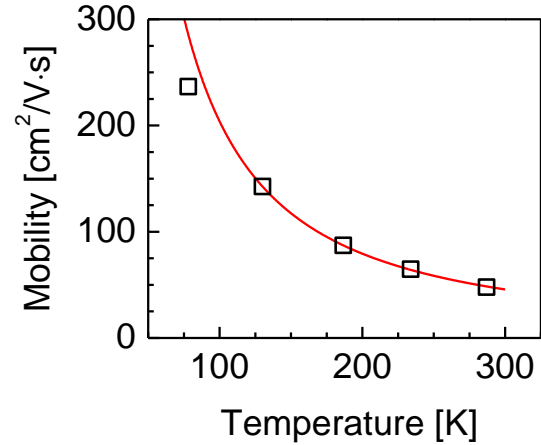


Figure 6: Mobility as a function of temperature extracted from the sheet resistance with power law fit $\mu \sim T^\gamma$, where $\gamma=1.36$, $T=130\text{K}$ and above.