

Advanced physical characterization of Si-passivated III-V semiconductors for low-power logic applications: Defect chemistry, band bending and surface photo-voltage

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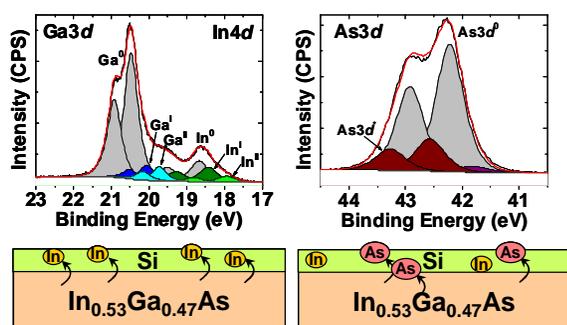
Replacing Si as a channel material in future metal-oxide-semiconductor field effect transistors (MOSFET) is today the leading option for advanced CMOS technology nodes. For n-FET, In_{0.53}Ga_{0.47}As (InGaAs) offers the best compromise between electron mobility and band-gap.

Over the last years, important progress has been made in the development of InGaAs based MOSFETs for future low-power logic application. Devices with promising characteristics have been reported. Integration path for self-aligned, laterally scaled devices has been shown [1]. Fabrication of similar devices on silicon substrate has been also demonstrated, for example using the transfer of high-quality, thin III-V heterostructures onto Si by direct wafer bonding [2]. Understanding and controlling the density of interface states (D_{it}) between the InGaAs channel and the gate dielectric remains however a key issue to obtain competitive device characteristics with respect to Si based devices.

An appropriate gate dielectric is characterized by three different metrics, namely i) a low capacitance-equivalent thickness (CET < 1.3 nm) with low enough leakage current, ii) $D_{it} < 10^{12} \text{ cm}^{-2}$ and iii) a thermal stability compatible with a gate first or gate last integration flow.

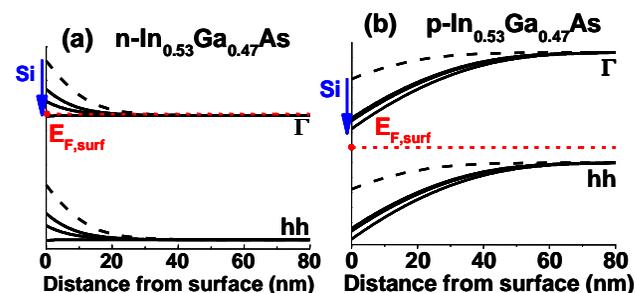
Using thin, passivating layers of amorphous Si onto InGaAs, sub-1.5 nm CET MOS capacitors and FETs with excellent mobilities have been demonstrated [3,4], in particular using a gate first integration flow. Better performance would require however to further reduce D_{it} . A detailed understanding of critical instabilities intimate to the physics and chemistry of the α -Si/InGaAs interface is therefore required. By combining laboratory and synchrotron X-ray photoelectron spectroscopy (XPS) data, we describe 1) the interface formation from the initial sub-monolayer reaction, 2) the evolution of its chemical composition, 2) Si-induced band bending and specific defect states in the InGaAs band gap, 3) surface photo-voltage effects on clean and Si-reacted interfaces.

The figure below summarizes two key features observed in the surface chemistry of InGaAs channels

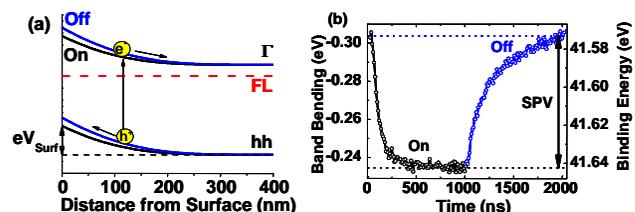


exposed to Si deposition. First the evolution of Ga 3d and In 4d core levels upon Si deposition is reported in the right figure. Each core level is featured by a doublet: Ga 3d_{5/2} and Ga 3d_{3/2}, In 4d_{5/2} and In 4d_{3/2}. Ga 3d⁰, In 4d⁰ are bulk doublets, while Ga 3d^I, Ga 3d^{II}, In 4d^I and In 4d^{II} are Si-induced components. The observation is consistent with an inherent instability between α -Si and InGaAs, leading to slight out-diffusion of Indium atoms. This phenomenon takes place starting from Si thicknesses as low as 0.3 nm. In addition, the evolution of the As 3d doublet upon 380°C anneal in UHV (left figure) reveal that this additional thermal treatment triggers a small out diffusion of arsenic. (As 3d⁰ is bulk doublet, As 3d^I is Si-induced component).

Simultaneously, the evolution of band bending can be extracted from the XPS data as illustrated in the figures below. The evolution upon Si deposition is schematized for both n- and p-type samples, respectively. With respect to the clean samples, Si deposition flattens the bands in n-type whereas an increased bending is induced in the opposite direction, creating surface inversion in p-type In_{0.53}Ga_{0.47}As. This observation seem to indicate that the deposition of Si onto the In_{0.53}Ga_{0.47}As surface does not completely “un-pin” the Fermi level, in the sense that interface states are not eliminated. In the latter case, flat band condition should be observed for all samples.



Finally, time resolved surface photo-voltage (SPV) experiments are performed to gain additional insight. In this pump-probe experiment, In_{0.53}Ga_{0.47}As samples are excited using a pulsed laser ($\lambda = 405 \text{ nm}$). Laser-induced band bending variation are followed by recording core level spectra. The mechanism depicted in the figure below (left) has been widely reported for Si. Electron-hole pairs photogenerated in the space charge region are quickly separated by the strong electric field: for upward (downward) band bending, holes (electrons) are swept towards the surface. This decreases the negative (positive) charge density trapped in the surface states and reduces the original band bending (right).



From these different observations, a consistent description of the interface defect chemistry between α -Si and InGaAs can be proposed, paving the way to solutions to further improve the quality for device operation.

- [1] L. Czornomaz et al., Proc. 2012 70th DRC (2012)
- [2] L. Czornomaz et al., IEDM Tech. Dig. (2012)
- [3] M. El Kazzi et al., APL **100**, 063505 (2012)
- [4] M. El Kazzi et al., APL **99**, 052102 (2011)