

Optical Measurements and Theory Test AlGa_N/Ga_N High Electron Mobility Transistor (HEMT) Models

by *Steven R. Kurtz*

Motivation—Ga_N-based electronics offer radical miniaturization of microwave power amplifier and radar circuitry. Ga_N's large bandgap, high breakdown field, high electron velocity, and excellent thermal properties have led to high electron mobility transistors (HEMTs) with up to ten times the power density of GaAs and other traditional semiconductors at frequencies up to 20 GHz.

Further contributing to the outstanding performance of Ga_N-based amplifiers is the highly conducting, 2-dimensional electron gas (2DEG) used for the HEMT channel. Intrinsic polarization and piezoelectric properties of Ga_N materials can produce a 2DEG at an AlGa_N/Ga_N interface with a sheet carrier concentration of $10^{13}/\text{cm}^2$, well in excess of that achievable in any other III-V material system. The physics and material science of the AlGa_N/Ga_N 2DEG are critical to the performance and future development of Ga_N-based electronics.

Accomplishment—To further reduce dislocation densities, we are evaluating alternative substrates to sapphire. Focusing on the properties of AlGa_N/Ga_N devices grown on SiC substrates, we found that overall material quality and transport properties of the 2DEG were much improved over structures grown on sapphire. Hall mobility versus electron density measurements revealed that 2DEG mobility was limited by scattering in the AlGa_N barrier, not dislocations, in devices grown on SiC.

Electronic properties of AlGa_N/Ga_N heterostructures and HEMTs on SiC were determined using a contacted electroreflectance technique. This optical probe augments conven-

tional electrical characterization of Ga_N-based field-effect transistors. By studying variations in the electroreflectance with applied electric field, spectral features associated with the AlGa_N barrier, the 2DEG at the interface, and bulk Ga_N are clearly identified. The 2DEG produced a broad, first-derivative-like electroreflectance feature. Changing bias voltage, the 2DEG electroreflectance narrowed and converged with the Ga_N band-edge. A first-principle, Golden Rule calculation of the dielectric function was developed which described the variation of 2DEG electroreflectance with voltage (see Fig. 1). The AlGa_N barrier displayed Franz-Keldysh oscillations (FKO), and the period of the FKO varied with bias voltage. Airy function lineshape fits provided accurate determinations of AlGa_N barrier composition and polarization electric field. Comparing measured AlGa_N electric fields with values predicted by a standard model of the AlGa_N/Ga_N heterostructure conduction band, we found < 10% discrepancy between the measured polarization field and that predicted by a standard model for devices grown on SiC (see Fig. 2). However, AlGa_N electric field measurements for devices grown on sapphire produced anomalous results, indicating the presence of trapped space charge.

Significance—An optical probe of AlGa_N/Ga_N heterostructures and transistors augments conventional electrical characterization tools. The electroreflectance technique and supporting calculations provide the first optical measurements of AlGa_N barrier electric field and composition and 2DEG Fermi energy and electron density. This new information allows us to test and refine basic models of AlGa_N/Ga_N heterostructures and transistors.

Sponsors for various phases of this work include: DOE Office of Basic Energy Sciences, Laboratory Directed Research & Development, and Cooperative Research & Development Agreement

Contact: Steven R. Kurtz, Semiconductor Physics, Dept. 1123

Phone: (505) 844-9637, Fax: (505) 844-3211, E-mail: srkurtz@sandia.gov

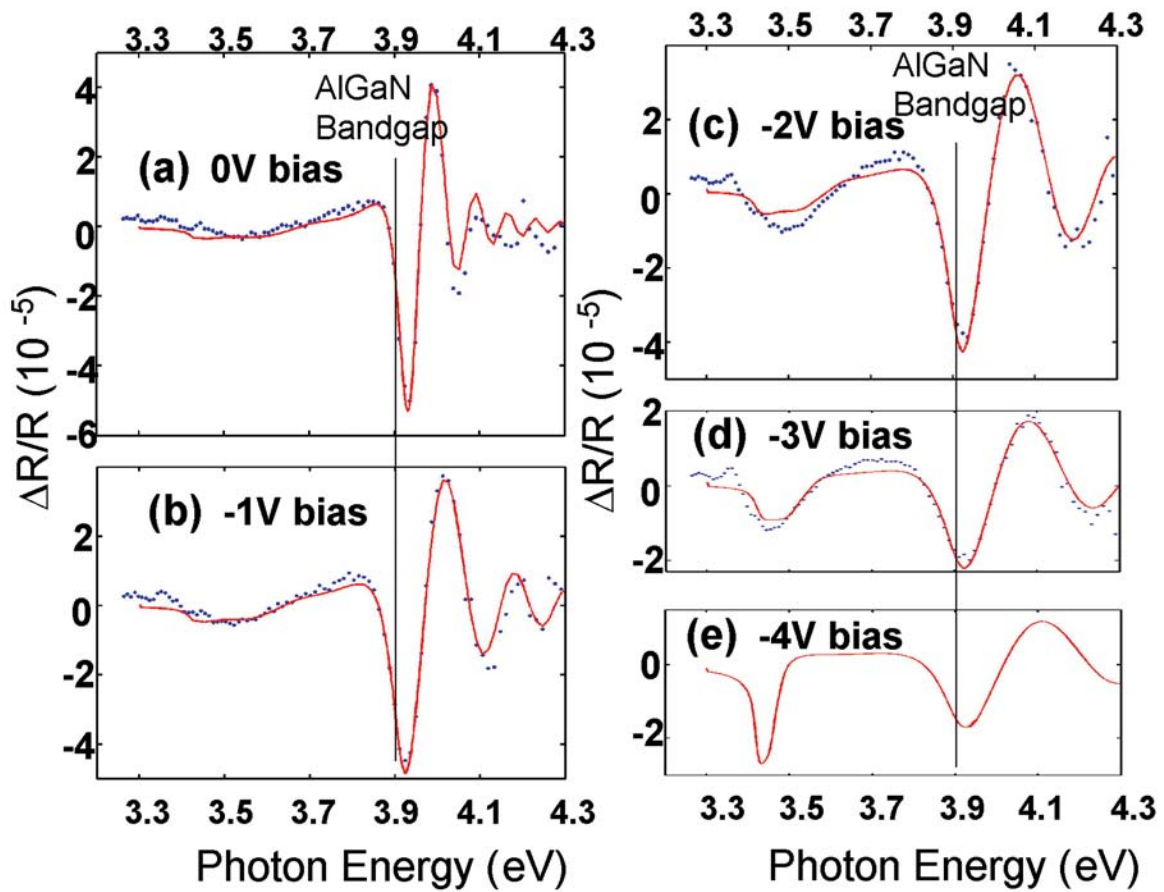


Figure 1. Electroreflectance spectra (300 K) for the $\text{Al}_{0.19}\text{Ga}_{0.81}\text{N}$ (320 Å)/GaN heterostructure at 0V (a), -1V (b), -2V (c), and -3V (d) gate bias. (same scale) Solid lines are the AlGaIn FKO lineshape fit plus the contribution from the 2DEG dielectric model. Based on the models, the -4V spectrum (e) was simulated.

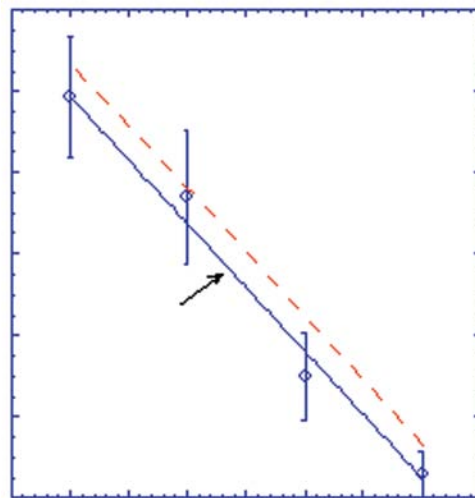


Figure 2. Electric field versus gate bias obtained from the electroreflectance spectra in 1(a)-(d). The electric field predicted by the "standard model" for an $\text{Al}_{0.19}\text{Ga}_{0.81}\text{N}$ (320 Å)/GaN heterostructure is indicated by the dotted line.