## Effects of H<sub>2</sub>O Pretreatment on the Capacitance–Voltage Characteristics of Atomic-Layer-Deposited Al<sub>2</sub>O<sub>3</sub> on Ga-Face GaN Metal–Oxide–Semiconductor Capacitors

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Atomic layer deposition (ALD) of  $Al_2O_3$  on Ga-face GaN is studied with respect to the effects of growth saturation, precursor injection sequence, and  $H_2O$ pretreatment. A metal-oxide-semiconductor capacitor (MOSCAP) structure is fabricated to measure the capacitance-voltage (*C*-*V*) characteristics. The origin of *C*-*V* hysteresis is explained by a model considering the different trapping behaviors of interface states and oxide border traps. The interface state density ( $D_{it}$ ) is extracted as a function of band bending using an ultraviolet (UV)-assisted method. It is found that  $H_2O$  pretreatment followed by saturated ALD growth produces the best interface quality, with a reduced  $D_{it}$ compared with growth without  $H_2O$  pretreatment.

**Key words:** GaN, Al<sub>2</sub>O<sub>3</sub>, atomic layer deposition, interface states, traps, hysteresis, MOSCAP

## INTRODUCTION

 $Al_2O_3$  has emerged as a suitable gate dielectric for III-nitride-based electronic devices.<sup>1-10</sup> It has a large bandgap and relatively high dielectric constant, and exhibits high thermal and structural stabilities.  $Al_2O_3$  can also be used as a passivation material to improve device output performance.<sup>1-3,7,9,11</sup> The passivation effects are often seen to be better than with the widely used SiN.<sup>1,11</sup> For both applications it is essential to have a low interface state density ( $D_{it}$ ) between  $Al_2O_3$  and III-nitride semiconductor, and a low oxide trap density within  $Al_2O_3$  itself.

Due to the great chemical versatility and the exceptional control over film uniformity and conformity, atomic layer deposition (ALD) has become the predominant method for depositing  $Al_2O_3^{12-17}$  and various other dielectrics. In typical thermal ALD of  $Al_2O_3$ , trimethylaluminum (TMA) and  $H_2O$  precursors are introduced alternately. Each half-cycle reaction is driven by the respective surface chemistry,<sup>12</sup> limiting growth to proceed in a layer-by-layer

fashion. Saturated growth is usually preferred in order to get the maximum growth rate and incorporation efficiencies, as well as better run-to-run consistencies. To achieve a low- $D_{it}$  interface at the initial stage of growth, the precursor injection sequence needs to be designed carefully by considering the surface termination and chemical properties of the underlying semiconductor.<sup>15–17</sup> For example, a TMA pretreatment step is found to reduce  $D_{it}$  in ALD of Al<sub>2</sub>O<sub>3</sub> on As-terminated InGaAs.<sup>15</sup> The As termination favors TMA adsorption and passivation,<sup>16</sup> and prefers the formation of Al–As bonds for the first layer growth.<sup>17</sup> On the other hand, for ALD of Al<sub>2</sub>O<sub>3</sub> on Ga-face GaN, it might be advantageous to use H<sub>2</sub>O pretreatment and passivate the surface with H<sub>2</sub>O adsorption.<sup>1,3</sup>

The Terman method is a standard high-frequency capacitance-voltage (C-V) technique for analyzing  $D_{it}$ .<sup>18,19</sup> It relies on the interface states to change occupancy along with the direct-current (DC) bias sweep. By making a point-by-point comparison between a measured C-V curve containing responses from the interface states and a calculated ideal reference curve assuming no interface states,  $D_{it}$ can be extracted as a function of band bending. However, the Terman method can grossly

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