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# Comparative study on performance of cubic $Al_xGa_{1-x}N/GaN$ nanostructures MODFETs and MOS-MODFETs



**Superlattices** 

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#### ABSTRACT

We report some comparative results on cubic Al<sub>x</sub>Ga<sub>1-x</sub>N/GaN nanostructures MODFET and MOS-MODFET. The drain current characteristics of cubic Al<sub>x</sub>Ga<sub>1-x</sub>N/GaN MODFET and MOS-MODFET are simulated by changing the different device parameters such as Al content x and the cubic GaN buffer layer thickness using 2D next**nano<sup>3</sup>** numerical simulation software. Drift-diffusion model has taken for simulating the proposed device. These results clearly indicate that the transistor simulation with 5 nm isolator SiO<sub>2</sub> layer thickness under the gate, Al content of x = 25% and 200 nm cubic GaN buffer laver thickness shows the tremendous I-V characteristics. Also, this structure shows an increase of the drain saturation current and a decrease in the threshold voltage. Moreover, our simulation results exhibited lower threshold voltage and higher drain current density of MOS-MODFET is a factor 30% higher than the same current of a conventional MODFET. The MODFET with 5 nm isolator SiO<sub>2</sub> layer thickness has been much better performance. To avoid current flow through the high conductive 3C-SiC substrate a 150 nm p-doped cubic GaN layer is deposited. A comparison between our experimental and simulation results are shown to be in good agreement for cubic Al<sub>0.25</sub>Ga<sub>0.75</sub>N/GaN nanostructures MOS-MODFET. The demonstrated MOS-MODFET will be attractive for the next-generation microwave and high power switching application fields.

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#### 1. Introduction

Beside the impressive applications of group III nitrides in optoelectronics, the number of electronic devices based on wide gap III-nitrides is presently increasing. Due to the large band gap of GaN and AlN and their ternary compounds they are good candidates for high frequency and high power devices and tremendous research activities have been undertaken to grow multilayer group III-nitride stacks in still improving crystalline perfection. Most of these activities have been devoted to group III-nitrides with wurtzite (hexagonal) crystal structure.

However, group III-nitrides can also been grown in the less energetically favorable cubic (Zinc blende) structure, which due to the absence of built-in electric fields have some potential advantages over their hexagonal counterparts where built-in electric fields arising from spontaneous and piezoelectric polarization may set limits to the performance of some electronic devices. Thus the cubic structure may be advantageous for some device applications like normally-off field effect transistors. Also phonon scattering is lower in the cubic phase due to the higher crystallographic symmetry. As a result the mobility of electrons and holes in the cubic phase is expected to be larger than in hexagonal structures.

 $Al_xGa_{1-x}N/GaN$  hetero-junction field-effect transistors (HFETs) are of interest for high-power and high-frequency amplifiers. This is motivated by their potential in commercial and military applications, e.g. in communication systems, radar, wireless stations, high-temperature electronics and high-power solid-state switching. Currently, state of the art HFETs are fabricated on c-plane wurtzite  $Al_xGa_{1-x}N/GaN$  heterostructures. Their inherent polarization fields produce extraordinary large sheet carrier concentrations at the  $Al_xGa_{1-x}N/GaN$  hetero-interface which are advantageous for normally-on type transistors [1–3].

However, for switching devices and digital electronics field-effect transistors (FETs) with normallyoff characteristics are desirable. It was discussed by Abe et al. [4] that the use of cubic III-nitrides would offer fabricating HFETs without parasitic polarization fields and with equal electrical properties for all gate orientations. Further, with cubic nitrides the same technology for the production of normally-on and normally-off devices can be applied. A first demonstration of a c-nitride HFET with normally-off characteristics with a threshold voltage of  $V_{th} = 0.6$  V has recently been reported by Tschumak et al. [5]. The critical issue in the operation of this device was its high gate leakage current, which is undesirable for high power and low noise applications and severely reduces the device performance. Therefore, the use of metal/insulator layers instead of a Schottky gate, leading to the metalinsulator-semiconductor heterostructures field effect transistors (MIS-HFET), may improve the device characteristics [6,7]. Zado et al. [8] have fabricated different MIS structures on cubic GaN.

Moreover,  $Al_xGa_{1-x}N/GaN$  based MOS-HEMT structure was fabricated by Yue et al. [9] is used to validate the present model. For MOS devices, the gate insulator plays an important role, since the good quality dielectric layers would yield these devices with outstanding performances. Metal–Oxide–Semiconductor **Mo**dulation-Doped-Field-Effect-Transistor (MOS-MODFET) structure with a high quality insulating dielectric layer, inserted between the gate metal and  $Al_xGa_{1-x}N$  is expected to offer reduced leakage current. Many dielectrics, including native oxide, SiO<sub>2</sub> [10] have been used as the insulating dielectric in  $Al_xGa_{1-x}N/GaN$  based MOS-MODFET [11].

However, the absence of spontaneous and piezoelectric fields in the cubic AlGaN/GaN system allows the fabrication of normally-on and normally-off heterojunction field-effect transistors (HFETs). However, the critical issue in the operation of the fabricated HFETs was the insufficient insulation of the gate contact. To improve the gate characteristics an insulating layer is required. Low interface trap density metal-insulator-semiconductor (MIS) interfaces are essential for high-power and high-frequency devices [12].

Therefore, the aim of our work is comparative study of the electrical characteristics of cubic  $Al_xGa_{1-x}N/GaN$  nanostructures MODFET [13] and MOS-MODFET switching devices using 2D next**nano**<sup>3</sup> numerical simulation software [14], changing the different device parameters such as the Al content *x* and the cubic GaN buffer layer thickness at  $V_{DS} = 0 V$  for the output characteristics and at  $V_{GS} = 0 V$  for the transfer characteristics of the MODFET and MOS-MODFET. Al<sub>x</sub>Ga<sub>1-x</sub>N/GaN

MOS-MODFET has received great attention due to their lower gate leakage current and larger gate voltage swing compared to the conventional Schottky gate HFETs (SG-HFET) [15–17].

Furthermore, we have assumed 5 nm gate oxide thicknesses. The stability of the device parameters determines the maximum acceptable operating voltage below 5 nm of gate oxide [18–22] with x = 25% Al content of the barrier layer and 200 nm of cubic GaN buffer layer thickness have been improving the performances of cubic Al<sub>x</sub>Ga<sub>1-x</sub>N/GaN nanostructures MODFET, with decreasing of buffer leakage by thinner c-GaN buffer layer. However, this device offered high parallel conductivity through the thick c-GaN buffer [23]. This is decreasing the trap states at the interface of hetero-structures results the reduction scattering processes, whereas decreasing the parasite source resistance to avoid gate leakage current and current collapse. Thus, our simulation results shown that the MOS-MODFET superior performance to that of conventional MODFET, compared with the results similar which obtained by Narihiko Maeda et al. for the design of insulator/Al<sub>x</sub>Ga<sub>1-x</sub>N structures in MOS Al<sub>x</sub>Ga<sub>1-x</sub>N/GaN HFET for higher device performance [24] and by Ki-Yeol Park et al. for the comparative study on Al<sub>x</sub>Ga<sub>1-x</sub>N/GaN nanostructures based MODFET and MOS-MODFET.

#### 2. Simulation results

In this work, we have simulated the *I–V* characteristics of two different devices such as cubic  $Al_xGa_{1-x}N/GaN$  nanostructures MODFET and MOS-MODFET by using 2D next**nano**<sup>3</sup> numerical simulation software. We have assumed in our calculations 10  $\mu$ m for a long the devices structures with 2  $\mu$ m gate length.and without strain.

#### 2.1. Design considerations

Fig. 1 shows the device structure of cubic  $Al_xGa_{1-x}N/GaN$  nanostructures MOS-MODFET with *p*-type doping sample consisting of 150 nm *p*-doped (UID) cubic GaN buffer layer (doping concentration of  $N_A = 1 \times 10^{17}$  cm<sup>-3</sup>), where deposited on the 3C-SiC substrate in (001) direction. The 3C-SiC substrate is highly doped with a free electron concentration of  $3 \times 10^{18}$  cm<sup>-3</sup>. On top 50 nm of the UID cubic GaN buffer layer with a net doping concentration of  $N_D = 1 \times 10^{17}$  cm<sup>-3</sup>, followed by the growth of a 4 nm of UID cubic  $Al_xGa_{1-x}N$  spacer layer with  $N_D = 5 \times 10^{17}$  cm<sup>-3</sup>, 6 nm of cubic  $Al_xGa_{1-x}N$ . Si with  $N_D = 5 \times 10^{18}$  cm<sup>-3</sup>, and with 5 nm of



Fig. 1. Schematic cross-section of cubic Al<sub>x</sub>Ga<sub>1-x</sub>N/GaN nanostructures MOS-MODFET with p-type doping.

the SiO<sub>2</sub> insulator thickness. To avoid current flow through the high conductive 3C-SiC substrate a 150 nm *p*-doped cubic GaN layer is deposited. This results in two p-n junctions at the SiC/*p*-GaN, and the *p*-GaN/*n*-GaN interfaces forming high insulating layers. Finally, the MOS-MODFET sample device has simulated using two-dimensional numerical simulation next**nano<sup>3</sup>** software for the calculations of the electrical output and transfer characteristics.

#### 2.2. Results and discussion

For the validation of our model, we present comparison between the simulation and experimental results. Fig. 2 shows the experimental data and simulation results of the transfer characteristics of cubic  $Al_{0.25}Ga_{0.75}N/GaN$  nanostructures MOS-MODFET were compared, for the purpose of calibrating and validating the simulation results. The model parameters is adjusted to get a perfect matching between experimental and simulation results is shown in Fig. 2 for the transfer characteristics of cubic  $Al_xGa_{1-x}N/GaN$  nanostructures MOS-MODFET at  $V_{DS} = 0$  V and at room temperature. After matching is done, the model is then applied for simulating our model proposed of cubic  $Al_{0.25}Ga_{0.75}N/GaN$  nanostructures MOS-MODFET. The numerical simulator uses drift-diffusion model to solve the Poisson's equation self-consistently with carrier continuity equation. We have assumed the mobility-model-simba-2, if the exponents kappa<sup>n,p</sup> are temperature dependent then this equation is called Canali model (with  $\mu^{n,p}$  as the low field mobility) [14] with suitable modifications to precisely capture the non-equilibrium carrier transport:

$$\mu^{p,n}(\vec{E}) = \frac{\mu^{p,n}}{\left[1 + \left(\mu^{p,n} \frac{|\vec{E}|}{v_0^{p,n}}\right)^{\kappa^{p,n}}\right]^{1/\kappa^{p,n}}}$$
(1)

The device transfer characteristics are shown in Fig. 2, a good matching between the experimental and simulation results of the transfer characteristics at  $V_{DS} = 0$  V with the simulation results, thus the model is validated, extensive simulation of cubic Al<sub>x</sub>Ga<sub>1-x</sub>N/GaN nanostructures MOS-MODFET is then has been done.



**Fig. 2.** (Color lines) represent transfer characteristics of cubic  $Al_xGa_{1-x}N/GaN$  nanostructures MOS-MODFET fabricated at room temperature. The dotted brown line represents calculated transfer characteristics at VDS = 0 V. (For interpretation of the references to color in this figure legend, the reader is referred to the web version of this article.)



Fig. 3. The electrical output characteristics of cubic  $Al_xGa_{1-x}N/GaN$  nanostructures MOS-MODFET with different Al content are calculated at room temperature.



Fig. 4. The electrical transfer characteristics of cubic  $Al_xGa_{1-x}N/GaN$  nanostructures based MOS-MODFET at different Al content are calculated at room temperature.

#### 2.2.1. Variation in Al content x within the $Al_xGa_{1-x}N$ barrier layer

Figs. 3 and 4 show the simulated electrical characteristics of the cubic  $Al_xGa_{1-x}N/GaN$  nanostructures MOS-MODFET, for different values of Al content *x* in the cubic  $Al_xGa_{1-x}N$  barrier layer, (25%, 30%, 35%, 40%, and 45%) are calculated at room temperature. At *x* = 25% (black line), the MOS-MODFET reaches a maximum saturation drain current density is shown in Fig. 3. On the other hand, Fig. 4 shows the electrical transfer characteristics at  $V_{DS} = 0$  V, where the MOS-MODFET reaches a low value of threshold voltage at *x* = 25%. Thus, the current flowing in the device depends strongly on the Al contents of the top layer. This is essentially due to increase in the current drain saturation of the normally-off MOS-MODFET.



**Fig. 5.** The current–voltage characteristics of cubic Al<sub>0.25</sub>Ga<sub>0.75</sub>N/GaN nanostructures MOS-MODFET with different thickness of cubic GaN buffer are calculated at VGS = 0 V and at room temperature.



Fig. 6. The electrical output characteristics of cubic Al<sub>0.25</sub>Ga<sub>0.75</sub>N/GaN nanostructures MODFET and MOS-MODFET are calculated at room temperature.

#### 2.2.2. Variation in thickness of cubic GaN buffer layer

Fig. 5 shows the simulated  $I_D$ – $V_{DS}$  characteristics of the model of cubic Al<sub>0.25</sub>Ga<sub>0.75</sub>N/GaN nanostructures MOS-MODFET with different thickness of cubic GaN buffer layer varied from (200 to 800) nm in a step of 200 nm are calculated at  $V_{GS} = 0$  V. Our simulation results indicate that the fabrication of the MOS-MODFET nanostructures by using a thinner cubic GaN buffer layer successfully decreasing the buffer leakage current, which is critical to obtain a maximum drain current. Thus, the current flowing in the device increases with reduction of the impurity scattering in the buffer layer, thereby increasing electron mobility within the channel and yielding a high device transconductance.



Fig. 7. The electrical transfer characteristics of cubic  $Al_{0.25}Ga_{0.75}N/GaN$  nanostructures MODFET are calculated at room temperature.



Fig. 8. The electrical transfer characteristics of cubic Al<sub>0.25</sub>Ga<sub>0.75</sub>N/GaN nanostructures MOS-MODFET are calculated at room temperature.

## 2.2.3. Comparative study on device characteristics of cubic $Al_xGa_{1-x}N/GaN$ nanostructures MODFET and MOS-MODFET

Fig. 6 shows the comparison of electrical output characteristics when the gate bias  $V_{GS}$  is at 0 V. Cubic Al<sub>0.25</sub>Ga<sub>0.75</sub>N/GaN nanostructures MODFET with and without SiO<sub>2</sub> gate insulator have  $I_{DSS}$  of 439 and 141 A/m, respectively at  $V_{DS}$  = 0.4 V. This result may be attributed to the increase in effective carrier concentration in the channel in MODFET with insulating gate layer. From a comparison of the saturation currents of this device, the different influences of the insulating cap layer become evident. Compared with a free-surface of MODFET without SiO<sub>2</sub> gate insulator, the structures with an insulating cap layer have a lower surface-state density that reduces the surface-related depletion effects. Moreover, the structure with a high-resistively capping layer reduces the electronic density of states at the interface with cubic Al<sub>0.25</sub>Ga<sub>0.75</sub>N. It is clear from the figure that for same value of gate and for same values of drain bias, the drain current is higher for Al<sub>0.25</sub>Ga<sub>0.75</sub>N/GaN MOS-MODFET than that a conventional counterpart. This leads to a reduction of surface-related depletion of the channel layer and increases the sheet carrier concentration [26] by about 32%. Figs. 7 and 8 show the transfer characteristics of the MODFET and MOS-MODFET, respectively. The gate current of MOS-MODFET is smaller than that of a conventional MODFET and therefore present high device transconductance. Our simulation results shown that the electrical performance can be enhanced through the surface gate insulation.

#### 3. Conclusions

The evaluation of the drain current characteristics of cubic  $Al_xGa_{1-x}N/GaN$  based MOS-MODFET has been presented to compare the device performance. The proposed model design can easily be used to predict the characteristics of MODFET as well. The MOS-MODFET structure shows clearly an advantage over their MODFET counterparts in terms of gate voltage swing and higher saturation currents. With 5 nm isolator SiO<sub>2</sub> layer thickness under the gate, Al content x = 25% of cubic Al<sub>x</sub>Ga<sub>1-x</sub>N layer and a thinner cubic GaN buffer layer successfully decreasing the buffer leakage current, which is critical to obtain a maximum drain current. The simulation results suggest that with a lower thick of  $SiO_2$ gate insulator exhibited excellent DC characteristics of cubic  $Al_xGa_{1-x}N/GaN$  FET. The MODFET with 5 nm thick SiO<sub>2</sub> insulator has much better performances. To avoid current flow through the high conductive 3C-SiC substrate a 150 nm p-doped cubic GaN layer is deposited. This results in two p-n junctions at the 3C-SiC/p-GaN, and the p-GaN/n-GaN interfaces forming high insulating layers. Our simulation results have exhibited a maximum saturation drain current with Al content x = 25% at the gate voltage  $V_{GS}$  = 0 V of SiO<sub>2</sub>/cubic Al<sub>x</sub>Ga<sub>1-x</sub>N/GaN nanostructures MOS-MODFET. This model provides a valuable tool for design, optimization, and performance prediction of microwave power devices and demonstrates clear superiority of MOS-MODFET over MODFET for application of optical devices.

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