

GROWTH AND CHARACTERIZATION OF ALGaN/GaN HETEROSTRUCTURES FOR ELECTRONIC DEVICES AND SENSORS

B. Paszkiewicz, M. Wosko, R. Paszkiewicz*, M. Tlaczala

The Faculty of Microsystem Electronics and Photonics, Wrocław University of Technology,
Janiszewskiego 11/17, 50-372 Wrocław, Poland

ABSTRACT

The influence of the Si dopant concentration and its distribution through the AlGaN barrier layer of the AlGaN/GaN heterostructures on their electrical properties was studied. Three types of the heterostructures were grown by MOVPE method on sapphire substrate. Electrical properties of the $\text{Al}_x\text{Ga}_{1-x}\text{N}/\text{GaN}$ heterostructures were determined by the impedance spectroscopy method. The carrier concentration in AlGaN/GaN heterostructure, the incremental sheet charge concentration and 2DEG sheet charge concentration were obtained and correlated with Si dopant distribution through the AlGaN barrier layer of the heterostructure.

INTRODUCTION

The AlGaN/GaN material system is attractive for numerous device applications [1]. The relatively well developed research areas are photonic and high-power electronic devices based on nitrides. Recently, in addition to these, new trends are emerging which indicate that chemical and biochemical sensor applications of AlGaN/GaN heterostructures could benefit from unique properties of nitrides such as their excellent chemical stability and inertness as well as the possibility of formation of high-density two dimensional electron gas (2DEG) on the AlGaN/GaN hetero-interface, even without intentional doping. Its existence near the heterostructure surface is extremely beneficial for highly sensitive detection of surface phenomena. Nitrides heterostructures are also one of the leading candidates for high frequency application up to THz range (regime). However, up to now the parameters of AlGaN/GaN heterostructures do not reached the predicted, theoretical, values. Because of the lack of commercially available bulk GaN substrates the AlGaN/GaN heterostructures are typically grown on sapphire, SiC or Si substrates. The very large mismatch of lattice parameters and thermal expansion coefficients between nitrides and these substrates causes many problems during the growth and influenced the quality of the heterostructures. The choice of the substrate results from the applications and it strongly determines the properties of AlGaN/GaN heterostructures. The

* e-mail: Regina.Paszkiwicz@pwr.wroc.pl

basic theoretical correlations between the thicknesses and composition of AlGa_xN barrier layer, its doping level and spontaneous and piezoelectric polarization are relatively well established. It was shown that the Al content in the AlGa_xN barrier of the heterostructure influenced mainly spontaneous and piezoelectric polarization and its breakdown voltage [2]. The optimization of MOVPE (Metalorganic Vapour Phase Epitaxy) growth process parameters optimization was studied extensively [3-5]. It was established that the MOVPE system configuration as well as the details of the growth process strongly influenced the obtained results. Additionally, our previous study has shown that the 2DEG parameters at the AlGa_xN/GaN heterointerface are not influenced by the parameters of HT GaN buffer layer in a wide range of its carrier concentration [6]. It is in a good agreement with recently published results [7]. But till now it seems that many unidentified factors exist that influence the electrical properties of the AlGa_xN/GaN heterostructures. One of them is the distribution of Si dopant in AlGa_xN barrier which influences on the 2DEG parameters at AlGa_xN/GaN heterointerface. To study these phenomenon three sets of AlGa_xN/GaN heterostructures were grown by MOVPE in which the AlGa_xN layer of AlGa_xN/GaN heterostructure was divided into three sub-layer of different thickness and different Si dopant concentration. The laser interferometer with 635 nm laser diode was used for *in-situ* characterization of the whole process growth mechanism. The electrical parameters of the heterostructures were evaluated by impedance spectroscopy method using the procedure work-out by us and verified by others [8,9]. The electrical properties of AlGa_xN/GaN heterostructures such as the carrier concentration in AlGa_xN/GaN heterostructure and the sheet carrier concentration of 2DEG were evaluated and correlated with Si dopant distribution through the AlGa_xN barrier layer of the heterostructure.

EXPERIMENTAL

Si doped and un-doped AlGa_xN/GaN heterostructures were grown on c-plane sapphire substrates by using 3×2" Thomas Swan Close Coupled Showerhead MOVPE system at 100 mbar. Trimethylgallium, Trimethylaluminum and NH₃ were used as a Ga, Al, and N source respectively, and monosilane (SiH₄) was used as a n type dopant. H₂ was used as a carrier gas.

The growth was performed at six main stages: 1 - sapphire substrate annealing in hydrogen atmosphere, 2 - substrate nitridation in the mixture of H₂ and NH₃ (1:1), 3 - low temperature growth of GaN nucleation layer (LT-GaN NL) at 530°C, 4 - coalescence of LT - GaN NL during temperature ramping, 5 - growth of high temperature GaN buffer layer at 1035°C, 1000 V/III molar ratio, 6 - growth of undoped and/or Si-doped AlGa_xN layer at 1060°C, 1200 V/III molar ratio. The schematic drawings of the MOVPE process sequence for all sets of samples are presented in *Fig. 1*. The laser interferometer with 635 nm laser diode was used for *in-situ* characterization of the growth process that

allowed us to evaluate the thickness of individual layers. Three sets of samples, with different AlGa_{0.2}N sub-layers, were fabricated on 2,5 μm optimized HT-GaN buffer grown on LT-GaN nucleation layer [6]. The designed arrangement of AlGa_{0.2}N barrier sub-layer in individual heterostructure is shown in Fig. 2.

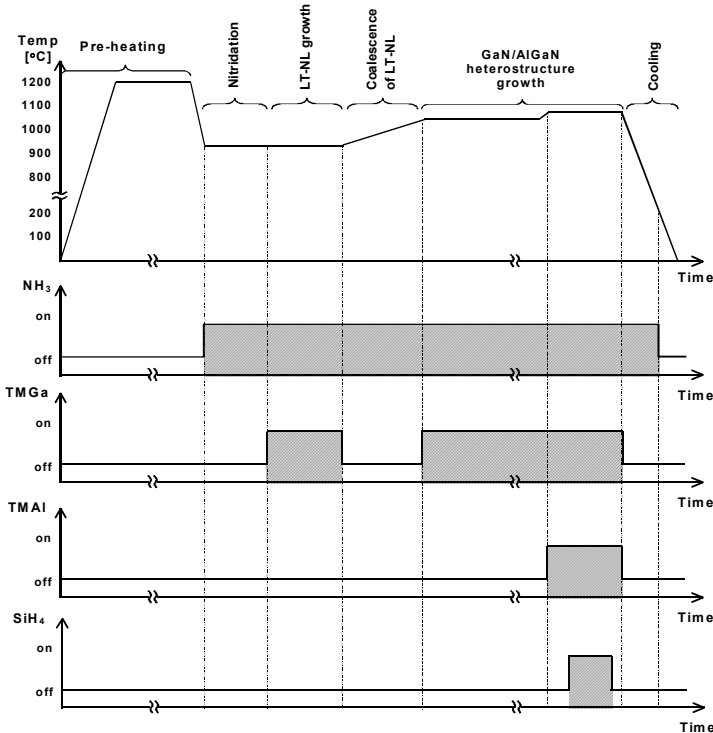


Fig. 1 – The schematic drawing of the MOVPE process sequence of AlGa_{0.2}N/GaN heterostructures grown on LT- GaN NL

The AlGa_{0.2}N/GaN heterostructures #045 consists of, looking from top to bottom, 25 nm thick undoped Al_{0.2}Ga_{0.8}N barrier grown on 2,5 μm thick undoped high resistive HT-GaN buffer. In Al_{0.2}Ga_{0.8}N heterostructures #042 barrier layer was divided in to two sub-layers, from top to bottom, 10 nm thick undoped Al_{0.2}Ga_{0.8}N sub-layer and 15 nm thick Si-doped Al_{0.2}Ga_{0.8}N sub-layer (Si concentration was $2,1 \cdot 10^{18} \text{ cm}^{-3}$). The barrier layer of Al_{0.2}Ga_{0.8}N heterostructures #044 was divided in to three sub-layers, from top to bottom, 5nm thick undoped Al_{0.2}Ga_{0.8}N sub-layer, 20 nm thick Si-doped Al_{0.2}Ga_{0.8}N sub-layer (Si concentration was $4.2 \cdot 10^{18} \text{ cm}^{-3}$) and 2 nm thick undoped Al_{0.2}Ga_{0.8}N sub-layer.

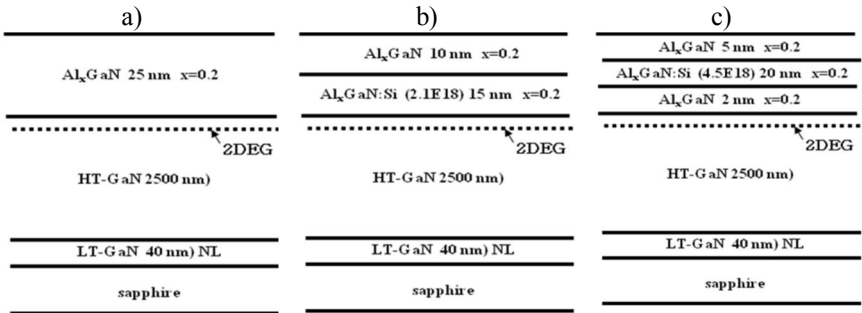


Fig.2. – The schematic cross section of AlGaIn/GaN heterostructures: # 045 (a), # 042 (b) and # 044 (c)

RESULTS AND DISCUSSION

Electrical properties of the $\text{Al}_x\text{Ga}_{1-x}\text{N}/\text{GaN}$ heterostructures were determined by the impedance spectroscopy method performed in the range of frequencies from 80 Hz to 10 MHz with a HP 4192A impedance meter, using a two contact mercury probe. The capacitance and conductance versus frequency characteristics of the Schottky contact to AlGaIn/GaN heterostructures were measured over a range of DC biases and the results were fitted to a worked-out model. The distributed elements equivalent circuit model regarding the series resistances and the Schottky junction admittance was used to evaluate the electrical properties of AlGaIn/GaN heterostructures [8]. The carrier concentration distribution through the epitaxial structure and the incremental sheet charge concentration versus bias voltage were obtained for every set of the samples (Fig. 3, Fig. 4 and Fig. 5).

In Fig. 3 the carrier concentration distributions through the AlGaIn/GaN heterostructures versus the distance from their surfaces are shown. In all heterostructures the characteristic spikes of the concentration at AlGaIn/GaN interface, caused by 2DEG formation, were observed. For detailed analysis the excerpt of the initial part of the Figure 3 is presented in Fig. 4. The incremental sheet charge concentration in AlGaIn/GaN heterostructures versus bias voltage is shown in Fig. 5 that illustrates the depletion of the channel with voltage. Based on data shown in Fig. 3 and Fig. 5 the 2DEG sheet carrier concentration and pinch off voltage of every AlGaIn/GaN heterostructure could be evaluated. In the Table 1 the values of incremental sheet charge concentration (Δn_s), the 2DEG sheet charge concentration (n_s), pinch of voltage (U_p), average mobility ($\bar{\mu}$) and maximum mobility (μ_{\max}) measured and evaluated for all the studies AlGaIn/GaN heterostructures are summarized.

Table 1 – Electrical characteristic of the heterostructures

N°	Δn_s [cm ⁻²]	n_{s2DEG} [cm ⁻²]	U_p [V]	R_s [Ω]	$\bar{\mu}$ [cm/V*S]	μ_{max} [cm/V*S]
045	$2.0 \cdot 10^{12}$	$2.0 \cdot 10^{12}$	-0.9	1732	1708	2400
042	$3.7 \cdot 10^{12}$	$3.1 \cdot 10^{12}$	-1.6	1592	1055	1150
044	$6.1 \cdot 10^{12}$	$3.6 \cdot 10^{12}$	-3.0	1423	719	1155

It was observed that in type #044 heterostructures the edge of depletion layer is located at AlGa_{0.3}N barrier and that the initial carrier concentration in AlGa_{0.3}N barrier (below 20 nm from the surface) is equal to the Si dopant concentration. The increase in bias voltage up to -1.1 V caused the removal of electrons from the barrier layer (Fig. 5). Above the -1.1 V bias voltage the removal of 2DEG electrons starts and at voltage bias equal to -3 V the channel is pinched off. From charge balance it could be seen that for type #044 heterostructures a part of formally evaluated as 2DEG electrons, with concentration equal to $6,1 \cdot 10^{12}$ cm⁻², are electrons at AlGa_{0.3}N barrier with concentration $2,5 \cdot 10^{12}$ cm⁻². The mobility of these electrons is low; they are not useful for carrier transport at the electron devices.

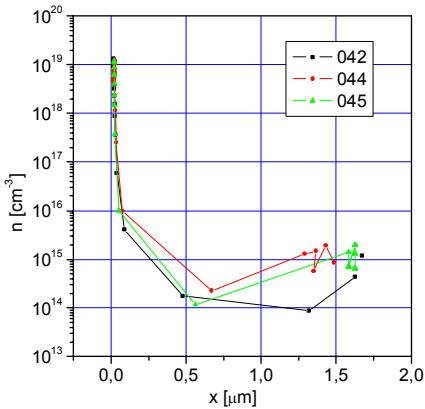


Fig. 3 – The carrier concentration distribution through the AlGa_{0.3}N/GaN heterostructures.

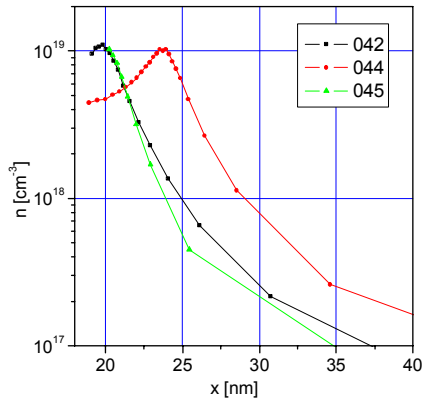


Fig. 4 – The excerpt of the initial part of the Fig. 3.

In #042 AlGa_{0.3}N/GaN heterostructures the total dose of the Si dopant was lower. In Figure 4 an initial increase in the carrier concentration could be observed. The value of charge at AlGa_{0.3}N barrier equal to $0.6 \cdot 10^{12}$ cm⁻² was measured. For this heterostructure the sheet carrier concentration of $3.6 \cdot 10^{12}$ cm⁻² was obtained. It means that the 2 DEG sheet carrier concentrations was $3.0 \cdot 10^{12}$ cm⁻². It seems that for these MOVPE process conditions it is the maximum value

of 2 DEG concentration that could be obtained in Si doped $\text{Al}_{0.2}\text{Ga}_{0.8}\text{N}/\text{GaN}$ heterostructure with Al content in the barrier equal 20%.

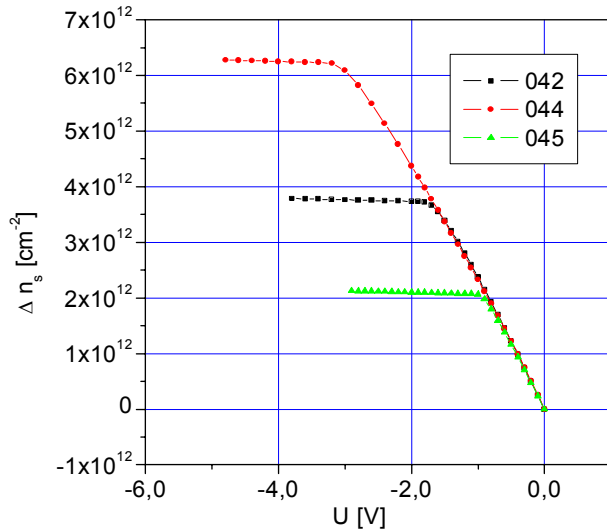


Fig. 5 – The incremental sheet charge concentration versus voltage bias in AlGaIn/GaN heterostructures

In undoped AlGaIn/GaN heterostructures #045 the measured sheet carrier concentration, equal to $2.0 \cdot 10^{12} \text{ cm}^{-2}$, arising from the equilibrium between the charges of surface states, the charges in the channel and the charges induced by spontaneous and piezoelectric polarization.

The performed experiments showed that for $\text{Al}_{0.2}\text{Ga}_{0.8}\text{N}/\text{GaN}$ heterostructure the maximum 2 DEG sheet carrier concentration could be obtained in the range from $2.0 \cdot 10^{12} \text{ cm}^{-2}$ (undoped heterostructures) to $3.6 \cdot 10^{12} \text{ cm}^{-2}$ (Si-doped heterostructures). The increase in the Si dopant concentration above $\sim 2.0 \cdot 10^{18} \text{ cm}^{-3}$ is not efficient. It results in the increase of concentration of the carrier with low mobility that is useless for the electronic devices.

CONCLUSION

The influence of the Si dopant concentration and its distribution through the AlGaIn barrier layer of the MOVPE AlGaIn/GaN heterostructure on carrier concentration in the channel was studied using impedance spectroscopy measurement. It was found that for appropriate Al concentration at AlGaIn barrier the optimum concentration of Si dopant exists which allows us to obtain the heterostructures with good electrical properties. Too high concentration of the Si-dopant could result in decreasing the 2DEG carriers' mobility.

Acknowledgements

This work was co-financed by the European Union within European Regional Development Fund, through grant Innovative Economy (POIG.01.01.02-00-008/08), National Centre for Science under the grant no. N N515 495740, by Wrocław University of Technology statutory grant S10019 and Slovak-Polish International Cooperation Program no. SK-PL-0017-09.

REFERENCES

- [1] B. Boratyński, B. Paszkiewicz, R. Paszkiewicz, M. Tlaczala, *Act. Phys. Pol. A*, 116, 5, (2009), 800.
- [2] S. Heikman, S. Keller, Y. Wu, SP Den Baars, and UK Mishra, *J.App. Phys.*, 93, (2003), 1014.
- [3] S. Heikman, S. Keller, T. Mates, SP Den Baars, and UK Mishra, *J. Cryst. Growth*, 248, (2003), 513.
- [4] Y.S. Cho, N. Kaluza, V. Guzenko, Th. Schäpers, H. Hardtdegen, H.-P. Bochem, U. Breuer, M. R. Ghadimi, M. Fecioru-Morariu, B. Beschoten, *phys. stat. sol. (a)*, 204, 1, (2007), 72.
- [5] K. Köhler, S. Müller, R. Aidam, P. Waltereit, W. Pletschen, L. Kirste, H. P. Menner, W. Bronner, A. Leuther, R. Quay, M. Mikulla, O. Ambacher, R. Granzner, F. Schwier, C. Buchheim, and R. Goldhahn, *J. Appl. Phys.* 107, (2010), 053711.
- [6] B. Paszkiewicz, M. Wosko, R. Paszkiewicz, M. Tlaczala, *Proceedings of European Workshop on Metalorganic Vapour Phase Epitaxy*, 5-8 June, Wrocław, Poland, C11, (2011), 195.
- [7] S.M. Hubbard, G. Zhao, D. Pavlidis, W. Sutton, E. Cho, *J. Cryst. Growth*, 284 (2005) 297.
- [8] B. Paszkiewicz, *J. Cryst. Growth*, 230, 3-4, (2001), 590.
- [9] J.C. Martinez-Orozco, L.M. Gaggero-Sager, Stoyan J. Vlaev, *Solid-State Electron.*, 48, (2004), 2277.