

Novel Si stacking on β -Ga₂O₃ MOSFET structure enabling 0.59 m Ω cm² of Ron,sp at 600V blocking voltage

Takashi Miida¹, Ikuo Kurachi², Kazuo Tanimoto¹, Kazuhiro Kawasaki¹

1. EASTWIND, LLC, Kobe, Hyogo, Japan, Email; miida@trio.plala.or.jp
2. D&S, Inc, Hachioji, Tokyo, Japan

Abstract—Ga₂O₃ base vertical trench MOSFET achieving the normally-off operation is newly proposed, in which thin Si-MOSFET is stacking on n-type β -Ga₂O₃ drift layer. The state-of-the-art “Blocking layer” in p-Si is the key to relax the E-field in p-Si at Off-state and to obtain the current stability and the optimum V_{th} at On-state. Using this architecture, the normally-off operation with optimum V_{th} as 4V and an extremely low Ron,sp as 0.59 m Ω cm² are estimated at 600V blocking voltage device.

I. INTRODUCTION

Ga₂O₃ has emerged being thought as a promising material for the power switching devices among the wide band gap (WBG) semiconductor such as SiC or GaN from the cost and performance point of view. Even though Ga₂O₃ has many advantages over Si, it is hard to enable the normally-off type vertical trench transistor, because there exists no p-type Ga₂O₃.

Furthermore, MOS channel mobility of every WBG is very low due to its high surface state density, which rises Ron higher. Due to the lack of p-type Ga₂O₃, FINFET structure [1] or another p-type material [2] has been proposed as the alternative to enable the normally-off operation for MOSFET. However, there has been currently no breakthrough to realize the vertical trench MOSFET that can achieve the high current density.

To overcome these drawbacks with maintaining the other advantages of Ga₂O₃, the novel MOSFET device architecture is proposed and verified by TCAD simulation.

II. DEVICE STRUCTURE

The proposed device is composed with Si-MOSFET stacking on 5 μ m n-type β -Ga₂O₃ drift layer (Nd= 2e16cm⁻³), as shown in Fig.1, where the direct bonding technology [3] can be utilized to form the p-n hetero-junction [4].

The direct bonding technology such as surface activated or fusion bonding enables the strong adhesion between different materials at RT with thermal stress free and an atomic level uniformity at the interface, that can maintain crystal periodicity as same quality as epitaxial growth. The Fig.2 shows the band diagram of Si (100)-Ga₂O₃ hetero-junction at the equilibrium condition. The trench gate penetrates the 1 μ m thickness into p-Si bulk and another 1 μ m depth to n-Ga₂O₃ drift region covering with 100nm gate Si-dioxide, where this film thickness will bring Qg*Ron FOM smaller. In order to protect the trench bottom from the E-field stress, Nitrogen ion implant is applied under the trench bottom, where the isolation layer is formed as the high resistive region [5]. Source (n⁺) and bulk diffusion layer (p⁺) are set on the p-Si surface, and the drain is set at the bottom of n-Ga₂O₃. In particular, the p-type “blocking layer” is partially (~0.3 μ m) formed at the bottom of p-Si bulk with higher doping level (Na= 1e17cm⁻³) than bulk to suppress the leakage between source and drain at “Off-state” and to regulate the channel current conduction at “On-state”. The cell pitch of the proposed device is 2 μ m where the mesa width is 1 μ m.

III. DEVICE OPERATION AND SIMULATED RESULTS

Since it is so important to see what is brought by these kind of emerging technologies is, TCAD is the key to predict the characteristics and potential performance of newly proposed

device structure. Here, AdvanceSoft’s device simulation tool is used applying the physical property of β -Ga₂O₃ cited from the past works [6] [7].

In “On-state”, it is so significant that the hetero junction is to be forced by the E-field from gate, thereby forcing the surface potential of n-Ga₂O₃, wherein Ec of Ga₂O₃ is 0.05eV higher than that of Si as in Fig. 2. It is also remarkable that the thickness of “blocking layer” is regarded as the “effective channel length” of MOSFET wherein the V_{th} and channel conductivity are determined by the thickness and the doping level of the p-type “blocking layer”. Therefore, it will bring the less p-Si bulk thickness dependency on the channel conductance, which ensures the good bonding manufacturability in this proposed device. As a result, 0.59 m Ω cm² of Ron,sp excluding the substrate resistance at Vg= 15V and 4.0V of threshold voltage at Vd= 1V are obtained using Yamaguchi mobility model [8] as shown in Fig.3 and Fig.4, respectively. This excellent performance is brought by the usage of fast channel mobility of Si-MOSFET in this proposed device.

Meanwhile, it is so important for “Off-state” to relax the electric field from Ga₂O₃ to Si less than 1/10. The solution is to narrowing the Mesa width in conjunction with the “blocking layer” explained above, these will prevent the electric field penetration into Si from the punch through or breakdown by the spatial effect and the doping control. Fig.5 shows the 2-D electric field distribution and potential in the whole device region simulated with Neumann condition for mirror boundary.

It is obvious that the averaged electric field within Si is small enough to bring the reciprocal of impact ionization factor α^{-1} [9] becomes 5 μ m depth in Si. Fig.6 shows the comparison for wide/narrow mesa width and with/without blocking layer, respectively. The E-field at the gate bottom oxide is still large, which is around 5MV/cm, though the isolation layer can prevent the leakage current perfectly like an insulator. The E-field distribution can be relaxed by reducing the doping level or extending for the drift layer thickness, those have a trade-off relationship against on Ron. If the gate width is getting narrower, it is obvious that Ron,sp becomes smaller as the cell pitch shrinks less than 2 μ m and below.

IV. SUMMARY

This proposed device structure will provide the fusion solution for Si- and WBG-MOSFET to achieve the high performance. To verify the new structures such as the bonding technology including hetero-junction, TCAD plays an important role to verify the device theory and determines the doping level and geometrical factors for the further trial chip production to optimize.

REFERENCES

- [1] Hu, et.al. IEEE EDL, Vol. 39, No.6, pp. 869-871, 2018
- [2] FLOSFIA Inc. et.al , Press release July, 2018
- [3] Takagi, et.al. ECS Transactions 86(5) pp. 169-174, 2018
- [4] Liang, et.al. ECS Solid State Letters, 4(11) pp. 55-57, 2015
- [5] Tetzner, et.al., Appl. Phys letters. October, 2018
- [6] Gibbons, et.al AIP Advances 8, 065011, 2018
- [7] Pearson, et.al Applied Physics Reviews 5, 011301, 2018
- [8] Yamaguchi, IEEE Electron Device Vol.30 pp. 658-663, 1983
- [9] MAES, et.al. Solid State Electronics, Vol.33, No.6, pp.705-718, 1990.

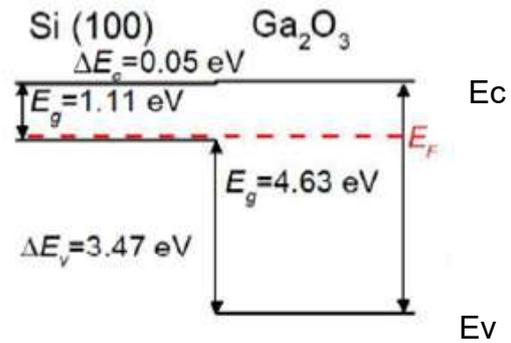
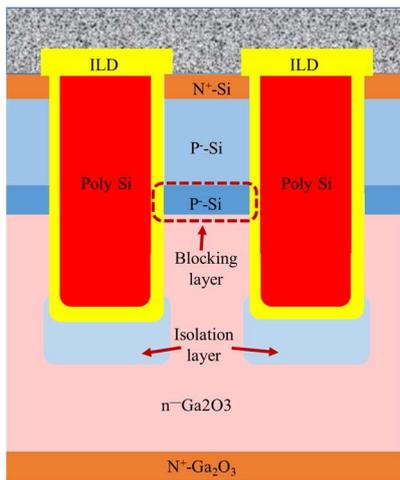


Fig.1 Cross section of the proposed device. SAB is applied to bond the Ga₂O₃ and Si at the interface with Ga₂O₃.

Fig.2 Band diagram of the proposed device illustrated at the equilibrium condition [6]. The electrons will get over the barrier height between Ga₂O₃ and Si by gate biasing.

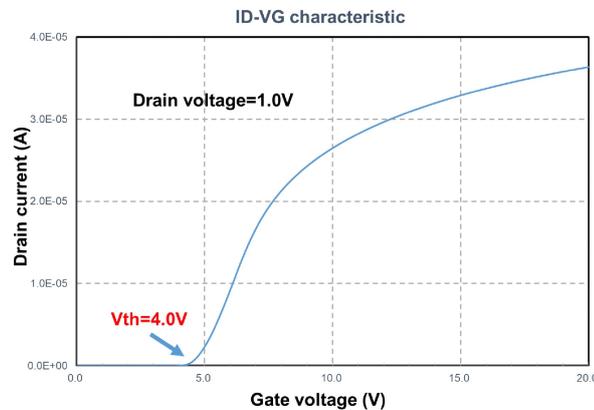
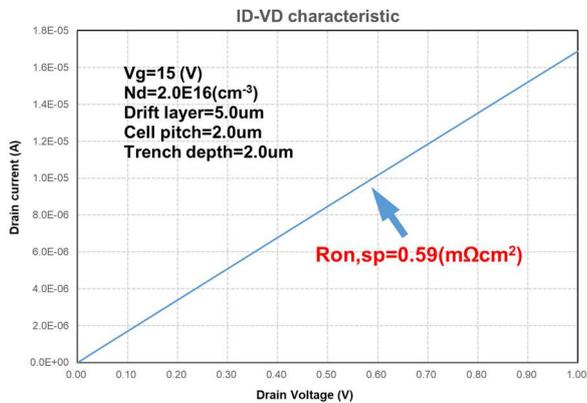


Fig.3 Id vs. V_d characteristic of the proposed device where 0.59mΩcm² of Ron,sp is achieved at Von(V_d)= 1V.

Fig.4. Id vs Vg characteristic of the proposed device, where Normally-off operation is obviously achieved.

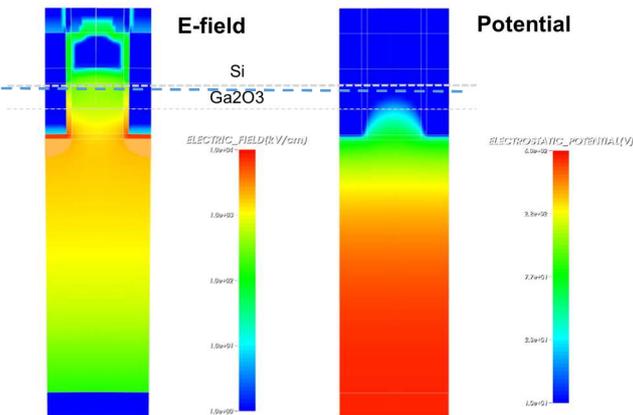


Fig.5 E-field and potential distribution at reversed biasing with 600V on drain.

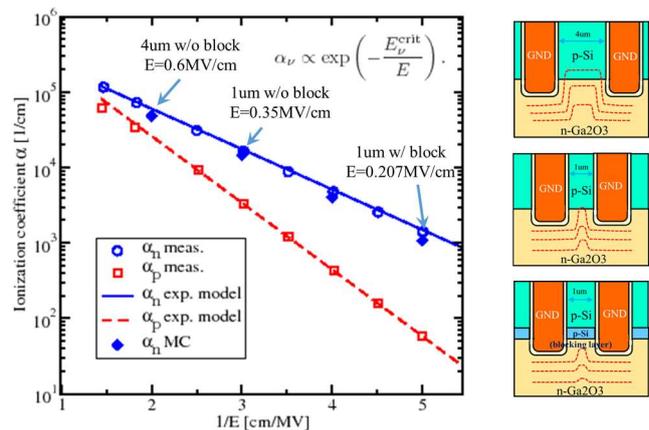


Fig.6. Impact ionization coefficient at the center of the proposed device. 1um channel width and block layer. α⁻¹ will become 5um which is long enough to suppress the impact ionization in Si [9].