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Numerical Approach of Single-Junction InGaN Solar Cell Affected by Carrier Lifetime and Temperature

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Abstract

The PC1D simulation and origin software were successfully used for the study of Carrier Lifetime and Temperature effect on InGaN Single-Junction Solar Cell. For the simulation, the total device area was 100 cm², dielectric constant 13.1, band gap 1.35eV, intrinsic constant is 1×10¹⁰ cm⁻³, doping concentration is 1×10¹⁷cm⁻³, electron number and hole number 1000 and 170 respectively, and the refractive index was 3.58. The optimized temperature and bulk recombination were 25°C and 1000μs respectively along with the efficiency of 18.258 % for both n and p – InGaN solar cell. Several graphs were plotted under the following conditions: a) bulk recombination time of p-InGaN and temperature are kept constant at 1000 μs and 25°C, the variation of bulk recombination time of n – type InGaN solar cell with base current and voltage, maximum current and voltage, and efficiency and maximum power were studied. b) bulk recombination time of n-InGaN and temperature are kept constant at 1000 μs and 25°C, the variation of bulk recombination time of p – type InGaN solar cell with base current and voltage, maximum current and voltage, and efficiency and maximum power were studied.

Keywords: InGaN solar cell, single junction, bulk recombination, PC1D.

Introduction

With the development of technologies, new materials has been developed especially in energy efficient devices. There are two aspects of energy management. Either using efficient devices or generation of energy sources both of which were adopted by us. For the development of energy efficient devices, we have conducted investigations on different materials like, Ferrites [1, 2, 3–10, 11–15,], MXenes(D. Parajuli, 2018; D. Parajuli et al., 2023; D. Parajuli, Murali, K. C, et al., 2022; D Parajuli et al., 2019; Deependra Parajuli & Samatha, 2022), electrodes (D. Parajuli, Murali, Samatha, et al., 2022; D. Parajuli, Taddesse, Murali, Veeraiah, et al., 2022) etc. For the energy sources investigation we have gone through different alternative energy approaches mainly in solar cells [21–23] etc. As our global energy expenditure increases exponentially, it is apparent that renewable energy solution must be utilized. Solar PV technology is the best way to utilize the unlimited solar energy. The InGaN is a recently developed novel solar cell material for its promising tunable band gap of 0.7 eV to 3.4 eV for the realization of high efficiency tandem solar cells in space and terrestrial applications(Akter, 2014). The III-nitride semiconductor material system, which consists of InN, GaN, AlN and their alloys, offers a substantial potential in developing ultra-high efficiency photovoltaics mainly due to its wide range of direct-band gap, and other electronic, optical and mechanical properties. However, this novel InGaN material system possesses challenges from theoretical, as well as technological standpoints, which are further extended into the performance of InGaN devices. The InxGa (1-x)N material system is a promising candidate for developing high-efficiency PV systems. Research involving InxGa (1-x) N photovoltaic devices is still in its initial stages. Since the first proposal in 2003 to use InxGa (1-x)N for solar cell applications, substantial efforts have already been made in this research space (Honsberg et al., 2004).

The first device showing PV response was reported in 2007 [26]. Although InGaN solar cells are still not fully developed, various theoretical models and numerical simulations have been conducted to investigate the performance of single- and multiple-junction InGaN solar cells (Feng et al., 2010; Hsu & Walukiewicz, 2008; Shen et al., 2008; Zhang et al., 2007). In 2011, S. Ben Machiche has achieved efficiency of 24.88% for single junction InGaN solar cell and efficiency of 34.34% and 37.15% for double junction and triple junction tandem solar cell respectively (Bouزيد & machiche, 2011).

There is greater loss in efficiency due to the recombination of charge carriers. The solar radiations reception also depends in many parameters including geographical values. The reflectance, thermodynamic efficiency, charge carrier separation efficiency charge carrier collection efficiency, conduction efficiency are key factors affecting the efficiency of a solar cell system(Green, 1981). The recombination losses are due to quantum efficiency, V_{OC} ratio, and fill factor values. Resistive losses are predominantly accounted for by the fill factor value (Kumar, 2017), but also contribute to the quantum efficiency and V_{OC} ratio values. The highest known efficiency of solar cell was 47.1% was achieved by the use of multi-junction concentrator solar cells, developed at National Renewable Energy Laboratory, Golden, Colorado, USA (Geisz et al., 2018; *Photovoltaic Cell Conversion Efficiency Basics*, 2014).

There are many simulation tools for the study of device simulations like Silvaco TCAD, SCAP-1D, GVPDM, AFORS-HIT, PC1D etc. In this work, PCID simulation is used for the study of InGaN single junction solar cell. The aim of this simulation work is to obtain the maximum conversion efficiency of InGaN single-junction solar cell with the best structure parameters. The effects of the carrier lifetime of each layer on the electrical parameters of the solar cell, such as the short circuit current (I_{sc}),

the open circuit voltage (V), the fill factor (FF), and the conversion efficiency (η), were investigated. Furthermore, the effects of temperature on the conversion efficiency of the single-junction InGaN solar cell were also studied.

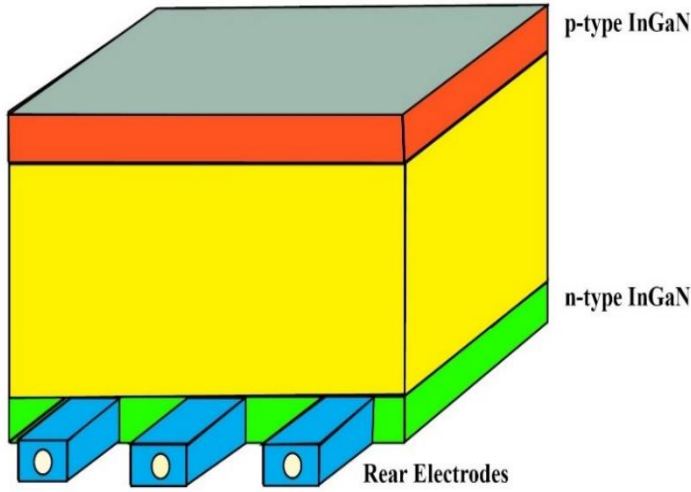


Fig. 1: A 3D schematic diagram of proposed p-n InGaN homojunction solar cell

Methods of Analysis

The PC1D simulation software is used for the computation of the solar cell parameters due to its free of cost, speed, user interface and updated cell model. PC1D is a three step process: 1. Setting up the simulation parameters which includes the device and material parameters, and the excitation to be applied to the device. 2. Running the simulation. 3. Examining the results. In addition, we have used ORIGIN for plotting graph. We have used the free version of this software. There is also the data analysis facility in origin. In the analysis, the statistics, signal processing and curve fitting option. It can imports, ASCII text, Excel, NITDM etc. and can export the images in JPEG, GIF, EPS, TIFF...etc. In this study, we have considered the effect of temperature, bulk combination in the efficiency of the InGaN solar cell. For device area, 100 cm^2 , different sets of reading can be taken and plotted graph between base current and base voltage for different temperatures. Similar set of graphs under the similar input parameters can be drawn for various bulk recombination (carrier lifetime).

The efficiency of a solar cell is determined as the fraction of incident power which is converted to electricity and is defined as:

$$P_{max} = V_{oc} I_{sc} FF$$

$$\eta = V_{oc} I_{sc} FF / P_{in}$$

Where: $FF = \frac{P_{MP}}{V_{oc} \times I_{sc}} = \frac{V_{MP} \times I_{MP}}{V_{oc} \times I_{sc}}$, V_{oc} is the open-circuit voltage; I_{sc} is the short-circuit current; FF is the fill factor and η is the efficiency.

The process of taking the highest efficiency for the study of other solar cell parameters is called optimization. In our case, the maximum efficiency was obtained at 25°C with bulk recombination in $1000 \mu\text{s}$ of P-InGaN and N-InGaN on single junction solar cell which is 18.358.

Result and Discussions

A schematic diagram (3D and 2D) of proposed p-n InGaN homojunction solar cell are shown in figure 1 and 2 respectively with P-layer and n-layer having thickness of 120 nm and 270 nm. At the time of simulation, the thicknesses are $0.5 \mu\text{m}$ and $5 \mu\text{m}$ respectively.

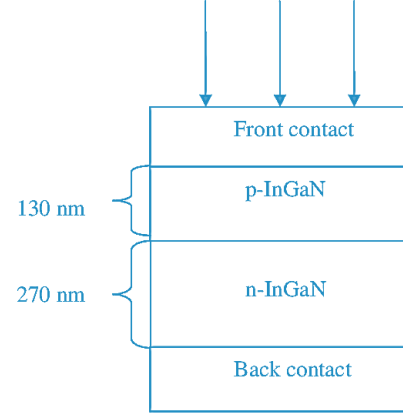


Fig.2: InGaN single junction solar cell

The parameters of the material that are used as the input values in PC1D simulation on InGaN solar cell are listed in table 1.

Table I. Material parameters

| Parameter | Value N- InGaN | Value P InGaN |
|---------------------|------------------------------------|------------------------------------|
| Thickness | $0.5 \mu\text{m}$ | $5 \mu\text{m}$ |
| Dielectric constant | 13.1 | 13.1 |
| Band gap | 1.35eV | 1.35eV |
| Intrinsic conc. | $1 \times 10^{10} \text{ cm}^{-3}$ | $1 \times 10^{10} \text{ cm}^{-3}$ |
| Refractive index | 3.58 | 3.58 |
| Doping | $1 \times 10^{17} \text{ cm}^{-3}$ | $1 \times 10^{17} \text{ cm}^{-3}$ |
| Electron number | 1000 | 1000 |
| Hole number | 170 | 170 |

Carrier concentration effect of n-InGaN Solar cell

We have taken the thickness and bandgap of the n region as $0.5 \mu\text{m}$ and 1.35eV respectively to find the optimized value. From our calculation, the maximum efficiency is obtained at 25°C temperature and the bulk recombination of n - InGaN is $1000 \mu\text{s}$ as optimized value. Different sets are taken for the value of n- InGaN to be 0.001, 0.01, 0.1, 1, 10, 100, $1000 \mu\text{s}$ respectively. Keeping temperature fixed at 25°C , we have recorded the changes of current with voltage. All the obtained values are plotted in a graph as shown in figure 3. The base current or the short circuit current (I_{sc}) seems independent with bulk recombination time, while the open circuit voltage is initially constant, drop suddenly to minimum at the bulk combination 10 and then again increases in the same way.

A graph of I_{\max} against V_{\max} is as shown in figure 4. The P_{\max} and its efficiency initially increases and becomes constant with the bulk recombination. The variation of bulk recombination time of n - InGaN was done at the constant value of temperature and p - InGaN bulk recombination.

Similarly, the graph of maximum power and efficiency with bulk recombination for n-InGaN solar cell is as shown in figure 5. In the figure, the power is increasing more rapidly with bulk recombination than the efficiency and finally both attains steady state. The bulk recombination at 10 (μ s) show significance change in all the graphs.

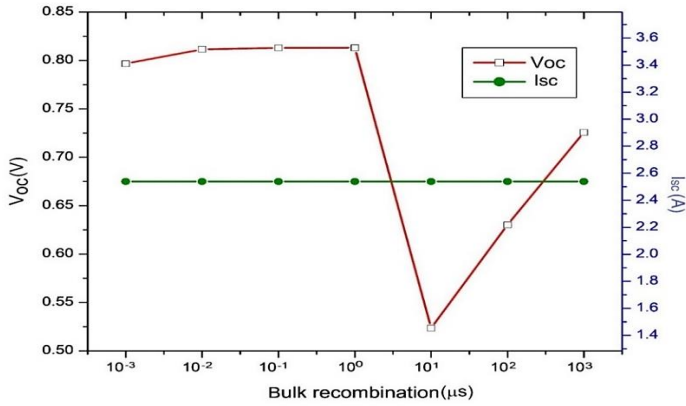


Fig. 3: Graph of variation short circuit and open circuit voltage with bulk recombination of n-InGaN

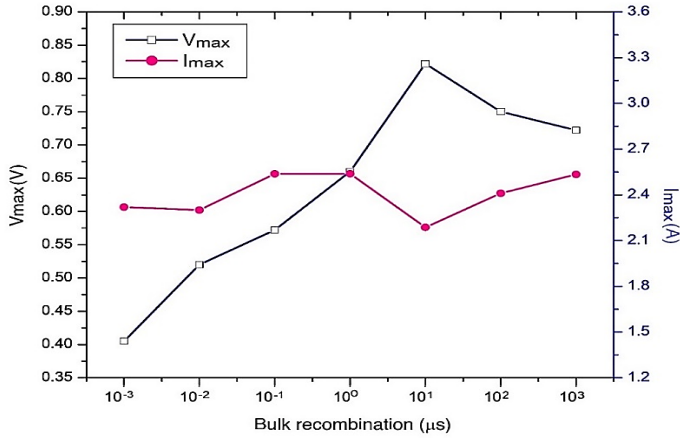


Fig. 4: Graph between I_{\max} and V_{\max} with bulk recombination of n-InGaN

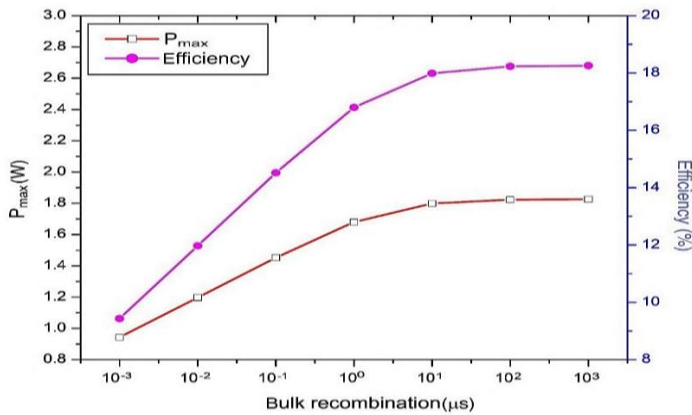


Fig. 5: Graph between P_{\max} and efficiency with bulk recombination of n-InGaN

Carrier concentration effect on p -InGaN Solar cell

The input values for the optimization of p-InGaN solar cell were 0.5 μ m and 1.35eV for thickness and bandgap of the p region respectively. For the optimization, it is assumed that proposed cell has maximum efficiency in the active region. As in the n-type carriers, the bulk recombination time for p-InGaN was varied from 0.001, 0.01, 0.1, 1, 10, 100, 1000 μ s respectively.

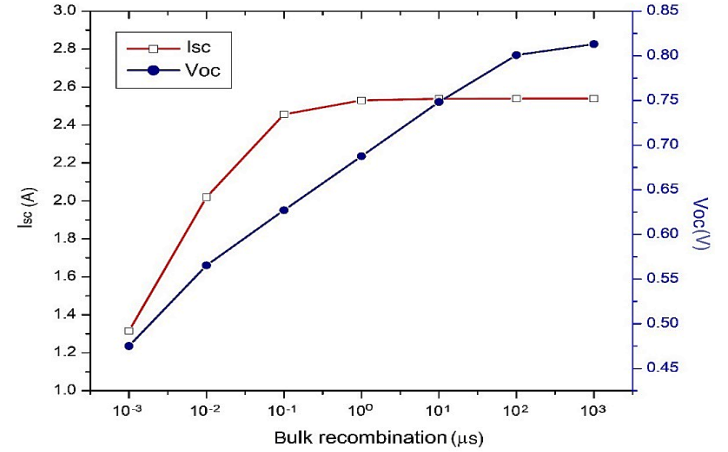


Fig. 6: Graph between the base voltage and current with bulk recombination variation

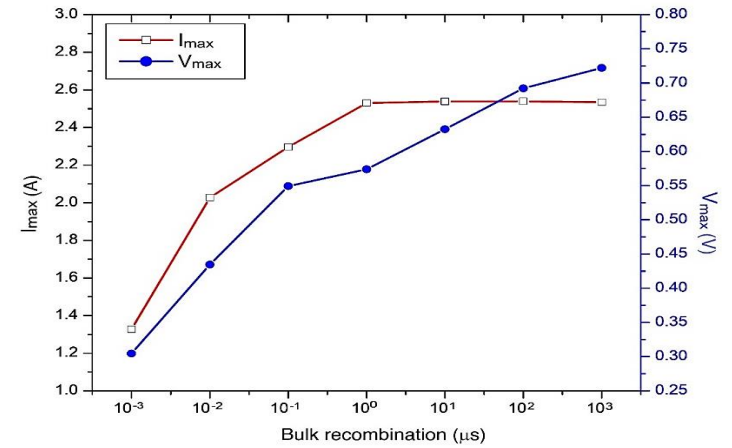


Fig. 7: Graph between I_{\max} and V_{\max} with bulk recombination of n-InGaN

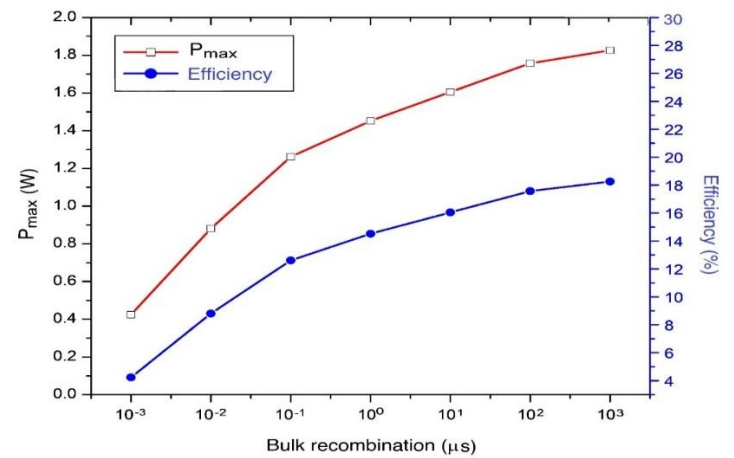


Fig. 8: Graph between P_{\max} and efficiency with bulk recombination in p-region

Similarly, the graph of bulk recombination time with maximum current and voltage at constant temperature and n-type bulk recombination is as shown in figure 7. The maximum current and maximum voltage show the same trend with bulk recombination as for I_{sc} and V_{oc} against bulk recombination.

In the same way, the graph between maximum power and efficiency with bulk recombination of p-type under the same condition is shown in figure 8. The power is increasing more rapidly than efficiency with bulk combination.

Impact of temperature

In this case, we have studied the change in various solar cell parameters with the variation in temperature. Here, both the n and p-type bulk recombination time are kept constant at 1000 μs . We have taken different values of short circuit current and open circuit voltage at various temperature ranging from 25 °C to 50°C. The obtained values were plotted in a graph as shown in figure 9. The open circuit voltage is decreasing and the short circuit current is increasing with the temperature. The increasing current with temperature show the semiconducting nature of the system.

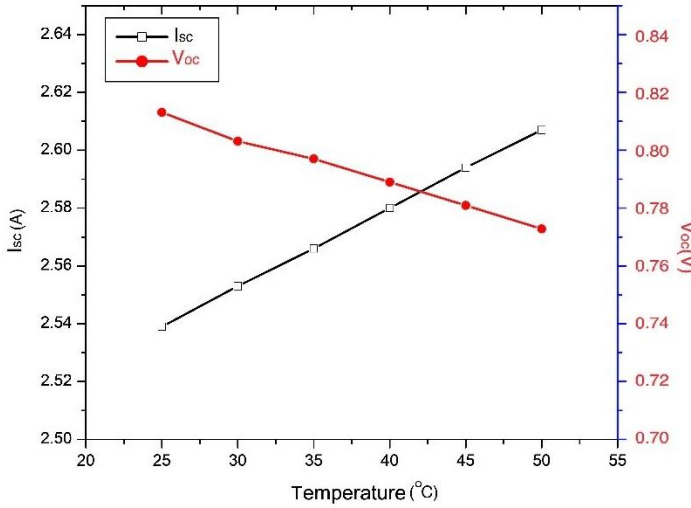


Fig. 9: Graph between I_{sc} and V_{oc} with temperature

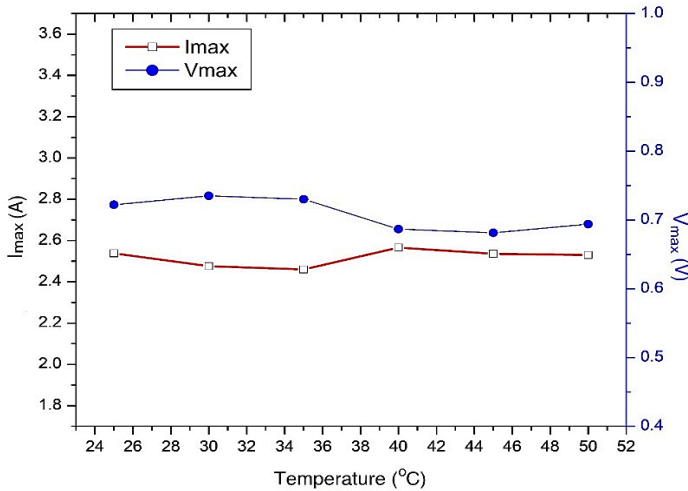


Fig. 10: Graph between I_{max} and V_{max} with temperature

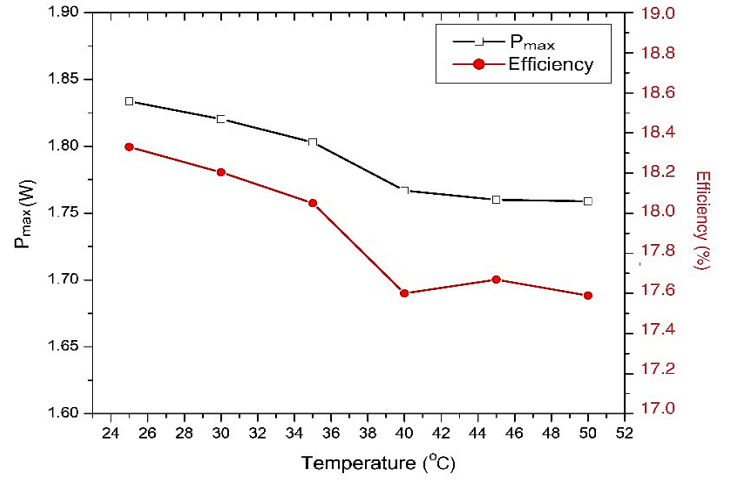


Fig.11: Graph between P_{max} and efficiency with temperature

Similarly, the variation of temperature with maximum voltage and current for p-type InGaN is as shown in figure 10. The current is increasing and voltage is decreasing with the temperature and finally both attain the steady values. The variation of temperature with maximum power and the efficiency is as shown in figure 11. Both the efficiency and maximum power are decreasing with the temperature indicating that the solar cell is not good at higher temperature.

Optimization

In this section, two steps are performed,

- 1) bulk recombination time of p-InGaN and temperature are kept constant at 1000 μs and 25°C, the variation of bulk recombination time of n – type InGaN solar cell with base current and voltage, maximum current and voltage, and efficiency and maximum power were studied.
- 2) bulk recombination time of n-InGaN and temperature are kept constant at 1000 μs and 25°C, the variation of bulk recombination time of p – type InGaN solar cell with base current and voltage, maximum current and voltage, and efficiency and maximum power were studied.

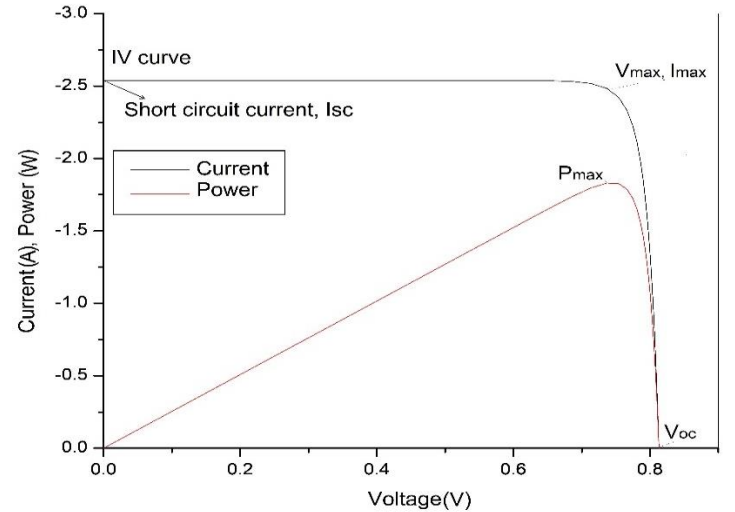


Fig. 12: Base I-V/Power curve

For the InGa_N simulation, the n - InGa_N layer thickness was 0.5 μm , the p - InGa_N layer thickness was 5 μm , the doping concentrations for these layers are $1 \times 10^{17} \text{ cm}^{-3}$, dielectric constant is 13.1, band gap 1.35 eV, intrinsic constant $1 \times 10^{10} \text{ cm}^{-3}$, refractive index 3.58, electron number and hole number 1000 and 170. The obtained results were extracted from the *I-V* curve characteristics for the model as shown in figure 12. The open-circuit voltage (V_{oc}), short-circuit current (I_{sc}), and the efficiency of this initial simulation were found to be 0.8131V, 2.539A, and 18.26 % respectively. The maximum power is obtained from the figure. After all, the optimized value of temperature and bulk recombination were found to be 25°C and 1000 μs respectively.

Conclusions

The PC1D simulation and origin software were successfully used for the study of Carrier Lifetime and Temperature effect on InGa_N Single-Junction Solar Cell. For the simulation, the total device area was 100 cm^2 , dielectric constant 13.1, band gap 1.35 eV, intrinsic constant is $1 \times 10^{10} \text{ cm}^{-3}$, doping concentration is $1 \times 10^{17} \text{ cm}^{-3}$, electron number and hole number 1000 and 170 respectively, and the refractive index was 3.58. The optimized temperature and bulk recombination were 25°C and 1000 μs respectively along with the efficiency of 18.258 % for both n and p - InGa_N solar cell. Several graphs were plotted under the following conditions:

- 1) bulk recombination time of p-InGa_N and temperature are kept constant at 1000 μs and 25°C, the variation of bulk recombination time of n - type InGa_N solar cell with base current and voltage, maximum current and voltage, and efficiency and maximum power were studied.
- 2) bulk recombination time of n-InGa_N and temperature are kept constant at 1000 μs and 25°C, the variation of bulk recombination time of p - type InGa_N solar cell with base current and voltage, maximum current and voltage, and efficiency and maximum power were studied.

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