Research Article

Numerical Simulation of Carbon Nanotubes/GaAs Hybrid PV Devices with AMPS-1D

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1. Introduction

The unique electronic properties of single-walled carbon nanotubes (SWCNTs) are intriguing for many research groups in the present semiconductor technology to incorporate CNTs into CNT/III–V semiconductor hybrid devices [1, 2]. On the other hand, GaAs has the superior electrical and optical properties owing to its diffusion length, well-developed industrial application compared with Si, and some better electrical performance in certain aspects like higher carrier mobility [3]. Thus, from the energy conversion point of view, CNT/GaAs hybrid devices could be a promising optoelectronic combination since both materials are direct band gap semiconductors [4]. The solar cell constructed from SWCNT/GaAs thin film could efficiently absorb light energy with a wide absorption spectrum, if the diameter-dependent bandgap of the SWCNT was controlled through growth conditions from 0 to 2 eV [5, 6]. Albeit the diode behaviour and photovoltaic response of the CNT/GaAs heterojunction interface have been studied, the device modelling and simulations are still under investigation. In this work, for the first time, we used the one-dimensional simulation program called Analysis of Microelectronic and Photonic Structures (AMPS-1D) device simulation tool for the modelling of the characteristics and performance parameters of a zigzag SWCNT/GaAs heterojunction solar cell. To our knowledge this is the first time that a hybrid heterojunction including a nanostructure material is going to be simulated in AMPS simulator or even in a simulator platform. This tool is widely used for the simulation of the thin film or silicon based photovoltaics [7–9].

2. Device Model

There are several methods to bring in contact the SWCNT absorber and GaAs (n-type substrate): chemical vapour deposition [10], followed by electron-beam lithography with assistance of a third material, that is, Si₃N₄ or SiO₂ [4].

A typical solar cell structure has been shown in Figure 1 where the Au/Pd contact, CNT layer, n-type GaAs substrate, and Au/Pd came in contact. Here we ignore any additional layer (e.g., SiC) which in practice is deposited on the GaAs layer to cap the CNT layer. To proceed with the simulation, the material parameters employed as the inputs are selected based on the reported literature values or constrained to
reasonable ranges. The key semiconductor properties of the layers as the input parameters for the simulations are given in Table 1. Although no attempt was made to match the simulation results with the experimental data, the purpose is to analyse the performance parameters of the cell versus the material parameters.

3. Simulation Results and Discussion

The schematic energy-band diagram of a typical hybrid SWCNT/GaAs solar cell has been illustrated in Figure 2. Based on this band diagram, the current-voltage characteristics of this structure have been calculated in the dark and light conditions under one sun at AM1.5 standard conditions (Figure 3) in order to show electrical behaviour of the proposed structure in different operation conditions. The dependence of SWCNT’s acceptor concentration (p) on the performance parameters of the cell has been investigated. In the simulations, shown in Figure 4, obviously there is an optimum value for the p acceptor concentration of about $10^{16}$ cm$^{-3}$ that has a better performance than those with low doping concentration. Increasing the p concentration increases the open circuit voltage, $V_{oc}$, negligibly where it stays almost constant. However, the short circuit current, $I_{sc}$, decreases and the energy conversion efficiency reduces. On the other hand, donor concentration in the n-type GaAs has the same role with optimum value of about $10^{16}$ cm$^{-3}$ and after that while lower than this value the $V_{oc}$ decreases.

In contrast, the $I_{sc}$ increases very slightly by $N_D$, but the efficiency and $FF$ decrease as shown in Figure 5. Higher donor concentration excites more electrons to the conduction band while higher acceptor concentration provides more holes to the valence band which creates better conductivity and photocurrent generation leading to increased efficiency. We need to note that the concentration of the acceptor and donor in the absorber layer is very effective on the performance parameters of the cell since the ratio of $N_A/N_D$ changes the Fermi levels of electron and holes, respectively. This is due to the fact that the positive charge density in the SWCNT absorber layer is higher than in the window layer. The $N_A$ increase at the junction of the window and absorber layer has no influence on the Fermi levels of
the absorber layer because the increase in the positive charges (holes) is screened by electrons in an accumulation region of the window layer at the space charge region.

In Figure 6, we considered the variation in the electron affinity of both semiconductor layers as well. The electron affinity, $\chi$, is the minimum energy needed to move an electron from the Fermi level into vacuum level. There is a very small increase in $V_{oc}$ with the increased $\chi$ of SWCNT but a slight decrease in $V_{oc}$ of the GaAs. However, in both cases, the performance parameters increase to an optimum value by increasing the $\chi$. We also considered the effect of the bandgap of each layer on the performance parameters. The efficiency and FF decrease after the optimum interval of 1.10 and 1.24 eV for SWCNT and GaAs layer, respectively. In this regard, in Figure 7 the simulated external quantum efficiency (QE) is presented for 1 sun illumination condition.

It is apparent that the drop of efficiency after 500 nm is associated with the fall of the FF. The enhancement in the short wavelength can be due to the improvement of the $J_{sc}$ which is obtained for lower $p$ and higher $n$ concentration levels. This is because the carrier collection will improve by additional electrons. In the SWCNT absorber layer as p-type, the electrons are the minority carriers and the transportation of this minority might be improved by doping to a proper level. Further consideration on the thickness of the layers gives the optimum values for the electrical parameters at the given values of the thickness in Table 1. The cell performance also varies with the operating temperature. For higher temperatures than 300 K, the cell showed lower stability with lower $V_{oc}$ and performance (not shown here). Once we have optimal values for doping concentrations and thickness of the layers (summarized in Table 1), we can simulate the solar cell structure shown in Figure 2. The semiconductors of our heterostructures may contain defect states within the band gap. If the defect states become charged, they add to the space charge of the heterostructures, similar as interface states. Thus, they influence the potential drop in the layers. Certainly, defect states in the band gap also give rise to carrier recombination. This is the reason why the $V_{oc}$ reduces. In the case that there is more than one deep defect level in the band gap, the recombination rates are added up as well as their contributions to the current densities. However, within the three-dimensional absorber (i.e., SWCNT) there may be distinct zones of particularly high defect densities: grain boundaries and regions of high strain, accumulated

Figure 4: The effect of acceptor concentration, ($p$) in the SWCNT layer, on the performance parameters with an optimum value of about $10^{16}$ cm$^{-3}$. 

![Graphs showing the effect of acceptor concentration on performance parameters](image-url)
impurities, or stoichiometry deviations. The performance of the solar cells is influenced by the quality of interfaces and formation of defects such as point defects, stacking faults, twins, dislocations, and grain boundaries. It is important to understand the defect physics so that appropriate methods may be developed to suppress the formation of harmful defects. This is the reason why the performance in Figure 8 is reducing by increased defect concentration.

Further considerations on the critical issues of this structure can be done to attain the high-efficiency heterojunction. For example, the effects of interface properties in the surface and interface defect layer on the performance parameter can be achieved. The simulation investigation along with the practical work on the growth and the preventing of damage to the SWCNT/GaAs wafer needs to be done in order to optimize the cell performance. Further consideration by AMPS tool can be done for the nanostructured solar cells.

For example nanoparticle solar cells can be modelled in this platform by adding a number of quantum dot layers in the active region and assuming the nanoparticles as defects which can be defined in this tool. Note that in these simulations we assumed a defect-free device to simplify the results. However, the more detailed studies can be done elsewhere.

4. Summary

A hybrid heterojunction solar cell based on carbon nanotube/GaAs has been proposed and simulated using the one-dimensional simulation program called Analysis of Micro-electronic and Photonic Structures (AMPS-1D). The SWCNT layer was added as the absorber layer and the window layer was taken from GaAs. The value of the parameters was taken from the experimental results of the papers published about this heterojunction. The results well simulate the band diagram and the effect of the thickness and carrier concentration on the performance parameters of the cell. The effects of acceptor and donor concentration, electron affinity, band gap, and thickness of the layers as well as operating temperature on the cell performance parameters have been investigated with a view to find a more efficient
Figure 6: The effect of electron affinity ($\chi$) on the performance parameters.

Figure 7: Quantum efficiency under one sun standard conditions.

cell. For this heterojunction, the maximum efficiency of 6.55% was achieved for a 300 nm and 3 $\mu$m thickness of nanotube and GaAs layer, which decreases for the other values compared to the optimum values. The results are close to the reality since the parameter values were selected from the practical reported data in the literature. This work opens a road to explore in novel nanostructure solar cell configurations for the simulation of the nanostructured materials with AMPS getting a better view of the cell operation [II].
Conflict of Interests

The authors declare that there is no conflict of interests regarding the publication of this paper.

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