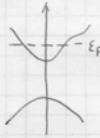
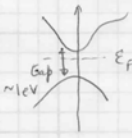
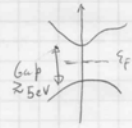


**EE 566 – Solid State Devices
Spring 2006 – Debdeep Jena
Assignment 1 – Solutions**

<< Graphics`

Material Type	Bandstructure/ E_F	Carrier conc.	ρ ($\Omega\text{-cm}$) ^(300K)	ϵ	Growth/Deposition
METALS		$n \sim 10^{23}/\text{cm}^3$	$\sim 10^{-6}$	Very large (dc)	E-beam deposition Electroplate, etc...
SEMICONDUCTORS		$p, n \sim 10^{13}/\text{cm}^3$ depends on doping.	$10^{-2} - 10^{-2}$	~ 10	Chemical, LPE, MBE, MOCVD
INSULATORS		~ 0	$\geq 10^6$	~ 4	Sputtering, epitaxy, etc...

PROBLEM 2

```

q = 1.6 * 10-19      (* Electron charge, Coulomb *)
hbar =  $\frac{6.63}{2 \pi} * 10^{-34}$  (* Reduced Planck's constant, J.s *)
kb = 1.38 * 10-23   (* Boltzmann constant, J/K *)
m0 = 9.1 * 10-31   (* Electron rest mass, Kg *)
meGaN = 0.2 * m0    (* Electron effective mass, CB *)
mhGaN = 1.4 * m0    (* Hole effective mass, VB *)

Nc[T_] = 2 *  $\left( \frac{meGaN * kb * T}{2 * \pi * hbar^2} \right)^{\frac{3}{2}} * 10^{-6}$  (* CB edge Effective DOS, cm-3 *)
Nv[T_] = 2 *  $\left( \frac{mhGaN * kb * T}{2 * \pi * hbar^2} \right)^{\frac{3}{2}} * 10^{-6}$  (* VB edge Effective DOS, cm-3 *)

F[η_] = Abs  $\left[ \frac{2}{\sqrt{\pi}} \int_0^{\infty} \frac{\sqrt{x}}{1 + \text{Exp}[x - \eta]} dx \right]$ 
(* Fermi-Dirac Integral of order j=1/2 *)

```

```

(* In energy scale, Ev is set to zero,
and Ec=Ev+Eg=3.4 eV. Ef is in eV too! *)
Eg = 3.4      (* Bandgap of GaN, eV. Also conduction band edge!*)
Ea = 0.16     (* Acceptor ionization energy, eV *)
Ed = Eg - 0.01 (* Donor activation energy, eV *)
ND = 1016    (* Donor atom volume density, cm-3 *)
NA = 1019    (* Acceptor atom volume density, cm-3 *)

```

```

NDp[Ef_, T_] =  $\frac{ND}{1 + 2 * \text{Exp}\left[\frac{q * (Ef - Ed)}{kb * T}\right]}$  (* Ionized donor density, cm-3 *)
NAm[Ef_, T_] =  $\frac{NA}{1 + 4 * \text{Exp}\left[\frac{q * (Ea - Ef)}{kb * T}\right]}$ 
(* Ionized acceptor density, cm-3 *)
n[Ef_, T_] = Nc[T] * F  $\left[ \frac{q * (Ef - Eg)}{kb * T} \right]$ 
(* Free electron density in cm-3 dependent on Ef *)
p[Ef_, T_] = Nv[T] * F  $\left[ \frac{q * (0 - Ef)}{kb * T} \right]$ 
(* Free hole density in cm-3 dependent on Ef *)

```

```
(*-----SET Directory TO Export Data-----*)
SetDirectory["C:\Documents and Settings\Debdeep Jena\My
  Documents\Transfer\PostPhD\Universities\NotreDame\Teaching\
  CourseMaterial\6_Spring_EE_566_SSD\Assignments\Mathematica"]
FermiLevelPlot = Table[{Ef, NDp[Ef, 300], NAm[Ef, 300],
  n[Ef, 300], p[Ef, 300], NDp[Ef, 300] + p[Ef, 300],
  NAm[Ef, 300] + n[Ef, 300]}, {Ef, -0.2, 3.6, 0.05}]
Export["FermiLevelPlot.txt", FermiLevelPlot, "Table"]
```

```
(* Numerical Solution of the charge-
  neutrality equation gives us the Fermi Level in eV. Note that you
  can do this for any general temperature and doping densities. *)
FindRoot[NDp[Ef, 300] + p[Ef, 300] - (NAm[Ef, 300] + n[Ef, 300]) == 0,
  {Ef, 0}]

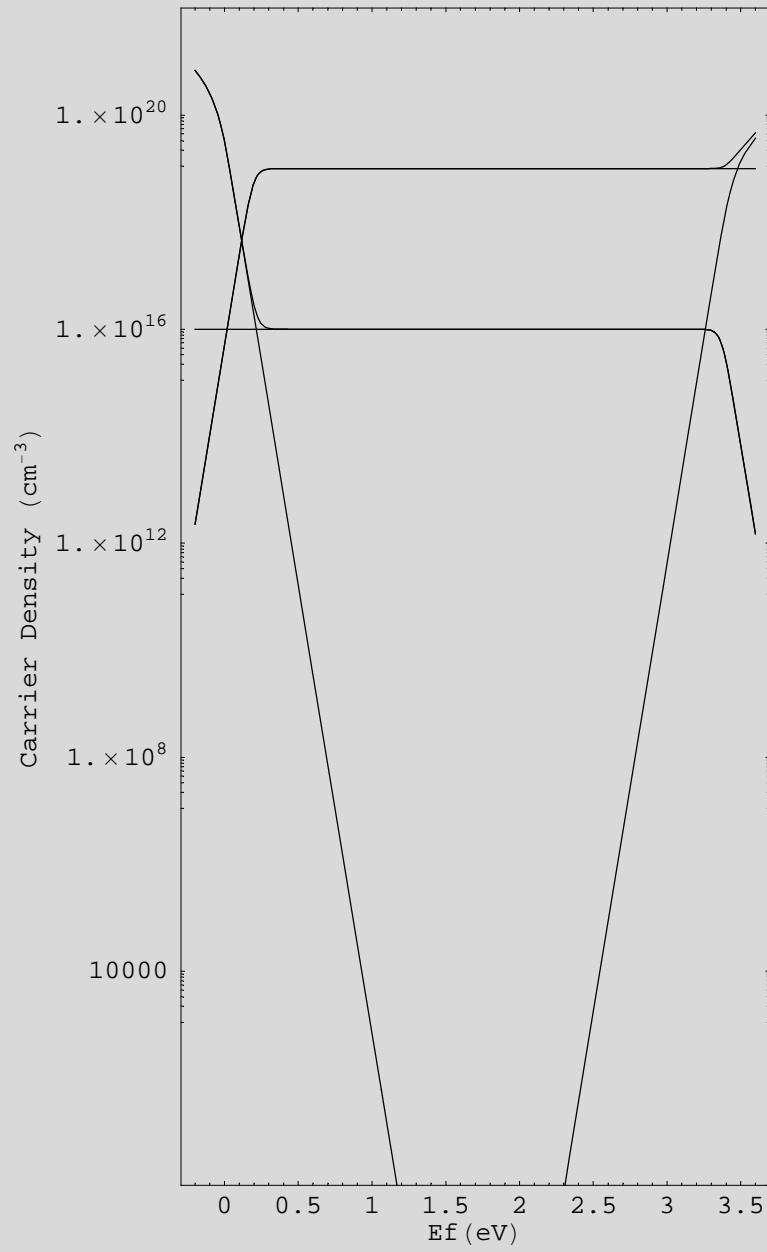
{Ef -> 0.117157}
```

```
root[T_] :=
  FindRoot[NDp[Ef, T] + p[Ef, T] - (NAm[Ef, T] + n[Ef, T]) == 0, {Ef, 0}]
r[T_] := Ef /. root[T]
```

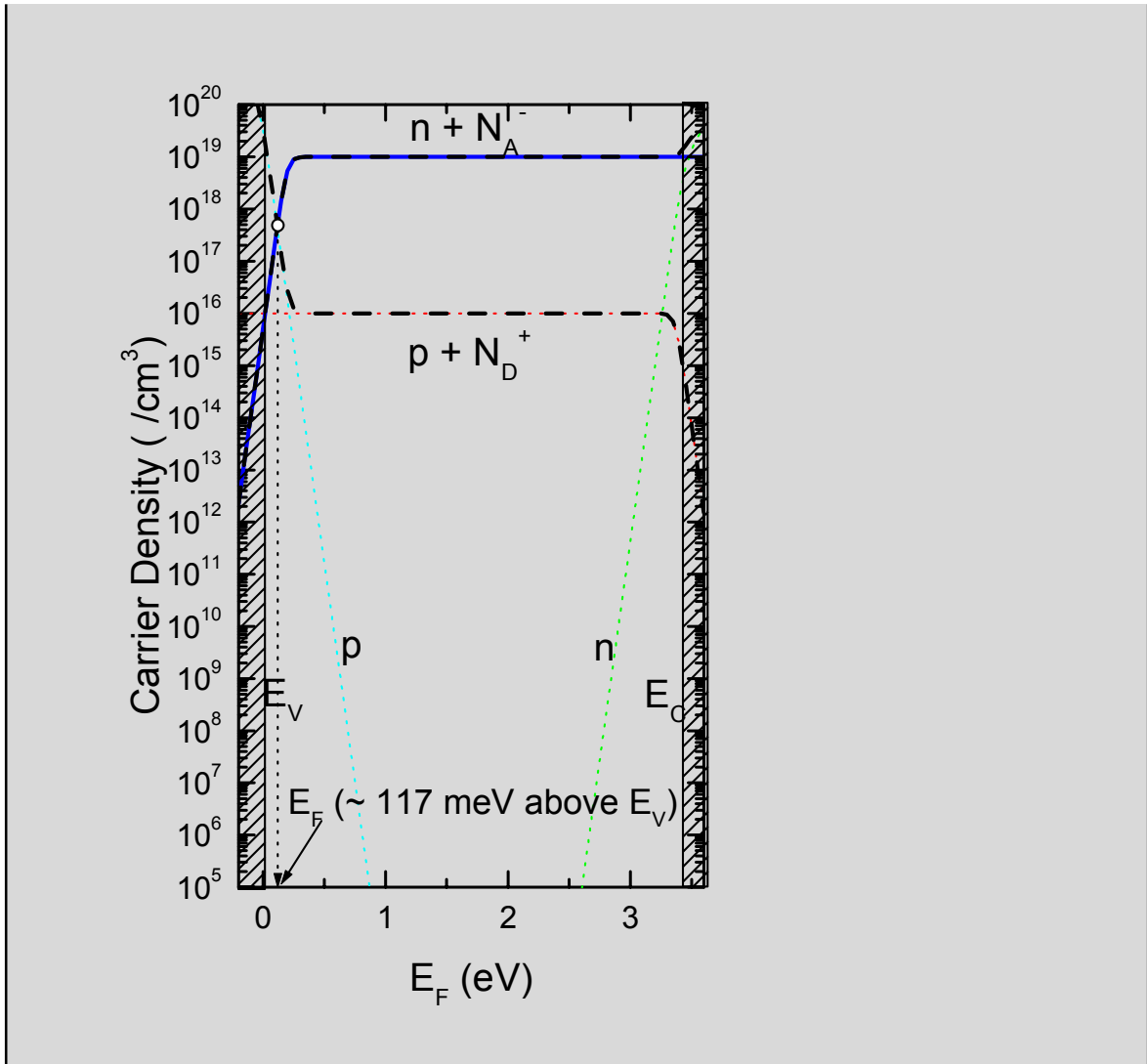
```
p[0.117157, 300]
(* This is the Hole concentration,
  the semiconductor is obviously p-type *)

 $4.45615 \times 10^{17}$ 
```

```
(*----You can do a Graphical Solution also -
  See the attached figures, which are basically the
  plot below labeled and suitably commented upon----*)
LogPlot[{NDp[Ef, 300], NAm[Ef, 300], n[Ef, 300], p[Ef, 300],
  NDp[Ef, 300] + p[Ef, 300], NAm[Ef, 300] + n[Ef, 300]},
  {Ef, -0.2, 3.6}, PlotRange -> {100, 1022}, Frame -> True,
  AspectRatio -> 2, FrameLabel -> {"Ef (eV)", "Carrier Density (cm-3)"}]
```



- Graphics -



PROBLEM 3

Simple problem - use Vegard's law to interpolate the bandgap for the alloy. Investigate what does one mean by the negative bandgap??

PROBLEM 4

Apply Gauss' s law (or Poisson equation) to the junction,
and calculate the charge - field -

band diagram. The charge profile should be neutral,
and thereafter the problem is simple. Note that the electric field profile

determines the profile of the potential $V(x)$. Since $V(x) = - \int_{-\infty}^x F(x) dx$,
the potential changes fastest where the electric field is the highest
in magnitude. Here, the field reaches it' s maximum value at the metal -
semiconductor interface, therefore the slope of $V(x)$ should be the
highest there. The slope of $V(x)$ should decrease continuously from $x =$
0 to $x = x_{d2}$. I use $F(x)$ for electric field in the following

$$F(x) = 0, \quad x < 0$$

$$F(x=0) = - |F_{\max}| = - \frac{[\rho_1 * x_{d1} + 2 * \rho_1 * (x_{d2} - x_{d1})]}{\epsilon_s} = - \frac{\rho_1 (2 x_{d2} - x_{d1})}{\epsilon_s},$$

$$F(x) = F(x=0) + \frac{1}{\epsilon_s} \int_{x=0}^x (\rho_1) dx =$$

$$- \frac{\rho_1 (2 x_{d2} - x_{d1})}{\epsilon_s} + \frac{(\rho_1 * x)}{\epsilon_s} = - \frac{\rho_1 [2 x_{d2} - (x_{d1} + x)]}{\epsilon_s} \quad \text{for } 0 < x < x_{d1},$$

$$F(x) = F(x_{d1}) + \frac{1}{\epsilon_s} \int_{x=x_{d1}}^x (2 \rho_1) dx =$$

$$- \frac{2 \rho_1 * (x_{d2} - x_{d1})}{\epsilon_s} + \frac{2 \rho_1 (x - x_{d1})}{\epsilon_s} = - \frac{2 \rho_1 (x_{d2} - x)}{\epsilon_s} \quad \text{for } x_{d1} < x < x_{d2},$$

$$F(x) = 0, \quad x > x_{d2}$$

Similarly, the potential can be evaluated as $V(x) = - \int_{-\infty}^x F(x) dx$ for each x .

