Funciones de Bessel

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In [2]:

In [3]:

```
# Funciones de Bessel (recursivo)
def Bessel(k,l,r):
    rho = k*r
    if rho==0.0:
        jl = 0
        if l==0:
            jl=1
        return jl
    j0 = \sin(rho) / rho
    j1 = \sin(rho)/(rho**2) - \cos(rho)/rho
    if l==0:
         return j0
    if l==1:
        return j1
    for j in range(1,1):
        jl = (2*l-1)/rho*j1 - j0
        j1 = jl
        j0 = j1
    return jl
```

In [4]:

```
# Funciones de Neumann (recursivo)

def Neumann(k,l,r):
    rho = k*r
    n0 = -cos(rho) / rho
    n1 = -cos(rho)/(rho**2) - sin(rho)/rho

if l==0:
    return n0

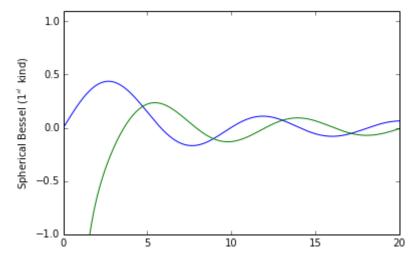
if l==1:
    return n1

for j in range(1,l):
    n1 = (2*l-1)/rho*n1 - n0
    n1 = n1
    n0 = n1

return nl
```

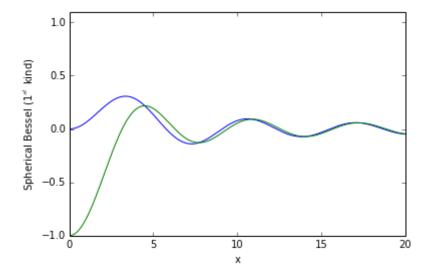
In [45]:

```
# Plot
# Parámetros
step=0.01
xmin=step
xmax=20
x=arange(xmin,xmax,step)
nsize = len(x)
ener= 0.3
k = sqrt(2*ener)
lq=1
jl=zeros(nsize)
nl=zeros(nsize)
for i in range(rx):
  jl[i]= Bessel(k,lq,x[i])
  nl[i]= Neumann(k,lq,x[i])
plot(x,jl[:])
plot(x,nl[:])
ylabel("Spherical Bessel (1$^{st}$ kind)")
axis([xmin,xmax,-1,1.1])
show()
```



In [46]:

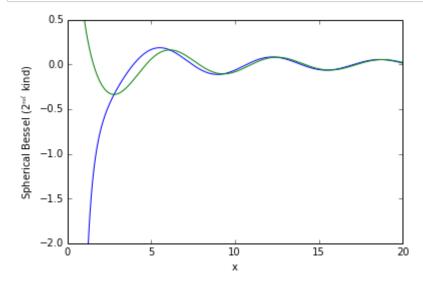
```
# Asymptotic jl
ener= 0.5
k = sqrt(2*ener)
lq=2
jl=zeros(nsize)
jlasy=zeros(nsize)
for i in range(nsize):
  jl[i]= Bessel(k,lq,x[i])
  jlasy[i] = sin(k*x[i] - lq*pi/2)/(k*x[i])
plot(x,jl[:])
plot(x,jlasy[:])
xlabel("x")
ylabel("Spherical Bessel (1$^{st}$ kind)")
axis([xmin,xmax,-1,1.1])
show()
```



In [47]:

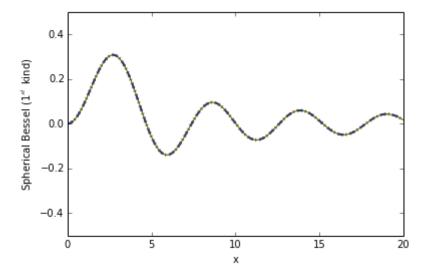
```
# Asymptotic nl
ener= 0.5
k = sqrt(2*ener)
lq=2
nl=zeros(nsize)
nlasy=zeros(nsize)

for i in range(nsize):
    nl[i]= Neumann(k,lq,x[i])
    nlasy[i] = -cos(k*x[i] - lq*pi/2)/(k*x[i])
plot(x,nl[:])
plot(x,nlasy[:])
xlabel("x")
ylabel("Spherical Bessel (2$^{nd}$ kind)")
axis([xmin,xmax,-2,0.5])
show()
```



In [48]:

```
# Comparison with Scipy Special Functions
import scipy.special as spl
import numpy as np
k=1.2294221585
lq=2
jl2=zeros(nsize)
jln=zeros(nsize)
for i in range(nsize):
    jl[i]= Bessel(k,lq,x[i])
    jl2[i] = np.sqrt(np.pi/(2*k*x[i])) * spl.jn(lq+0.5,k*x[i])
    A,B = spl.sph_jn(lq,k*x[i])
    jln[i] = A[lq]
plot(x,jl,color='blue',linestyle='dashed',linewidth='2')
plot(x,jl2,color='red',linestyle='dotted',linewidth='3')
plot(x,jln,color='green',linestyle='solid')
axis([xmin,xmax,-0.5,0.5])
xlabel("x")
ylabel("Spherical Bessel (1$^{st}$ kind)")
show()
```



In [9]:

```
# Solución por Diagonalización
```

In [50]:

```
from numpy import identity

def Laplacian(x):
    h = x[1]-x[0] # assume uniformly spaced points
    n = len(x)
    M = -2*identity(n,'d')
    for i in range(1,n):
        M[i,i-1] = M[i-1,i] = 1
    return M/h**2
```

In [51]:

```
from numpy import sqrt
# Normalización de las funciones
def Normalizate(U,x):
    h = x[1]-x[0] # assume uniformly spaced points
    n = len(x)
    for j in range(0,n):
        suma = 0.0
        for i in range(1,n):
             suma = suma + U[i,j]**2
        suma = suma*h
        rnorm = 1/sqrt(suma)
         print j,' integral (sin normalizar) =',rnorm
#
        Normalization
        rsign = 1
        if U[1,j] < 0:
            rsign = -1
        rnorm = rnorm * rsign
        for i in range(0,n):
            U[i,j] = U[i,j]*rnorm
    return U
```

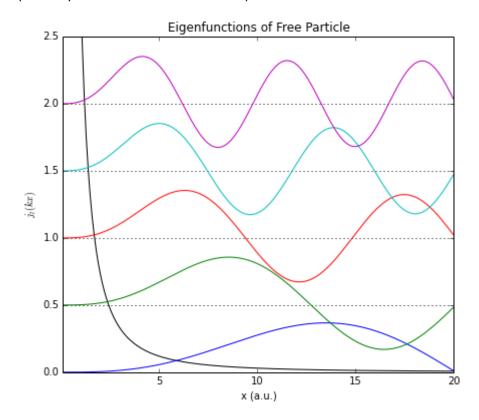
In [57]:

```
from numpy import diag, linspace, array
from numpy.linalg import eigh
from matplotlib.pyplot import axhline, xlabel, ylabel, plot, axis, \
                                figure, title, show
step=0.1
xmin=step
xmax=20
x=arange(xmin,xmax,step)
nsize = len(x)
# array defined above
nsize = len(x)
Dx = step
x = linspace(xmin,xmax,nsize)
T = array([nsize,nsize])
V = array([nsize,nsize])
H = array([nsize,nsize])
E = array([nsize])
# Kinetic (T) and Potential (V)
T = -0.5*Laplacian(x)
lq = 2
V = lq*(lq+1)/(2 * x * x)
# Hamiltonian
H = T + diag(V)
# Eigenvalues (E) and Eigenvectors (U)
E,U = eigh(H)
# Normalization
U=Normalizate(U,x)
```

In [58]:

```
# Plot
#define plot size in inches (width, height) & resolution(DPI)
fig = figure(figsize=(7, 6), dpi=100)
# Number of functions to plot
nfunctions = 5
# Plot the potential
plot(x,V,color='k')
# Plot wavefunctions
ebase=-0.5
for i in range(nfunctions):
    ebase=ebase+0.5
    # For each of the first few solutions, plot the energy level:
    axhline(y=ebase,color='k',ls=":")
    # as well as the eigenfunction, displaced the function
    # so they don't all pile up on each other:
    plot(x,U[:,i]+ebase)
    print("E=:",E[i])
axis([xmin,xmax,0,2.5])
title("Eigenfunctions of Free Particle")
xlabel("x (a.u.)")
ylabel("$j_l(k x)$")
show()
```

```
('E=:', 0.041106602998003935)
('E=:', 0.10235816665720565)
('E=:', 0.18788738170223074)
('E=:', 0.29777290406982221)
('E=:', 0.43201093493202958)
```



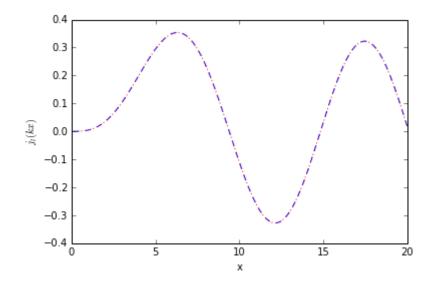
In [59]:

```
# Normalización de vector
def Normld(U,x):
    h = x[1]-x[0] # assume uniformly spaced points
    n = len(x)
    suma = 0.0
    for i in range(0,n):
        suma = suma + U[i]**2
    suma = suma*h
    rnorm = 1/sqrt(suma)
        Normalization
    rsign = 1
    if U[1] < 0:
        rsign = -1
    rnorm = rnorm * rsign
    for i in range(0,n):
        U[i] = U[i]*rnorm
    return U
```

In [64]:

```
# Comparison
nq = 2
ener= E[nq]
k = sqrt(2*ener)
lq=2
print k
jl=zeros(nsize)
for i in range(nsize):
  jl[i]= Bessel(k,lq,x[i]) *k * x[i]
jl = Norm1d(jl,x)
plot(x,jl,'blue',linestyle='dashed')
plot(x,U[:,nq],'red',linestyle='dotted')
xlabel("x")
ylabel("$j l(k x)$")
#axis([xmin,xmax,-0.5,0.5])
show()
```

0.613004700964



In []:

In []: