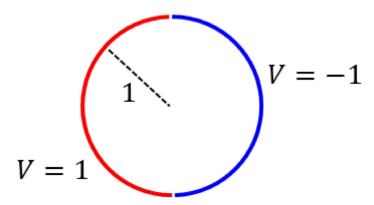
Question 2. By modifying the coefficients in the 2D Poisson solver, get the potential inside the circle of radius 1 in polar coordinates.



In [1]:

```
import numpy as np
import matplotlib.pyplot as plt
```

In [2]:

```
# Get some parameters
N = int(100)  # num. meshes
r=float(1)  # radius
imax=int(1000)  # max step to iterate
tol=float(1e-5)  # error tolerence. something like 1e-5
```

In [3]:

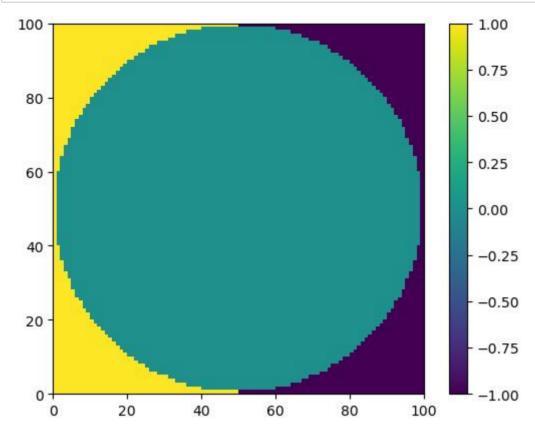
```
# Setup arrays and parameters.
dx=r/N  # mesh size in x
dy=r/N  # mesh size in y
x = np.linspace(-r, r, N)
y = np.linspace(-r, r, N)
X,Y=np.meshgrid(x,y) # Create a 2D mesh grid
u=0.0*X  # u-array, the potential
resid=0.0*X  # residue
```

In [4]:

```
# boundary condition
for i in range(N) :
    for j in range(N) :
        if X[i,j]**2+Y[i,j]**2 >= 1 :
            if X[i,j] < 0 : u[i,j] = 1.0
            elif X[i,j] > 0 : u[i,j] = -1.0
```

In [5]:

```
#plot initial potential
extn=[0,N,0,N]
cs=plt.imshow(u,extent=extn); plt.colorbar(cs); plt.show()
```



In [6]:

```
# Jacobi spectral radius
rjac=(np.cos(np.pi/N) + (dx/dy)**2*np.cos(np.pi/N))/(1+(dx/dy)**2)
```

In [7]:

```
# initial error
errf=sum(sum(np.fabs(u)))
err= errf # initial error
```

In [8]:

```
# Coefficients
a=1.0
b=1.0
c=(dx/dy)**2
d=(dx/dy)**2
e=-2.0-2.0*(dx/dy)**2
```

In [9]:

```
# initial overrelaxation factor omega=1
```

In [10]:

In [11]:

```
# print out run info.
print("Number of iteration=",i)
print("Relative Error=",err/errf)
```

Number of iteration= 98 Relative Error= 1.0

In [12]:

cs=plt.imshow(u,extent=extn); plt.colorbar(cs); plt.show()

