

ProblemSet2-Q2

January 21, 2019

0.0.1 Problem Set #2

MACS 30150, Dr. Evans

Due Monday, Jan. 21 at 11:30am

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2. Numerical integration exercises from Evans: Numerical Integration lab

```
In [1]: import numpy as np
import pandas as pd
from scipy.stats import norm
```

Exercise 2.1

```
In [2]: def g_x(x):
    x = 0.1 * (x ** 4) - 1.5 * x ** 3 + 0.53 * x ** 2 + 2 * x + 1
    return x

def integr_g_x(g, a, b, N, method):
    if method == "midpoint":
        #Calculate vector of N+1 bar bounds
        bin_cuts = np.linspace(a, b, N + 1)

        #Calculate vector of midpoints
        m = (b - a) / (2*N)
        midpoint = np.linspace(a + m, b - m, N)

        #Evaluate the function at the midpoint
        mid_val = g(midpoint)

        #Add up the area of the bins
        approx_integr = (b-a) / N * sum(mid_val)

    if method == "trapezoid":
        #Calculate vector of N+1 bar bounds
        bin_cuts = np.linspace(a, b, N + 1)
```

```

        #Evaluate the function at the bar bounds
        bin_val = g(bin_cuts)

        #Add up the area of the bins
        approx_integr = (b-a) / (2 * N) * (bin_val[0] + 2 * sum(bin_val[1 : N]) + bin_val[-1])

    if method == "Simpsons":
        #Calculate vector of N+1 bar bounds
        bin_cuts = np.linspace(a, b, 2*N + 1)

        #Evaluate the function at the bar bounds
        bin_val = g(bin_cuts)

        #Add up the value at bar bounds
        Sum = bin_val[0] + bin_val[2*N]
        for i in range(1, 2*N):
            if i % 2 == 0:
                Sum += 2 * bin_val[i]
            else:
                Sum += 4 * bin_val[i]

        #Add up the area of the bins
        approx_integr = (b-a) / (6 * N) * Sum

    return approx_integr

```

```

In [3]: Val = 0.02*(10**5-(-10)**5)+0.53/3*(10**3-(-10)**3)+20
        #Midpoint rule
        M = integr_g_x(g_x, -10, 10, 10000, "midpoint")
        print("The approxiamation value using midpoint rule is: ", M)
        print("The difference with true value is: ", M - Val)

The approxiamation value using midpoint rule is:  4373.333196466634
The difference with true value is:  -0.0001368666989947087

```

```

In [4]: #Trapezoid rule
        T = integr_g_x(g_x, -10, 10, 10000, "trapezoid")
        print("The approxiamation value using trapezoid rule is: ", T)
        print("The difference with true value is: ", T - Val)

The approxiamation value using trapezoid rule is:  4373.333607066684
The difference with true value is:  0.00027373335069569293

```

```

In [5]: #Simpson's rule
        S = integr_g_x(g_x, -10, 10, 10000, "Simpsons")
        print("The approxiamation value using Simpson's rule is: ", S)
        print("The difference with true value is: ", S - Val)

```

The approximation value using Simpson's rule is: 4373.33333333337
The difference with true value is: 3.728928277269006e-11

Comparing the values of these integrals to the true analytical value of the integral, we find that Simpson's rule is more accurate.

Exercise 2.2

```
In [6]: def appro_norm(mu, sigma, N, k):
        #Calculate vector of nodes
        z = np.linspace(mu-k*sigma, mu+k*sigma, N)

        #Initialize vector of weights
        weight = np.zeros(N)

        #Calculate vector of weights
        weight[0] = norm.cdf((z[0]+z[1])/2, loc=mu, scale=sigma)
        for i in range(1, N-1):
            weight[i] = integr_g_x(lambda x: norm.pdf(x, loc=mu, scale=sigma), (z[i-1]+z[i])/2, loc=mu, scale=sigma)
        weight[N-1] = 1-norm.cdf((z[N-2]+z[N-1])/2, loc=mu, scale=sigma)
        n = {'nodes' : z, 'weight' : weight}
        df = pd.DataFrame(n)

        return df

df = appro_norm(0, 1, 11, 3)
df
```

```
Out [6]:
```

	nodes	weight
0	-3.0	0.003467
1	-2.4	0.014397
2	-1.8	0.048943
3	-1.2	0.117253
4	-0.6	0.198028
5	0.0	0.235823
6	0.6	0.198028
7	1.2	0.117253
8	1.8	0.048943
9	2.4	0.014397
10	3.0	0.003467

Exercise 2.3

```
In [7]: def appro_log_norm(mu, sigma, N, k):
        #Calculate vector of nodes
        z = np.linspace(mu-k*sigma, mu+k*sigma, N)
        a = np.e**z
```

```

#Initialize vector of weights
weight = np.zeros(N)

#Calculate vector of weights
weight[0] = norm.cdf((z[0]+z[1])/2, loc=mu, scale=sigma)
for i in range(1, N-1):
    weight[i] = integr_g_x(lambda x: norm.pdf(x, loc=mu, scale=sigma), (z[i-1]+z[i])/2, loc=mu, scale=sigma)
weight[N-1] = 1-norm.cdf((z[N-2]+z[N-1])/2, loc=mu, scale=sigma)
n = {'node' : a, 'weight' : weight}
df = pd.DataFrame(n)
approxiamation = sum(df["node"]*df["weight"])
return approxiamation, df

appro, df = appro_log_norm(0, 1, 11, 3)
appro

```

Out [7]: 1.6639409483466545

In [8]: df

```

Out [8]:
   node  weight
0  0.049787  0.003467
1  0.090718  0.014397
2  0.165299  0.048943
3  0.301194  0.117253
4  0.548812  0.198028
5  1.000000  0.235823
6  1.822119  0.198028
7  3.320117  0.117253
8  6.049647  0.048943
9  11.023176  0.014397
10 20.085537  0.003467

```

Exercise 2.4

```

In [10]: appro, df = appro_log_norm(10.5, 0.8, 1000, 8)
EI = np.e**(10.5+0.5*(0.8**2))
difference = abs(appro - EI)

print("approximation of the expected value of income: ", appro)
print("expected value of income: ", EI)
print("difference between these two values: ", difference)

approximation of the expected value of income: 50011.429102086586
expected value of income: 50011.08700852173
difference between these two values: 0.34209356485371245

```

Exercise 3.1

```
In [13]: from scipy.special.orthogonal import p_roots
```

```
def gauss(f,a,b,n):
    # Find optimal weights and nodes using p_roots
    [node,weight] = p_roots(n)
    # Compute weight*node and convert the limits of [a, b] to [-1,1]
    G=0.5*(b-a)*sum(weight*f(0.5*(b-a)*node+0.5*(b+a)))
    return G

Gauss = gauss(lambda x: 0.1*x**4-1.5*x**3+0.53*x**2+2*x+1, -10, 10, 3)
Newton = integr_g_x(lambda x: 0.1*x**4-1.5*x**3+0.53*x**2+2*x+1, -10, 10, 10000, "Simple")
Val = 0.02*(10**5-(-10)**5)+0.53/3*(10**3-(-10)**3)+20
print("The result of Gaussian approximation is", Gauss, 'and the absolute error is', Gauss-Val)
print("The result of Newton-Cotes approximation is", Newton, 'and the absolute error is', Newton-Gauss)

#Reference : https://stackoverflow.com/questions/27115917/gauss-legendre-quadrature-is
```

The result of Gaussian approximation is 4373.333333333335 and the absolute error is 1.818989403

The result of Newton-Cotes approximation is 4373.333333333337 and the absolute error is 3.72892

Compared with the results in Exercise 2.1, Gaussian approximate is more accurate.

Exercise 3.2

```
In [14]: from scipy import integrate
```

```
GQ, err = integrate.quad(lambda x: 0.1*x**4-1.5*x**3+0.53*x**2+2*x+1, -10, 10)

print("The result of Python Gaussian approximate is", GQ, 'and the absolute error is', err)
```

The result of Python Gaussian approximate is 4373.333333333334 and the absolute error is 9.094

Exercise 4.1

```
In [15]: def circle(x, y):
    if x ** 2 + y ** 2 <= 1:
        return 1
    else:
        return 0

def MC(fn, omega, N):
    counter = 0
    xran = np.random.uniform(omega[0], omega[1], N)
    yran = np.random.uniform(omega[2], omega[3], N)
```

```

    for i in range(N):
        x = xran[i]
        y = yran[i]
        counter += fn(x, y)
    area = (omega[1] - omega[0]) * (omega[3] - omega[2])
    return area*counter/N

np.random.seed(seed=25)
N = 1
while round(MC(circle, [-1, 1, -1, 1], N), 4) != 3.1415:
    N += 1

print("The smallest number of random draws N is", N)

```

The smallest number of random draws N is 615

Exercise 4.2

```

In [16]: def isPrime(n):
    '''
    -----
    This function returns a boolean indicating whether an integer n is a
    prime number
    -----
    INPUTS:
    n = scalar, any scalar value

    OTHER FUNCTIONS AND FILES CALLED BY THIS FUNCTION: None

    OBJECTS CREATED WITHIN FUNCTION:
    i = integer in [2, sqrt(n)]

    FILES CREATED BY THIS FUNCTION: None

    RETURN: boolean
    -----
    '''
    for i in range(2, int(np.sqrt(n) + 1)):
        if n % i == 0:
            return False

    return True

In [17]: def primes_ascend(N, min_val=2):
    '''
    -----
    This function generates an ordered sequence of N consecutive prime

```

numbers, the smallest of which is greater than or equal to 1 using the Sieve of Eratosthenes algorithm.
(https://en.wikipedia.org/wiki/Sieve_of_Eratosthenes)

INPUTS:

N = integer, number of elements in sequence of consecutive prime numbers
min_val = scalar ≥ 2 , the smallest prime number in the consecutive sequence must be greater-than-or-equal-to this value

OTHER FUNCTIONS AND FILES CALLED BY THIS FUNCTION:

isPrime()

OBJECTS CREATED WITHIN FUNCTION:

primes_vec = (*N*,) vector, consecutive prime numbers greater than *min_val*
MinIsEven = boolean, =True if *min_val* is even, =False otherwise
MinIsGrtrThn2 = boolean, =True if *min_val* is greater-than-or-equal-to 2, =False otherwise
curr_prime_ind = integer ≥ 0 , running count of prime numbers found

FILES CREATED BY THIS FUNCTION: None

RETURN: *primes_vec*

```
'''
primes_vec = np.zeros(N, dtype=int)
MinIsEven = 1 - min_val % 2
MinIsGrtrThn2 = min_val > 2
curr_prime_ind = 0
if not MinIsGrtrThn2:
    i = 2
    curr_prime_ind += 1
    primes_vec[0] = i
i = min(3, min_val + (MinIsEven * 1))
while curr_prime_ind < N:
    if isPrime(i):
        curr_prime_ind += 1
        primes_vec[curr_prime_ind - 1] = i
    i += 2

return primes_vec
```

```
In [18]: def equidis(n,d,Type):
prime = primes_ascend(d)
if Type == 'Weyl':
    return list((i - i // 1) for i in np.sqrt(prime)*n)
elif Type == 'Haber':
```

```

        return list((i - i // 1) for i in np.sqrt(prime)*n*(n+1)/2)
    elif Type == 'Niederreiter':
        return list((i - i // 1) for i in [n*(2**(j/(n+1))) for j in range(1, d+1)])
    elif Type == 'Baker':
        return list((i - i // 1) for i in [n*np.exp(1/j) for j in range(1, d+1)])

```

In [19]: `equidis(10, 4, 'Weyl')`

Out[19]: [0.142135623730951,
0.32050807568877104,
0.36067977499789805,
0.4575131106459054]

In [20]: `equidis(10, 4, 'Haber')`

Out[20]: [0.7817459305202306,
0.2627944162882443,
0.9837387624884428,
0.5163221085524867]

In [21]: `equidis(10, 4, 'Niederreiter')`

Out[21]: [0.6504108943996272,
0.3431252219546259,
0.08089444404447121,
0.8666489800943182]

In [22]: `equidis(10, 4, 'Baker')`

Out[22]: [0.18281828459045002,
0.4872127070012837,
0.9561242508608956,
0.8402541668774148]

Exercise 4.3

```

In [23]: def g(x,y):
            if x**2+y**2<1:
                return 1
            else:
                return 0

np.random.seed(25)
def QMC(fn, omega, N, Type):
    x1, x2=omega[0],omega[1]
    y1, y2=omega[2],omega[3]
    if Type=='Weyl':
        xran=[2*np.array(equidis(i,2,'Weyl'))[0]-1 for i in range(N)]
        yran=[2*np.array(equidis(i,2,'Weyl'))[1]-1 for i in range(N)]

```



```

elif Type=='Haber':
    xran=[2*np.array(equidis(i,2,'Haber'))[0]-1 for i in range(N)]
    yran=[2*np.array(equidis(i,2,'Haber'))[1]-1 for i in range(N)]
elif Type=='Niederreiter':
    xran=[2*np.array(equidis(i,2,'Niederreiter'))[0]-1 for i in range(N)]
    yran=[2*np.array(equidis(i,2,'Niederreiter'))[1]-1 for i in range(N)]
elif Type=='Baker':
    xran=[2*np.array(equidis(i,2,'Baker'))[0]-1 for i in range(N)]
    yran=[2*np.array(equidis(i,2,'Baker'))[1]-1 for i in range(N)]
counter = 0
for i in range(N):
    x = xran[i]
    y = yran[i]
    counter += fn(x, y)
area = (omega[1] - omega[0])*(omega[3] - omega[2])
return area*counter/N

```

```

In [24]: def f(x, y):
        if x ** 2 + y ** 2 <= 1:
            return 1
        else:
            return 0

```

```

N = 1
while round(QMC(f, [-1, 1, -1, 1], N, "Weyl"),4) != 3.1415:
    N += 1
print("The smallest number of random draws N with Weyl sequence is", N)

```

The smallest number of random draws N with Weyl sequence is 1230

```

In [25]: N = 1
        while round(QMC(f, [-1, 1, -1, 1], N, "Haber"),4) != 3.1415:
            N += 1

        print("The smallest number of random draws N with Haber sequence is", N)

```

The smallest number of random draws N with Haber sequence is 2064

```

In [27]: N = 1
        while round(QMC(f, [-1, 1, -1, 1], N, "Niederreiter"),4) != 3.1415:
            N += 1
            if N > 5000:
                print("The smallest number of random draws N with Niederreiter sequence exceed")
                break
        else:
            print("The smallest number of random draws N with Niederreiter sequence is", N)

```

The smallest number of random draws N with Niederreiter sequence exceeds 5000.

```
In [29]: N = 1
         while round(QMC(f, [-1, 1, -1, 1], N, "Baker"),4) != 3.1415:
             N += 1

         print("The smallest number of random draws N with Baker sequence is", N)
```

The smallest number of random draws N with Baker sequence is 205

From the above results, we can see that Baker sequence is the fastest one to converge to 3.1415, Weyl sequence is the second one, Haber sequence is the third one and Niederreiter sequence is the slowest one.

```
In [ ]:
```