

Student Names and IDs:

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Homework 5

Part 1: Numerical Differentiation

Problem 1.1

Write a formula for the Jacobian J_{P2C} of this transformation.

Answer

$$J_{P2C} = \begin{bmatrix} \cos\phi & -r\sin\phi \\ \sin\phi & r\cos\phi \end{bmatrix}$$

Problem 1.2

Write functions with headers

```
def P2C(z):
```

and

```
def JacobianP2C(z):
```

that compute the transformation above and its Jacobian matrix. Inputs and outputs should be numpy arrays.

Show your code. You will test it in a later problem.

Answer

```
In [647]: import numpy as np
```

```
In [648]: def P2C(z):
    '''
    This function transforms from polar to cartesian. Concretely, from r
    and phi to x and y
    Inputs: z, a numpy array of r and phi
    Outputs: c, a numpy array of x and y
    '''
    r = z[0]
    phi = z[1]
    a = np.zeros(z.shape)
    a[0] = r*np.cos(phi)
    a[1] = r*np.sin(phi)

    return a
```

```
In [649]: def JacobianP2C(z):
    '''
    This function returns the Jacobian of a transformation from polar to
    cartesian
    Input: numpy arrays, z = (2,)
    Output: numpy arrays
    '''
    r = z[0]
    phi = z[1]
    a = np.zeros((2,2))
    a[0][0] = np.cos(phi)
    a[0][1] = -r*np.sin(phi)
    a[1][0] = np.sin(phi)
    a[1][1] = r*np.cos(phi)
    return a
```

Problem 1.3

Write a Python function with header

```
def Jacobian(f, z, delta=1e-5):
```

that takes a function f from \mathbb{R}^d to \mathbb{R}^e , a numpy vector z with d entries, and an optional value for δ and returns a numpy array with the Jacobian of f at z , using the central difference formula given above.

Show your code. You will test it in the next problem.

Answer

```
In [650]: def Jacobian(f, z, delta=1e-5):
          '''
          Input: f, function
          z, numpy vector with d entries
          delta, optional value for small distance
          Output: numpy array of Jacobian of f at z using central difference t
          hat is (e,d)
          '''
          e = f(z).size          # length of f(z)
          d = z.size             # length of z
          j = np.zeros((e, d))   # (e,d)
          eye = np.identity(d)   # (d,d) identity matrix

          for i in range(d):
              e_i = eye[i]
              df = (f(z + e_i*delta) - f(z - e_i*delta)) / (2*delta) # e
              j[:, i] = df

          return j
```

Problem 1.4

Show the result of running the tests below. This will happen automatically once your functions `Jacobian`, `P2C`, and `JacobianP2C`, are defined correctly (and you run the cell below).

Answer

```
In [651]: def compare(a, f, b, delta=1e-5):
    def a2s(a):
        def n2s(x):
            return '{:g}'.format(round(x, 4))

        try:
            return '[' + ';' .join([' , '.join([n2s(y) for y in row]) for
row in a]) + ']'
        except TypeError:
            try:
                return '[' + ', '.join([n2s(y) for y in a]) + ']'
            except TypeError:
                return '[]' if a.size == 0 else n2s(a)

    aName, fName, bName = a.__name__, f.__name__, b.__name__
    msgBase = '{:s}({:s}, {{:s}}) = {{:s}}\n{:s}({:s}) = {{:s}}'
    msg = msgBase.format(aName, fName, bName)
    zs = np.array([[0, 0], [1, 0], [2, 1], [2, 2]])
    for z in zs:
        print(msg.format(a2s(z), a2s(a(f, z, delta)), a2s(z), a2s(b(z
))), end='\n\n')

    try:
        compare(Jacobian, P2C, JacobianP2C)
    except NameError:
        pass
```

```
Jacobian(P2C, [0, 0]) = [1, 0; 0, 0]
JacobianP2C([0, 0]) = [1, 0; 0, 0]
```

```
Jacobian(P2C, [1, 0]) = [1, 0; 0, 1]
JacobianP2C([1, 0]) = [1, -0; 0, 1]
```

```
Jacobian(P2C, [2, 1]) = [0.5403, -1.6829; 0.8415, 1.0806]
JacobianP2C([2, 1]) = [0.5403, -1.6829; 0.8415, 1.0806]
```

```
Jacobian(P2C, [2, 2]) = [-0.4161, -1.8186; 0.9093, -0.8323]
JacobianP2C([2, 2]) = [-0.4161, -1.8186; 0.9093, -0.8323]
```

Problem 1.5

Use the fact that the Hessian is the Jacobian of the gradient to write a Python function with header

```
def Hessian(f, x, delta=1e-5):
```

that uses your gradient function to compute the Hessian of f at x . Show your code.

Answer

```
In [652]: def Hessian(f, x, delta=1e-5):  
    d = x.size  
    H = np.zeros((d, d))          # (d,d)  
    J = Jacobian(f, x, delta=1e-5) # (1,d) with first derivatives  
    J = J.reshape((d,))          # reshape so that we can reuse J  
    H = Jacobian(lambda x: Jacobian(f, x, delta=1e-5), x, delta=1e-5)  
    return H
```

Problem 1.6

Show the result of running the tests below. This will happen automatically once your function `Hessian` is defined correctly (and you run the cell below).

Answer

```
In [653]: shift, scale = [2, 1], 10
```

```
def f(z):  
    d = z - shift  
    return np.array(np.inner(z, z) / scale + np.exp(-np.inner(d, d)))  
  
def gradientF(z):  
    d = z - shift  
    return 2 * (z / scale - d * np.exp(-np.inner(d, d)))  
  
def HessianF(z):  
    I = np.eye(2)  
    d = z - shift  
    return 2 * (I / scale + (2 * np.outer(d, d) - I) * np.exp(-np.inner(d, d)))  
  
try:  
    compare(Jacobian, f, gradientF)  
    compare(Hessian, f, HessianF)  
except NameError:  
    pass
```

```
Jacobian(f, [0, 0]) = [0.027, 0.0135]  
gradientF([0, 0]) = [0.027, 0.0135]
```

```
Jacobian(f, [1, 0]) = [0.4707, 0.2707]  
gradientF([1, 0]) = [0.4707, 0.2707]
```

```
Jacobian(f, [2, 1]) = [0.4, 0.2]  
gradientF([2, 1]) = [0.4, 0.2]
```

```
Jacobian(f, [2, 2]) = [0.4, -0.3358]  
gradientF([2, 2]) = [0.4, -0.3358]
```

```
Hessian(f, [0, 0]) = [0.2943, 0.0539; 0.0539, 0.2135]  
HessianF([0, 0]) = [0.2943, 0.0539; 0.0539, 0.2135]
```

```
Hessian(f, [1, 0]) = [0.4707, 0.5413; 0.5413, 0.4707]  
HessianF([1, 0]) = [0.4707, 0.5413; 0.5413, 0.4707]
```

```
Hessian(f, [2, 1]) = [-1.8, 0; 0, -1.8]  
HessianF([2, 1]) = [-1.8, 0; 0, -1.8]
```

```
Hessian(f, [2, 2]) = [-0.5358, -0; -0, 0.9358]  
HessianF([2, 2]) = [-0.5358, 0; 0, 0.9358]
```

Problem 1.7

Write one clear and concise sentence to describe which results are good and which are not in the tests below.

```

In [654]: try:
            delta = 1e-9
            compare(Jacobian, f, gradientF, delta)
            compare(Hessian, f, HessianF, delta)
        except NameError:
            pass

Jacobian(f, [0, 0]) = [0.027, 0.0135]
gradientF([0, 0]) = [0.027, 0.0135]

Jacobian(f, [1, 0]) = [0.4707, 0.2707]
gradientF([1, 0]) = [0.4707, 0.2707]

Jacobian(f, [2, 1]) = [0.4, 0.2]
gradientF([2, 1]) = [0.4, 0.2]

Jacobian(f, [2, 2]) = [0.4, -0.3358]
gradientF([2, 2]) = [0.4, -0.3358]

Hessian(f, [0, 0]) = [0.2943, 0.0539; 0.0539, 0.2135]
HessianF([0, 0]) = [0.2943, 0.0539; 0.0539, 0.2135]

Hessian(f, [1, 0]) = [0.4707, 0.5413; 0.5413, 0.4707]
HessianF([1, 0]) = [0.4707, 0.5413; 0.5413, 0.4707]

Hessian(f, [2, 1]) = [-1.8, 0; 0, -1.8]
HessianF([2, 1]) = [-1.8, 0; 0, -1.8]

Hessian(f, [2, 2]) = [-0.5358, -0; -0, 0.9358]
HessianF([2, 2]) = [-0.5358, 0; 0, 0.9358]

```

Answer

All of the results are good.

Part 2: Steepest Descent

Problem 2.1

Using the imports and definition in the cell below, write a function with header

```
def lineSearch(f, g, z0):
```

that performs line search on the function f , whose gradient is computed by the function g , starting at point z_0 . If the starting point z_0 is in \mathbb{R}^d , then functions f and g have the following signatures:

$$f : \mathbb{R}^d \rightarrow \mathbb{R} \quad \text{and} \quad g : \mathbb{R}^d \rightarrow \mathbb{R}^d$$

Show your code, and the result of running the function with the function f and value z_0 defined below. Defining the corresponding gradient g is your task.

```
In [655]: from scipy import optimize as opt
import numpy as np
import math
import matplotlib.pyplot as plt

small = math.sqrt(np.finfo(float).eps)

f, z0 = lambda z: -np.sin(z), 0
```

Answer

```
In [656]: # gradient function
def g(z):
    return -np.cos(z)
```



```
In [657]: def lineSearch(f, g, z0):
'''
    This function performs line search on the function f, whose gradient
    is computed by the function g, starting at point z0.
    a_k, the distance we travel on the line h(a_k)
    z_k+1, the input of the minimum of f(z) found using line search (the
    next step to take)
    Outputs: z, the next step of steepest descent
'''
mag_g = np.linalg.norm(g(z0)) # magnitude of gradient of f at z0
if mag_g < small:
    return z0
def hi(c): # function h from R to R
    return f(z0 + c*(-g_i))

# Find the third element c of the bracketing triplet
g_i = g(z0)
c = small # starting value for c
h_prev = f(z0 + c*(-g_i)) # initial h(c)
c *= 1.2
h = f(z0 + c*(-g_i)) # next h(c)
while h < h_prev:
    c *= 1.2
    h_prev = h
    h = f(z0 + c*(-g_i))
    if c > 100:
        print("Error: c > 100")
        return z0
# now we should have a c in which h_prev(c) < h(c)
a = 0
bracket = (a,c)
b = opt.minimize_scalar(hi, bounds = bracket, method = 'bounded')
z_next = z0 - (b.x * g_i)

return z_next
```

```
In [658]: answer = lineSearch(f, g, z0)
print("Running lineSearch with f and z0 returns a value of z =",answer)
```

Running lineSearch with f and z0 returns a value of z = 1.5707961264726962

Problem 2.2

Write a function with header

```
def steepest(f, g, z0, maxK=10000, delta=small, remember=False):
```

that uses your `lineSearch` function to implement steepest descent with line search.

Show your code and the value of \mathbf{z} that results from minimizing the provided function Rosenbrock with `steepest`. Start the minimization at $\mathbf{z}_0 = (-1.2, 1)$ (defined below), and use the default values for the keyword arguments.

```
In [659]: def Rosenbrock(z):  
           return 100 * (z[1] - z[0] ** 2) ** 2 + (1 - z[0]) ** 2  
  
           def RosenGrad(z):  
               return np.array([400 * (z[0] ** 3 - z[0] * z[1]) + 2 * z[0] - 2,  
                                200 * (z[1] - z[0] ** 2)])  
  
           z0, zStar = np.array([-1.2, 1]), np.array([1, 1])
```

Answer

```
In [660]: def steepest(f, g, z0, maxK=10000, delta=small, remember=False):
    '''
    uses lineSearch function to implement steepest descent with line search
    '''
    count = 0
    steps = []
    steps.append(z0)
    z_prev = lineSearch(f, g, z0) # this is scalar, but I want my line search to return a R^d vector z
    steps.append(z_prev)
    count += 1
    z_current = lineSearch(f, g, z_prev)
    count += 1
    steps.append(z_current)

    while True:
        count += 1
        if np.linalg.norm(z_current - z_prev) < delta or count > maxK:
            break
        z_prev = z_current
        z_current = lineSearch(f, g, z_prev)
        steps.append(z_current)
    m = len(steps)
    d = z0.size
    history = np.zeros((m,d))

    for i in range(m):
        history[i,:] = steps[i]
    if remember == True:
        return (z_current, history)
    elif remember == False:
        return z_current
```

```
In [661]: answer = steepest(Rosenbrock, RosenGrad, z0, maxK=10000, delta=small, remember=False)
print("Running steepest with Rosenbrock and z0 returns a value of z =", answer)
```

```
Running steepest with Rosenbrock and z0 returns a value of z = [0.99995379 0.99990749]
```

Problem 2.3

Now run steepest as follows (the try/except is there so that the notebook will still run if steepest is undefined).

```
In [662]: try:
          [zHat, history] = steepest(Rosenbrock, RosenGrad, z0, maxK=1000, rem
          ember=True)
except NameError:
          pass
```

Make a contour plot of the Rosenbrock function using 10 levels drawn as thin gray lines (use `colors='grey'`, `linewidths=1` in your call to `matplotlib.pyplot.contour` to this effect) for $-1.5 \leq z_1 \leq 1.5$ and $-0.5 \leq z_2 \leq 1.5$. To make the contour plot, sample this rectangle of values with 100 samples in each dimension.

Superimpose a plot of the path recorded in `history` on the contour plot. Also draw a blue dot at \mathbf{z}_0 and a red dot at \mathbf{z}^* , the true minimum.

Show code and plot.

Answer

```
In [663]: x = np.linspace(-1.5, 1.5, 100)
          y = np.linspace(-0.5, 1.5, 100)
          X, Y = np.meshgrid(x,y)
          Z = np.zeros((100,100))
          Z[0][0] = 1
          for i in range(100):
              for j in range(100):
                  Z[i][j] = Rosenbrock(np.array([x[i], y[j]]))
          x_hist = history[:,0]
          y_hist = history[:,1]
```

```
In [664]: plt.contour(grid[0,:,:), grid[1,:,:), Z, 10, colors='grey', linewidths=1)
plt.plot(z0[0], z0[1], 'bo')
plt.plot(zStar[0], zStar[1], 'ro')
plt.plot(x_hist, y_hist)
plt.xlabel("z1")
plt.ylabel("z2")
plt.title("Steepest descent for m=1000")
```

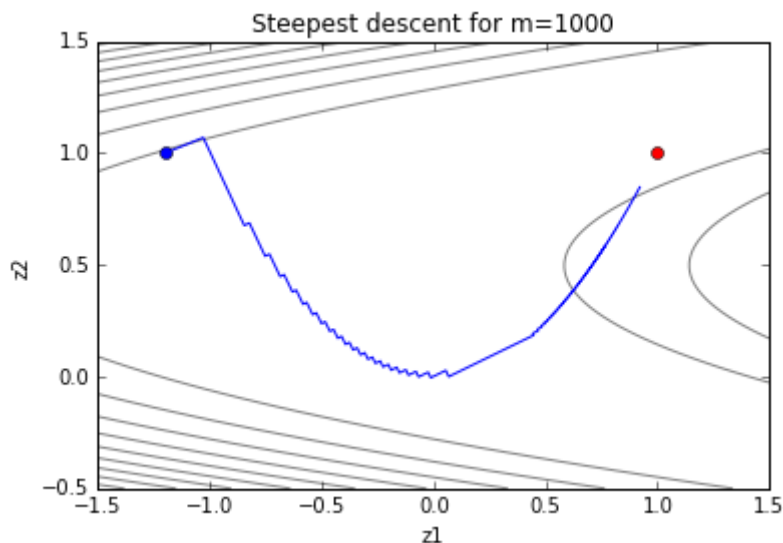
/anaconda3/lib/python3.6/site-packages/numpy/ma/core.py:6434: MaskedArrayFutureWarning: In the future the default for ma.minimum.reduce will be axis=0, not the current None, to match np.minimum.reduce. Explicitly pass 0 or None to silence this warning.

```
return self.reduce(a)
```

/anaconda3/lib/python3.6/site-packages/numpy/ma/core.py:6434: MaskedArrayFutureWarning: In the future the default for ma.maximum.reduce will be axis=0, not the current None, to match np.maximum.reduce. Explicitly pass 0 or None to silence this warning.

```
return self.reduce(a)
```

```
Out[664]: <matplotlib.text.Text at 0x11960fef0>
```



Problem 2.4

Convergence slows down as \mathbf{z}^* is approached, and even after 1000 iterations there is still a way to go. To see this slowdown more clearly, plot a vector distance with the Euclidean distances between each of the points in history (obtained in the previous problem) and \mathbf{z}^* . Label the plot axes meaningfully.

Show code and plot.

Answer ¶

```
In [665]: distances = []  
         for i in range(len(history)):  
             d = np.linalg.norm(history[i] - zStar)  
             distances.append(d)
```

```
In [666]: plt.plot(range(len(history)), distances)  
         plt.xlabel("mth step")  
         plt.ylabel("distance between z_m and z*")  
         plt.title("Decreasing step sizes in steepest descent")
```

Out[666]: <matplotlib.text.Text at 0x119595080>

