Convex Optimization Homework #03

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Problem 1:

Given, f ∈ CL(pn)

Then, & 2, d & comf

 $f(y) \leq f(x) + \nabla f(x)^{T}(y-x) + \frac{1}{2} ||y-x||_{2}^{2}$ Let, $y = \chi(x) + \frac{1}{2} ||y-x||_{2}^{2}$ and $\chi = \chi(x)$

Then,

$$f(x^{k} + t^{(k)} \Delta x^{(k)}) \leq f(x^{(k)}) + \nabla f(x^{(k)}) (x^{(k)} + t^{(k)}) + \frac{1}{2} \|x^{(k)} + t^{(k)} \Delta x^{(k)} - x^{(k)} \|^{2}$$

Pant 1:

For DX(w) to be a descent direction, we need to have,

from (1)

$$t^{(k)} \nabla f(x^{(k)})^T \Delta x^{(k)} + t^{(k)} \frac{1}{2} ||\Delta x^{(k)}||_2^2$$

Now in these two terrors,

$$\frac{S_0}{\text{for } 2^{+0}} \Rightarrow \forall f(x^{(u)})^{\top} \Delta x^{(u)} < 0$$

$$\frac{S_0}{\text{for } 2^{+0}} \Rightarrow \forall f(x^{(u)})^{\top} \Delta x^{(u)} < 0 ; [t^{(u)}, 0]$$

so, $Dx^{(k)}$ will be a of descent direction if and only if it makes are acute andle with $-\nabla \dot{f}(x^{(k)})$

But we also have to choose the carefully so that the positive term in ex. 2;

the Land overweigh the regative term; the positive term; the positive term is the carefully the regative term; the Theory of the regative term; the Theory of the positive term;

Part: 2

From ex 2 +(h) x +(x(h)) Dx(h) + +(w) - 1 1 2 1 Dx(h) 1 2 <0

=>
$$t^{(k)} < \frac{-\nabla f(x^{(k)})}{\frac{L}{2}} || \Delta x^{(k)} ||_{2}^{2}$$

50,

$$0 < \frac{(k)}{L} < -\frac{2}{L} \frac{\Delta f(x(w))}{\|\Delta x(w)\|_{2}^{2}}$$

Prodem 2:

from chain rule,

$$\begin{cases} J(x) & \text{where } x \in \mathbb{R}^n \\ D(x) & \text{where } x \in \mathbb{R}^n \end{cases}$$

for given function,

- Vous Vill be lipschitz continous iff 11 2 for - 2 for 1, < [12 21] > A 2 Egod 1 17 Jay - x Jay 1 = 114 11x 12 - x - 4114 12 7 1/2 =411 112/2 - 11/11/2 112. How, 11x112 - 62/1211 + 62/1211 - 2 2/1211 = R 2/1211 - 1/2/11x11 = 11x112(2-4) + (1|x112-11x11=) Using triansular inequality [11 at 611 < 11 all +11611

(because, ||cv|| = |c|. ||v|| Now, inverse triangular inequality 11a-611 > 11a11 -11611 112/12 - 111/2 | 111/2 < 11x - 7/1 (112/1/2 + 111/2) 117/1/2 comphing 3,965,

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= 112-4112 (112112 + (112112 + 112112) 114112)

Now, a, y & domf.

80, 112112 < 1

11711, ~ < 1

112112 + 114112 62

€ 3 112-311₂

From (1)

117 f(w) - 7 f(v)||2 < 12 ||2-11|2

So, \$1(a) fullfills the condition of lipschitz continous and Companing this to condition for lipschitz continous,

L = 12

Problem 3:

Strategy: I have completed the code in these five steps explained below.

Step_1: Defining the gradient descent function

At first, I have defined the gradient descent function with the inputs mentioned in the question, namely:

- -A function computing the gradient of the objective.
- An initialization point x(0)
- A flag specifying whether to use a fixed step size or a variable step size.
- A step size value (for fixed step size).
- A maximum number of iterations.
- A tolerance ϵ for the stopping criterion.

At first, two things are checked in the function,

- 1. Positive step size: Gradient descent algorithm is defined with a positive step size. Zero step size will not make any progress toward the minimum, and it will waste computational resources.
- 2. Step size type: Step size is kept fixed for this problem.

So if positive and fixed step size is not maintained, the function will not work.

Then, until we reach the max number of iterations, as defined in the question, the function stops when the L_2 norm of the gradient falls below the tolerance value, so the stopping criteria is defined in the function as such.

And the iterate was updated using the gradient descent algorithm:

$$\mathbf{x}^{(k+1)} = \mathbf{x}^{(k)} - \boldsymbol{\alpha} \cdot \nabla \mathbf{f}(\mathbf{x}^{(k)})$$

Also, a global iteration count variable was declared to see exactly how many iterations it took in every case to reach the tolerance limit.

So my function for step 1 looks like this:

```
def gradient descent(grad func, x0, step flag, alpha,
max iteration, tolerance):
   global iteration count # Declare the global variable first
   if alpha <= 0:
      print('Step size alpha must be greater than zero.')
      return None
  if step flag != 'fixed':
       print('Variable step size is currently not supported.')
      return None
  X = x.reshape(-1, 1) \# Store iterates as columns
   for k in range(max iteration):
      grad = grad func(x) # Compute gradient using the
      grad norm = 0.0
       for g in grad:
          grad norm += g**2
       grad norm = grad norm**0.5 # Take the square root
```

```
if grad_norm <= tolerance:
    iteration_count = k + 1  # Update the global variable
    break

# Gradient Descent Update
    x = x - alpha * grad

# Store iterate
    X = np.hstack((X, x.reshape(-1, 1)))

# If the loop completes without breaking, set iteration_count
to max_iteration
if grad_norm > tolerance:
    iteration_count = max_iteration

return X
```

Step_2: Taking the inputs as mentioned in question

Matrix Q1 and Q2 as mentioned in the question and their corresponding step sizes are taken.

Step_3: Function and gradient function

The given quadratic function was:

$$\mathbf{f}(\mathbf{x}) = \frac{1}{2} \mathbf{x}^{\mathrm{T}} \mathbf{Q} \mathbf{x}$$
So, $\nabla \mathbf{f}(\mathbf{x}) = \mathbf{Q} \mathbf{x}$

This is implemented in code as:

```
#Step_3: Function and gradient function

def grad_func1(x):

   return Q1 @ x # Gradient for Q1
```

```
def grad_func2(x):
    return Q2 @ x # Gradient for Q2

# Functions for quadratic function values
def f1(x):
    return 0.5 * x.T @ Q1 @ x # Quadratic function for Q1

def f2(x):
    return 0.5 * x.T @ Q2 @ x # Quadratic function for Q2
```

Step_4: Hyperparameters

The remaining hyperparameters in the design were initial point, maximum iterations and the stopping criteria tolerance. They were chosen as:

```
#Step_4: Hyperparameters
# Initial point
x0 = np.array([1, 1])
# Maximum iterations and tolerance
max_iter = 2000
tolerance = 1e-7
```

Step 5: Run Gradient Descent and Plot Results

Finally the gradient descent function was run on the given step size and matrix and the result (Contour plot of f(x), Function value and Gradient norm) were plotted in accordance.

Full code for my implementation:

```
import numpy as np
import matplotlib.pyplot as plt
# Global variable to store the number of iterations
iteration count = 0
#Step 1: Defining the gradient descent function
def gradient descent(grad func, x0, step flag, alpha,
max iteration, tolerance):
  global iteration count # Declare the global variable first
   #Check if a non-zero step size is given.
  if alpha <= 0:</pre>
       print('Step size alpha must be greater than zero.')
      return None
   # Check if step type is supported
  if step flag != 'fixed':
      print('Variable step size is currently not supported.')
      return None
  # Initialize
  x = x0
  X = x.reshape(-1, 1) # Store iterates as columns
   for k in range(max iteration):
       grad = grad func(x) # Compute gradient using the
provided grad func
       # Check stopping criterion
```

```
grad norm = 0.0
       for g in grad:
           grad norm += g**2
       grad norm = grad norm**0.5 # Take the square root
       # Check if the gradient norm is below the tolerance, break
if not.
       if grad norm <= tolerance:</pre>
           iteration count = k + 1 \# Update the global variable
           break
       # Gradient Descent Update
       x = x - alpha * grad
       # Store iterate
      X = np.hstack((X, x.reshape(-1, 1)))
   # If the loop completes without breaking, set iteration count
to max iteration
  if grad norm > tolerance:
       iteration count = max iteration
  return X
#Step 2: Taking the inputs as mentioned in question
# Define two different Q matrices
Q1 = np.array([[1, 0], [0, 1]])  # Case 1
Q2 = np.array([[10, 0], [0, 1]]) # Case 2
# Step sizes
alphas Q1 = [0.1, 0.5]
alphas Q2 = [0.01, 0.05]
```

```
#Step 3: Function and gradient function
# Functions for gradients
def grad func1(x):
  return Q1 @ x # Gradient for Q1
def grad func2(x):
  return Q2 @ x # Gradient for Q2
# Functions for quadratic function values
def f1(x):
  return 0.5 * x.T @ Q1 @ x # Quadratic function for Q1
def f2(x):
   return 0.5 * x.T @ Q2 @ x # Quadratic function for Q2
Q matrices = [Q1, Q2]
grad funcs = [grad func1, grad func2]
funcs = [f1, f2]
alpha sets = [alphas Q1, alphas Q2]
Q labels = ['Q1', 'Q2']
#Step 4: Hyperparameters
# Initial point
x0 = np.array([1, 1])
# Maximum iterations and tolerance
\max iter = 2000
tolerance = 1e-7
#Step 5: Run Gradient Descent and Plot Results
# Initialize figure counter
```

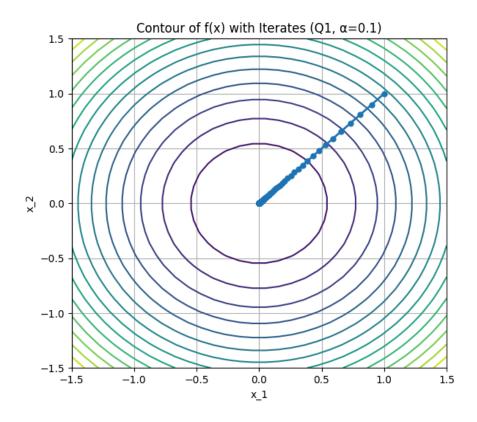
```
figure number = 1
for i in range(len(Q matrices)):
  Q = Q matrices[i]
  grad func = grad funcs[i]
  f = funcs[i]
  alphas = alpha sets[i]
  Q name = Q labels[i]
  for j in range(len(alphas)):
      alpha = alphas[j]
       # Reset the global variable before each run
      iteration count
       iteration count = 0
      # Run gradient descent
      X = gradient descent(grad func, x0, 'fixed', alpha,
max iter, tolerance)
       # Print the number of iterations for this case
      print(f'Case: {Q name}, Step size (α): {alpha},
Iterations: {iteration count}')
       # Compute function values and gradient norms
      num iters = X.shape[1]
       f vals = np.zeros(num iters)
      grad norms = np.zeros(num iters)
       for k in range(num iters):
           f vals[k] = f(X[:, k])
           grad norms[k] = np.linalg.norm(grad func(X[:, k]), 2)
```

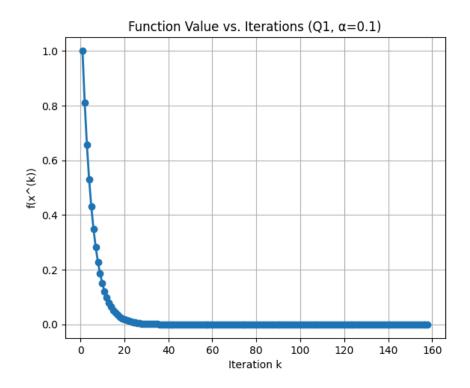
```
# (a) Contour Plot with Iterates
       plt.figure(figure number)
       figure number += 1
       X1, X2 = np.meshgrid(np.linspace(-1.5, 1.5, 30)),
np.linspace(-1.5, 1.5, 30))
       F \text{ vals} = np.array([[f(np.array([x1, x2])) for x1, x2 in ]))
zip(np.ravel(X1), np.ravel(X2))]]).reshape(X1.shape)
       plt.contour(X1, X2, F vals, 20)
       plt.plot(X[0, :], X[1, :], '-o', linewidth=2,
markersize=5)
       plt.title(f'Contour of f(x) with Iterates ({Q name},
\alpha = \{alpha\})')
       plt.xlabel('x 1')
      plt.ylabel('x 2')
       plt.grid(True)
       plt.show()
       # (b) Function Value vs. Iterations
       plt.figure(figure number)
       figure number += 1
       plt.plot(range(1, num iters + 1), f vals, '-o',
linewidth=2)
       plt.title(f'Function Value vs. Iterations ({Q_name}),
\alpha = \{alpha\})')
       plt.xlabel('Iteration k')
       plt.ylabel('f(x^{(k))'})
       plt.grid(True)
       plt.show()
       # (c) Gradient Norm vs. Iterations
       plt.figure(figure number)
       figure number += 1
```

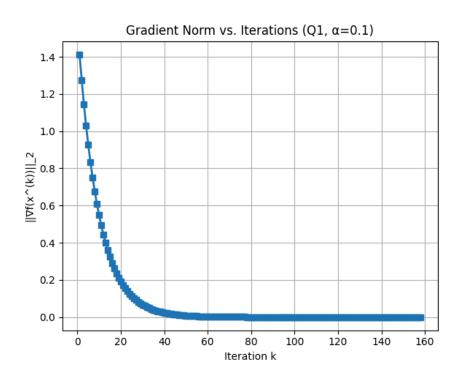
```
plt.plot(range(1, num_iters + 1), grad_norms, '-s',
linewidth=2)
    plt.title(f'Gradient Norm vs. Iterations ({Q_name}),
α={alpha})')
    plt.xlabel('Iteration k')
    plt.ylabel('||∇f(x^(k))||_2')
    plt.grid(True)
    plt.show()
```

Labeled output results for each case:

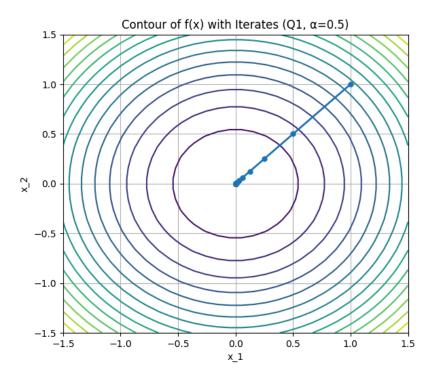
Case 1:
$$Q = [1 \ 0, 0 \ 1], \ \alpha = 0.1$$

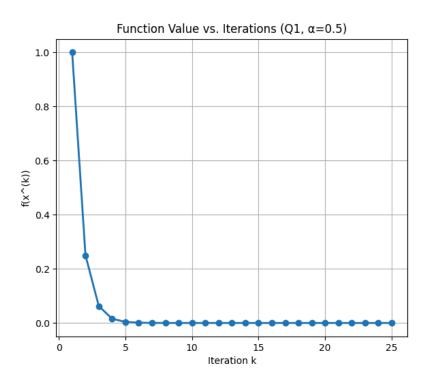


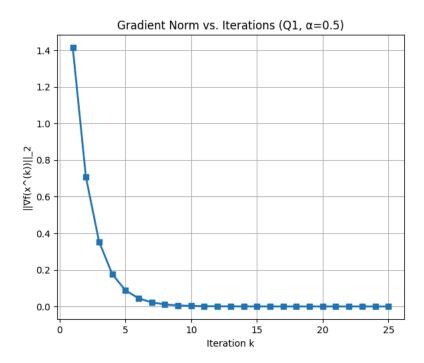




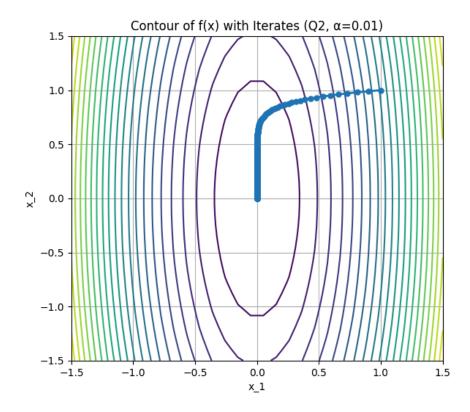
Case 2 : $Q = [1 \ 0 \ , 0 \ 1], \ \alpha = 0.5$

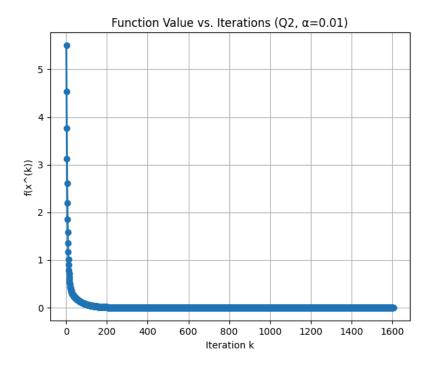


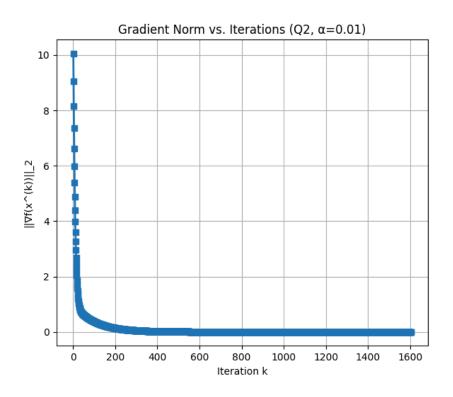




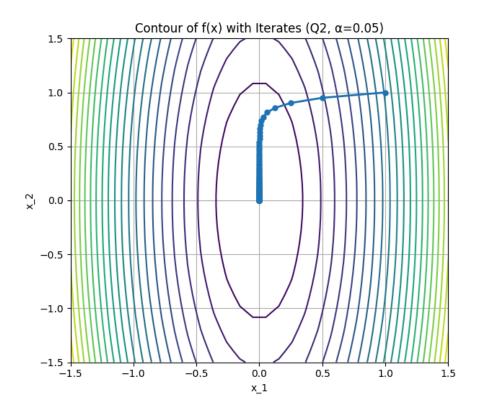
Case 3: $Q = [10 \ 0, 0 \ 1], \ \alpha = 0.01$

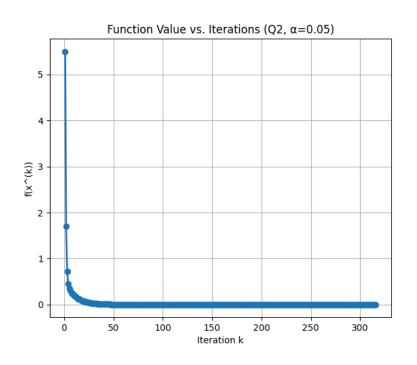


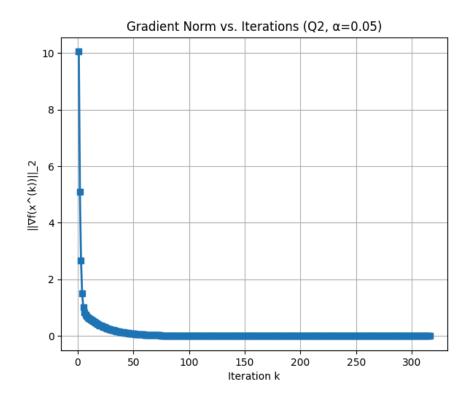




Case 4 : $Q = [10 \ 0, 0 \ 1], \ \alpha = 0.05$







Questions based on my results:

From my results and the plots:

Case: Q1, Step size (α): 0.1, Iterations: 158

Case: Q1, Step size (α): 0.5, Iterations: 25

Case: Q2, Step size (α): 0.01, Iterations: 1605

Case: Q2, Step size (α): 0.05, Iterations: 316

Q1. How does the choice of step size affect convergence behavior?

As we can see from the plots, for the smaller step sizes (like α =0.01 for Q2), the algorithm converges very slowly. It was expected because a small step size means the algorithm takes tiny steps toward the minimum, which slows down convergence.

On the other hand, for bigger step sizes (like α =0.5 for Q1), the algorithm converges very quickly. It was seen because a larger step size means the algorithm is now taking giant steps toward the minimum, which speeds up convergence.

However, this may not always be the case as the function can overshoot for very large step sizes.

For Moderate Step Size (like α =0.1 for Q1 or α =0.05 for Q2), the algorithm converges relatively quickly. Here, fewer iterations are needed to reach the minimum compared to a small step size. This moderate step size balances between convergence speed and stability.

Q2. How does changing the matrix Q affect convergence?

Here, The given quadratic function was:

$$\mathbf{f}(\mathbf{x}) = \frac{1}{2} \; \mathbf{x}^{\mathrm{T}} \, \mathbf{Q} \; \mathbf{x}$$

So, changing Q changes the shape of the overall quadratic function f(x), which eventually affects the convergence.

From the contour plot, for Q1, the contour plot shows circular contours. This indicates that f(x) is equally scaled in all directions for Q1. And from the function and gradient plots for Q1, we can see that gradient descent converges quickly and smoothly for both step sizes (α =0.1 and α =0.5). Most importantly, as evident from the contour plot, the algorithm takes a direct path toward the minimum for Q1, with larger step sizes leading to faster convergence.

On the other hand, for Q2, the contour plot shows elliptical contours. This indicates that f(x) is not equally scaled in all directions for Q2. And from the function and gradient plots for Q1, we can see that gradient descent converges slowly for both step sizes (α =0.01, α =0.05). Here, the algorithm takes a curved path toward the minimum for Q2, with larger step sizes leading to faster convergence, but overall slower convergence compared to Q1.

So, for a symmetric matrix like Q1, the algorithm converges quickly and smoothly. And for stretched matrices like Q2, the algorithm converges more slowly and may zigzag toward the minimum.