

Homework 3

Megi Dervishi

Exercise 1

We can re-write the LASSO problem as follows by introducing a new variable:

$$\begin{aligned} \min_{v, w} \quad & \frac{1}{2} \|Xw - y\|_2^2 + \lambda \|w\|_1 \\ \text{s.t.} \quad & v = Xw - y \end{aligned}$$

The Lagrangian is:

$$\begin{aligned} \mathcal{L}(v, w, \mu) &= \frac{1}{2} \|v\|_2^2 + \lambda \|w\|_1 + \mu^T (v - Xw + y) \\ &= \frac{1}{2} \|v\|_2^2 + \mu^T v + \lambda \|w\|_1 - (X^T \mu)^T w + y^T \mu \end{aligned}$$

And the dual function is:

$$g(\mu) = \inf_{v, w} \mathcal{L}(v, w, \mu) = y^T \mu + \underbrace{\inf_v \left(\frac{1}{2} \|v\|_2^2 + \mu^T v \right)}_{\text{a) convex and differentiable.}} + \underbrace{\inf_w \left(\lambda \|w\|_1 - (X^T \mu)^T w \right)}_{\text{b) use conjugate of } L_1 \text{ from thw2.}}$$

a) $\nabla_v \mathcal{L}(v, w, \mu) = 0 \Rightarrow v = -\mu$

Hence $\min_v \frac{1}{2} \|v\|_2^2 + \mu^T v = \frac{1}{2} \|\mu\|_2^2 - \mu^T \mu$
 $= -\frac{1}{2} \|\mu\|_2^2$

b) $\inf_w \lambda \|w\|_1 - (X^T \mu)^T w = \lambda \sup_w \left(\frac{1}{\lambda} X^T \mu \right)^T w - \|w\|_1$
 $= \lambda \left\| \frac{1}{\lambda} X^T \mu \right\|_1^* = \left\| \frac{1}{\lambda} X^T \mu \right\|_1^*$ since $\lambda > 0$

$$= \begin{cases} 0 & \text{if } \left\| \frac{1}{\lambda} X^T \mu \right\|_\infty \leq 1 \\ +\infty & \text{otherwise.} \end{cases} \quad \text{from } \|\cdot\|_1^*(u) = \begin{cases} 0 & \text{if } \|u\|_\infty \leq 1 \\ +\infty & \text{otherwise} \end{cases}$$

So we have that: $g(\mu) = y^T \mu - \frac{1}{2} \|\mu\|_2^2 + \left\| \frac{1}{\lambda} X^T \mu \right\|_1^*$

And we have:

$$\begin{aligned} \max_v \quad & y^T \mu - \frac{1}{2} \|\mu\|_2^2 \quad \text{which can be reformulated as:} \\ \text{s.t.} \quad & \left\| \frac{1}{\lambda} X^T \mu \right\|_\infty \leq 1 \end{aligned}$$

$$\left\| \frac{1}{\lambda} X^T \mu \right\|_\infty \leq 1 \Leftrightarrow \max_i \left| \frac{1}{\lambda} X^T \mu \right|_i \leq 1 \Leftrightarrow \forall_i -\lambda \leq X^T \mu \leq \lambda \Leftrightarrow \begin{pmatrix} X^T \mu \\ -X^T \mu \end{pmatrix} \leq \lambda \mathbf{1}_{2d}$$

Hence we have for $p = -y$, $Q = \frac{1}{2} \text{Id}$, $A = \begin{pmatrix} X^T \\ -X^T \end{pmatrix}$, $b = \lambda \mathbf{1}_{2d}$

$$\begin{aligned} \min_v \quad & \mu^T Q \mu + p^T \mu \\ \text{s.t.} \quad & A \mu \leq b. \end{aligned}$$

Exercise 2

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

#Question 2
b = 10

def g(Q,p,A,t,v):
    return t*(v.T @ Q @ v + p.T @ v) - np.sum(np.log(b - A @ v))

def backtrack_linesearch(Q,p,A,t,v,grad,newton_step):
    step_t = 1
    alpha = 0.05
    beta = 0.95
    def condition(step_t):
        c = g(Q,p,A,t,v+step_t * newton_step) > g(Q,p,A,t,v) + alpha*step_t*grad.T@newton_step
        return c
    while condition(step_t) or (any(b-A@(v+step_t*newton_step)<=0)):
        step_t = beta*step_t
    return step_t

#Centering step
def centering_step(Q,p,A,b,t,v0,eps):
    v = v0
    prev_v = v*np.inf
    v_seq = [v0]
    while np.linalg.norm(prev_v - v) > eps:
        h = 1/(b - A @ v)
        grad = t * (2*Q @ v + p) + A.T @ h
        hessian = 2 * t * Q + A.T @ np.diag(h.reshape(-1))*2 @ A
        newton_step = -np.linalg.pinv(hessian) @ grad
        step_t = backtrack_linesearch(Q,p,A,t,v,grad,newton_step)
        prev_v = v
        v = v + step_t * newton_step
        v_seq.append(v)
    return v_seq

def f(Q,p,v):
    return v.T @ Q @ v + p.T @ v
```

```
#Barrier method
def barr_method(Q,p,A,b,v0,eps,u):
    t = 1
    c = A.shape[0] #number of constraints
    v = v0
    v_seq = [v0]
    precision_seq = [c/t]
    fv = [f(Q,p,v)[0,0]]
    while (c/t > eps):
        v = centering_step(Q,p,A,b,t,v0,eps)[-1]
        v_seq.append(v)
        t = t*u
        precision_seq.append(c/t)
        fv.append(f(Q,p,v)[0,0])
    return v_seq, precision_seq, abs(fv-fv[-1])
```

Exercise 3.

```
#Question 3
lambda = 10
X = 0.4*np.random.randn(2, 100)
y = 4*np.random.randn(2).reshape(-1,1)
Q = np.eye(2)/2
p = -y
A = np.concatenate((X.T, - X.T), axis=0)
v0 = np.zeros(2).reshape(-1,1)+0.01
eps = 1e-8

# colorplot
dstep = 0.1
pmin, pmax = -1.2*lambda, 1.2*lambda
cory, corx = np.mgrid[pmin:pmax+dstep:dstep, pmin:pmax+dstep:dstep]
potential = np.zeros(cory.shape)
for i in range(cory.shape[0]):
    for j in range(cory.shape[1]):
        cur_point = np.array([corx[i, j], cory[i, j]])
        #print(cur_point.shape, Q.shape, p.shape)
        if (A @ cur_point <= b).all():
            potential[i, j] = cur_point.T @ Q @ cur_point + p.T @ cur_point
        else:
            potential[i, j] = np.inf

print(potential.shape)

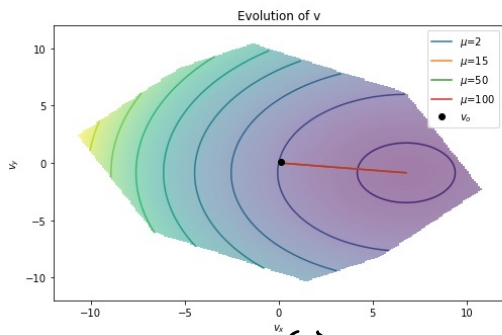
v_seq_u = dict((u, []) for u in [2,15,50,100])
precision_u = dict((u, []) for u in [2,15,50,100])
gap_u = dict((u, []) for u in [2,15,50,100])

for u in [2,15,50,100]:
    v_seq, precision, gap = barr_method(Q, p, A, v0, eps, u)
    v_seq_u[u] = np.array(v_seq).squeeze()
    precision_u[u] = precision
    gap_u[u] = gap
```

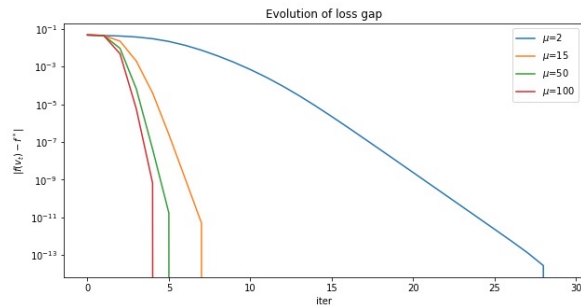
```
#plots
plt.figure(figsize =(10,5))
plt.contour(corx, cory, potential)
plt.pcolor(corx, cory, potential, alpha=0.5, snap=True)
colorbar = plt.colorbar()
colorbar.set_alpha(0.5)
for u in v_seq_u.keys():
    plt.plot(v_seq_u[u][:,0], v_seq_u[u][:,1], label=f'$\mu$={u}')
    plt.plot(0.1, 0.1, 'o', color='black', label='$v_o$')
plt.title('Evolution of v')
plt.xlabel('$v_x$')
plt.ylabel('$v_y$')
plt.legend()
plt.show()

plt.figure(figsize =(10,5))
for u in precision_u.keys():
    plt.semilogy(precision_u[u], label=f'$\mu$={u}')
plt.title('Evolution of precision')
plt.xlabel('iter')
plt.ylabel('$\epsilon_c/t$')
plt.legend()
plt.show()

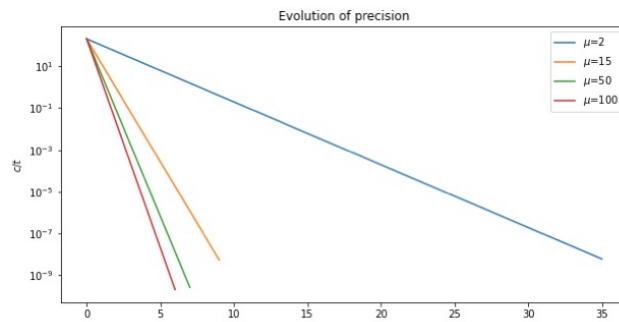
plt.figure(figsize =(10,5))
for u in precision_u.keys():
    plt.semilogy(loss_u[u], label=f'$\mu$={u}')
plt.title('Evolution of loss gap')
plt.xlabel('iter')
plt.ylabel('$|f(v_t)-f^*|$')
plt.legend()
plt.show()
```



(a)



(b)



(c)

We can see from Plot (a) that all solutions converge, but having $\mu=50$ or $\mu=100$ takes less backtracking (outer) iterations to reach the solution than with $\mu=2$. So an appropriate μ could be $\mu=50$.