

# CONVEX Optimisation: HW3\_LASSO

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## PART 1: Derive the dual problem of LASSO and format it as a general Quadratic Problem

CHAUVIN Paul 1) We have  $\min_w \frac{1}{2} \|Xw - y\|_2^2 + d \|w\|_1$

on introduit  $r = Xw - y$  :

$$\min_{r, w} \frac{1}{2} \|r\|_2^2 + d \|w\|_1$$

Nous avons alors un problème d'optimisation convexe et nous pouvons écrire son Lagrangien

$$\mathcal{L}(w, r, v) = (d \|w\|_1 + v^T Xw) + \left(\frac{1}{2} \|r\|_2^2 - v^T r\right) - (v^T y)$$

$$g(v) = \inf_{w, r} \mathcal{L}(w, r, v) \quad (\text{Derive the dual})$$

$$\Rightarrow g(v) = \inf_w (d \|w\|_1 + v^T Xw) + \inf_r \left(\frac{1}{2} \|r\|_2^2 - v^T r\right) - y^T v$$

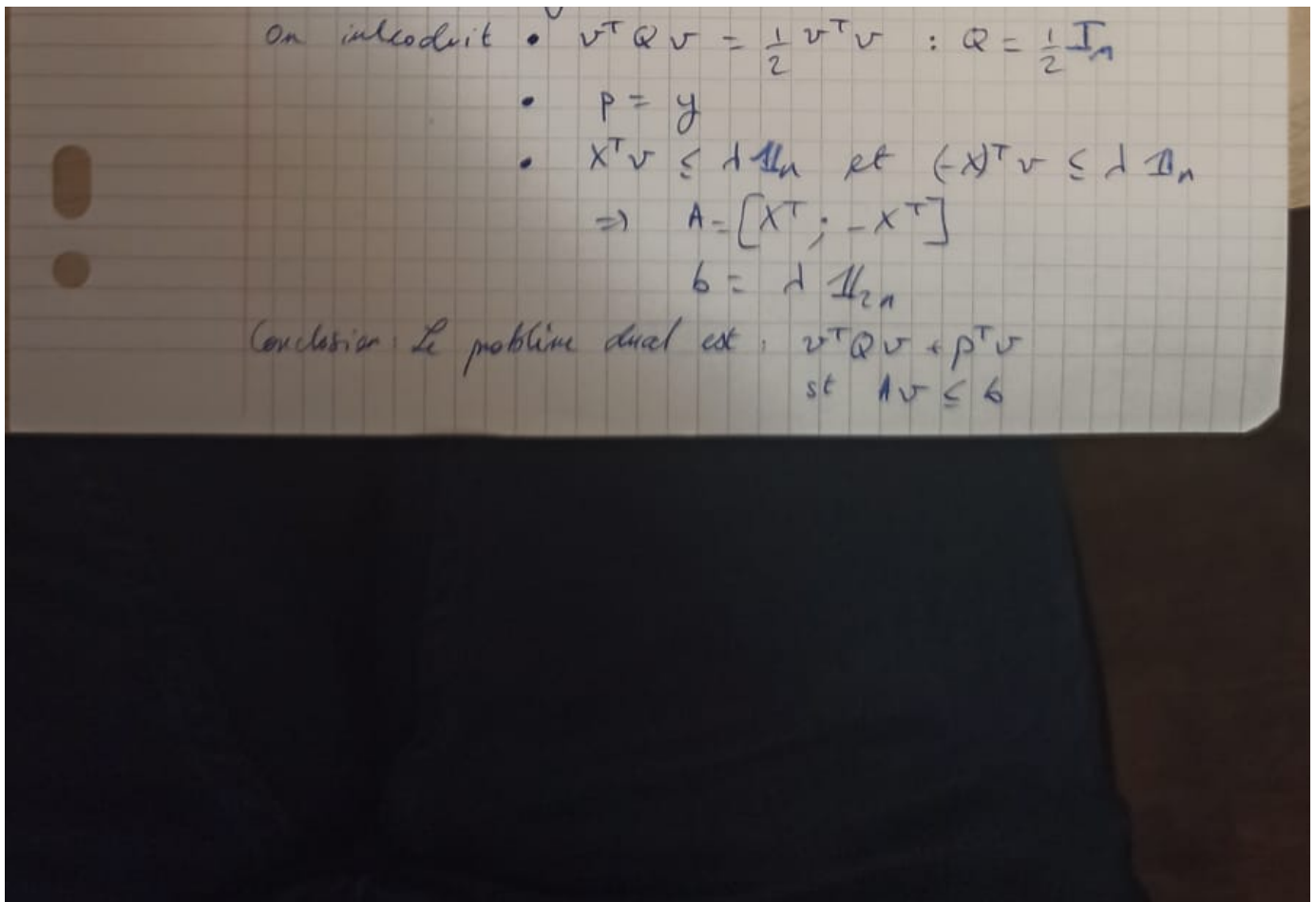
$$\neq \inf_w (d \|w\|_1 + v^T Xw) = -\sup_w (-d \|w\|_1 - v^T Xw)$$

$$= \begin{cases} 0 & \text{si } 0 \leq \|X^T v\|_\infty \leq d \\ -\infty & \text{sinon} \end{cases}$$

$$\neq \inf_r \left(\frac{1}{2} \|r\|_2^2 - v^T r\right) = -\sup_r \left(v^T r - \frac{1}{2} \|r\|_2^2\right) = -\frac{1}{2} v^T v$$

$$\Rightarrow g(v) = -\frac{1}{2} v^T v - y^T v \quad \text{tel que } \|X^T v\|_\infty \leq d$$

Or maximiser  $g$  revient à minimiser (QP)



## ▼ Part 2: Implement the barrier method to solve QP

### ▼ Environment

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

### ▼ Parameters

```
alpha = 0.1
beta = 0.7
mu = 5
n = 20
d = 15
eps = 10e-6
lambda = 10
```

### ▼ Useful functions

```

def function(v, Q, p, t0):
    return t0*(np.dot(v.T, np.dot(Q, v)) + np.dot(p.T, v)) - sum([np.log(b[i]-np.do

def function_true(v, Q, p):
    return np.dot(v.T, np.dot(Q, v)) + np.dot(p.T, v)

def line_search(Q, p, v, t, df, dx, t0):
    if function(v + t*dx, Q, p, t0) <= (function(v, Q, p, t0) + alpha*t*np.dot(df.T
        return v + t*dx
    else:
        return line_search(Q, p, v, beta*t, df, dx, t0)

```

## ▼ Barrier method

```

def centering_step(Q, p, A, b, t, v0, eps, numb_iter=0):
    df = t*(2*np.dot(Q, v0) + p) + \
        sum([A[i, np.newaxis].T/(b[i]-np.dot(A[i], v0)) for i in range(b.shape[0])])
    d2f = 2*t*Q + \
        sum([(np.outer(A[i, np.newaxis].T, A[i, np.newaxis].T)) / ((b[i] - np.dot(A
    dx = -1*np.dot(np.linalg.inv(d2f), df)
    l2 = np.dot(df.T, np.dot(np.linalg.inv(d2f), df))
    if l2/2 <= eps:
        return v0, numb_iter
    v1 = line_search(Q, p, v0, t=1, df=df, dx=dx, t0=t)
    return centering_step(Q, p, A, b, t, v1, eps, numb_iter+1)

def barr_method_inter(Q, p, A, b, v0, eps, t, mu, numb_iter=0, numb_newton = [], v_
    v_center, numb_iter_inter = centering_step(Q, p, A, b, t, v0, eps)
    numb_newton.append(numb_iter)
    v_seq.append(v_center)
    f_seq_true.append(function_true(v_center, Q, p)[0][0])
    if b.shape[0]/t < eps:
        return v_center, numb_newton, v_seq, f_seq_true
    else:
        t = mu*t
        return barr_method_inter(Q, p, A, b, v0, eps, t, mu, numb_iter+numb_iter_in

def barr_method(Q, p, A, b, v0, eps, mu):
    numb_newton = [0]
    v_seq = [v0]
    f_seq_true = [function_true(v0, Q, p)[0][0]]
    return barr_method_inter(Q, p, A, b, v0, eps, 1, mu, 0, numb_newton, v_seq, f_s

```

## ▼ Part 3: Test our functions

## ▼ Initialization

```
X = np.random.rand(n,d)
y = np.random.rand(n,1)
Q = 0.5*np.eye(n)
p = -y
A = np.vstack((X.T,-X.T))
b = lambd*np.ones((2*d,1))

v0 = np.zeros((n,1))

mu_list = [2, 5, 10, 15, 20, 25, 30, 40, 50, 70, 85, 100, 130, 170, 200, 250]
mu_list_restricted = [2, 5, 15, 50, 100, 200]
w_center_list = []
f_true_list = []
```

## ▼ Plot results

```
plt.figure()
for mu in mu_list:
    v_center, numb_newton, v_seq, f_seq_true = barr_method(Q, p, A, b, v0, eps, mu)
    w_center = np.linalg.lstsq(X,-v_center-y)[0]
    w_center_list.append(w_center)
    f_true_list.append(f_seq_true[-1])
    if mu in mu_list_restricted:
        plt.step(numb_newton, f_seq_true - f_seq_true[-1], label='mu = '+str(mu))
plt.legend()
plt.ylabel("Norm of f(mu) - f(mu)*")
plt.xlabel("Centering step number")
plt.semilogy()
plt.show()
```

```

/usr/local/lib/python3.7/dist-packages/ipykernel_launcher.py:4: FutureWarning:
To use the future default and silence this warning we advise to pass `rcond=Nc
after removing the cwd from sys.path.
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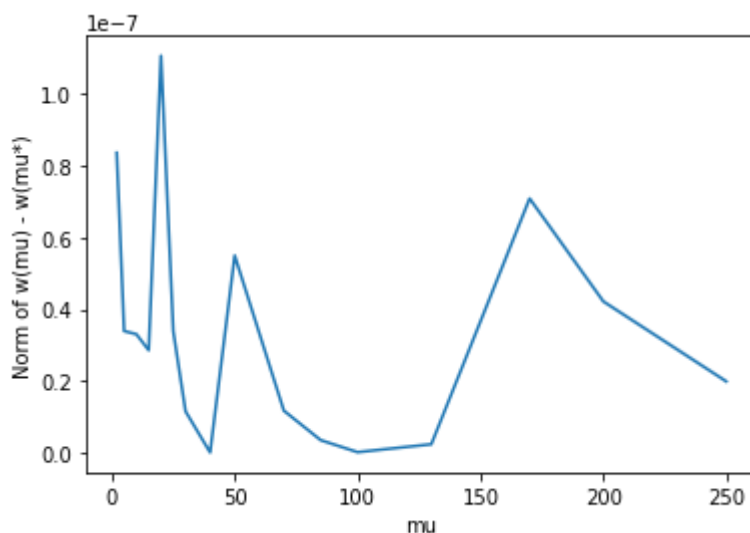
## ▼ Plot W in function of mu

10<sup>4</sup>

```

mu_min = np.argmin(f_true_list)
plt.figure()
w_diff_norm = [np.linalg.norm(w-w_center_list[mu_min]) for w in w_center_list]
plt.plot(mu_list, w_diff_norm)
plt.ylabel("Norm of w(mu) - w(mu*)")
plt.xlabel("mu")
plt.show()

```



```

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```

When  $\mu$  is too small, we need to many iterations to converge. Whereas with smaller  $\mu$ , it's faster, we need less iterations, however, each step is going to be more costly. An appropriate value of  $\mu$  would be around 40.

Produits payants Colab - Résilier les contrats ici

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