

# Computational Statistics - TD1

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(The notebook of all implementations can be found [here](#))

## Exercise 1: Lasso

1)

Let  $z = Xw - y$ . (LASSO) can be reformulated as:

$$\begin{aligned} & \text{minimize } \frac{1}{2} z^T z + \lambda \|w\|_1 \\ & \text{s.t. } Xw - y = z \end{aligned}$$

The Lagrangian is:

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

$\mathcal{L}$  is a quadratic form with respect to  $z$ , minimized for in  $v$ . To minimize with respect to  $w$  we use what was done for the previous homework:

$$\begin{aligned} \inf_w \left( \lambda \|w\|_1 - \left( - (X^T v)^T w \right) \right) &= -\lambda \sup_w \left( -\frac{1}{\lambda} (X^T v)^T w - \|w\|_1 \right) \\ &= -\lambda f_* \left( -\frac{1}{\lambda} X^T v \right) \end{aligned} \quad \text{With } f_* \text{ the dual of } \|\cdot\|_1$$

This add the constraint  $\|X^T v\|_\infty \leq \lambda$

Thus the dual problem is

$$\begin{aligned} & \text{maximize } -\frac{1}{2} v^T I v - v^T y \\ & \text{s.t. } \|X^T v\|_\infty \leq \lambda \end{aligned}$$

The inequality constraint can be rewritten as a linear system by writing that

$$\forall i \in \llbracket 1, d \rrbracket, -\lambda \leq (X^T v)_i \leq \lambda$$

with  $(X^T v)_i$  the  $i$ -th row of  $X^T v$ , leading to

$$\forall i \in \llbracket 1, d \rrbracket, -X_i^T v \leq \lambda \quad \text{and} \quad X_i^T v \leq \lambda$$

We thus have  $2d$  inequalities, boiling down to

$$\begin{pmatrix} X^T \\ -X^T \end{pmatrix} v = \begin{pmatrix} \lambda I_d \\ \lambda I_d \end{pmatrix}$$

We find the dual problem under its expected formulation with

$$A = \begin{pmatrix} X^T \\ -X^T \end{pmatrix} \in \mathbb{R}^{2d \times n} \quad B = \begin{pmatrix} \lambda I_d \\ \lambda I_d \end{pmatrix} \in \mathbb{R}^{2d} \quad Q = \frac{1}{2} I_n \quad p = y$$

2)

We first define some utility functions for the rest of the exercise

```
def f0(Q, p, v):
    return v@Q@v + p@v

def f(Q, p, A, b, t, v):
    in_log = (b - A@v)
    if np.any(in_log<0):
        return np.inf
    return t*(v@Q@v) + t*p@v - np.sum(np.log(in_log))

def grad_f(Q, p, A, b, t, v):
    denom = 1 / (b - (A@v)) # (2p,)
    return 2*t*(Q@v) + t*p + np.einsum('ij,i->j', A, denom)

def hess_f(Q, p, A, b, t, v):
    denom = 1 / (b - A @ v)**2 # (2p,)
    A_matrices = np.einsum('ij,ik->ijk', A, A) # (2p, n, n)
    sum_matrices = np.sum(denom[:,None,None] * A_matrices, axis=0) # (n, n)
    return 2*t*Q + sum_matrices
```

Below are provided the codes for the centering step and the barrier method

```
def centering_step(Q, p, A, b, t, v0, eps):
    v_arr = [v0]
    v = v0.copy()

    grad = grad_f(Q, p, A, b, t, v) # (n,)
    H_inv = np.linalg.inv(hess_f(Q, p, A, b, t, v)) # (n, n)
    delta_v = -H_inv @ grad
    lamb_2 = -grad@delta_v

    # Backtracking line search
    def back_line_search(delta_v, beta=0.9, alpha=0.1, max_iter=500):
        step = 1
        while f(Q, p, A, b, t, v + step*delta_v) > f(Q, p, A, b, t, v) \
            -alpha*step*grad@delta_v:
            step *= beta
        return step

    while lamb_2 / 2 > eps:
        best_step = back_line_search(delta_v)
        v += best_step * delta_v
        v_arr.append(v)

        grad = grad_f(Q, p, A, b, t, v) # (n,)
        H_inv = np.linalg.inv(hess_f(Q, p, A, b, t, v)) # (n, n)
        lamb_2 = -grad @ delta_v
        delta_v = -H_inv @ grad

    return v_arr
```

```

def barr_method(Q, p, A, b, v0, eps, mu):
    m = A.shape[0]
    v_arr = [v0]
    v = v0.copy()
    t = 1
    while m/t >= eps:
        v_center = centering_step(Q, p, A, b, t, v, eps)
        v_arr+=v_center
        t *= mu
        v = v_center[-1]
    return np.array(v_arr)

```

3)

$X$  and  $y$  were generated using with normal distribution. We set  $\varepsilon = 10^{-7}$ ,  $n = 20$ ,  $d = 300$  and  $\mu \in \{5, 15, 50, 100, 500, 400, 2000, 5000\}$ . The random seed is fixed to 0.

Here is the code used to generate data and plots

```

def init_params(n, d, l):
    """Utility function to initialize the problem"""

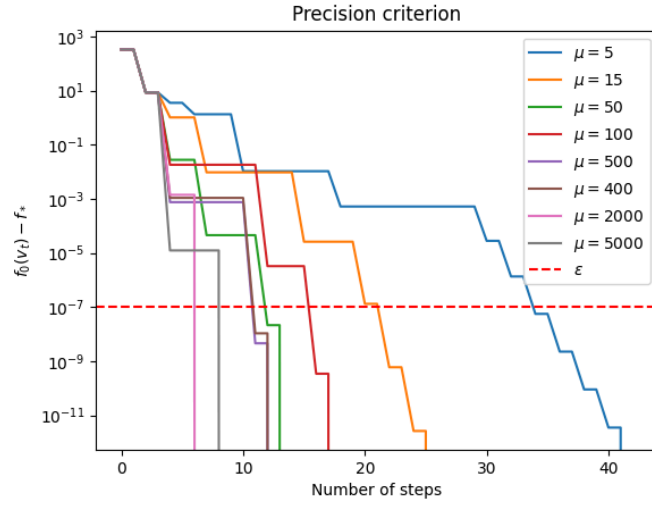
    X = np.random.normal(size=(n, d))
    Q = 0.5 * np.eye(n)
    p = np.random.normal(size=n) # =y
    A = np.vstack([X.T, -X.T])
    b = l * np.ones(2*d)
    v0 = np.random.normal(size=n)
    eps = 1e-7
    return Q, p, A, b, v0, eps

n = 10
d = 300
lamb = 10
Q, p, A, b, v0, eps = init_params(d, n, lamb)

all_v_arr = []
mu_vals = [5, 15, 50, 100, 500, 400, 2000, 5000]
for mu in tqdm(mu_vals):
    v_arr = barr_method(Q, p, A, b, v0, eps, mu)
    all_v_arr.append(v_arr)

plt.figure()
plt.title('Precision criterion')
for mu, v_arr in zip(mu_vals, all_v_arr):
    f_vals = []
    for v in v_arr:
        f_vals.append(f0(Q, p, v))
    plt.plot(np.array(f_vals)-np.min(f_vals), label=f'$\mu=${mu}')
    plt.ylabel('$f_0(v_t)-f_*$')
    plt.xlabel('Number of steps')
    plt.yscale('log')
plt.axhline(eps, ls='--', c='r', label='$\epsilon$')
plt.legend()
plt.show()

```



We notice that up to a certain value of  $\mu$ , the higher the faster. However,  $\mu = 2000$  requires fewer steps than  $\mu = 5000$ .

In this case  $\mu = 2000$  is an appropriate value