

# Exercise 6

Machine Learning I

## 6A-1.

First, rearrange through Bayes theorem:

$$p(\mathbf{t}|\mathbf{X}) = \frac{p(\mathbf{t}|\mathbf{w}, \mathbf{X})p(\mathbf{w}|\mathbf{X})}{p(\mathbf{w}|\mathbf{t}, \mathbf{X})}$$

Decompose:

$$\log p(\mathbf{t}|\mathbf{X}) = \log p(\mathbf{t}|\mathbf{w}, \mathbf{X}) + \log p(\mathbf{w}|\mathbf{X}) - \log p(\mathbf{w}|\mathbf{t}, \mathbf{X})$$

We already determined the densities for the likelihood, posterior and prior in the previous exercises:

$$p(\mathbf{t}|\mathbf{w}, \mathbf{X}) = N(\mathbf{t}; \boldsymbol{\Phi}\mathbf{w}, \tau_\epsilon^{-1}\mathbf{I})$$

$$p(\mathbf{w}|\mathbf{X}) = N(\mathbf{w}; \mathbf{0}, \tau_0^{-1}\mathbf{I})$$

$$p(\mathbf{w}|\mathbf{t}, \mathbf{X}) = N(\mathbf{w}; \tau_e \boldsymbol{\Sigma}_N \boldsymbol{\Phi}^T \mathbf{t}, (\tau_0 \mathbf{I} + \tau_e \boldsymbol{\Phi}^T \boldsymbol{\Phi})^{-1})$$

No need to recompute everything again. Because above has to be valid for any  $\mathbf{w}$ , we pick  $\mathbf{w} = \boldsymbol{\mu}_N$  for convenience. Let us first convert only the exponent

$$\begin{aligned} \log p(\mathbf{t}|\mathbf{X}) &\propto -\frac{1}{2}(\mathbf{t} - \boldsymbol{\Phi}\mathbf{w})^T \tau_\epsilon \mathbf{I} (\mathbf{t} - \boldsymbol{\Phi}\mathbf{w}) - \frac{1}{2} \mathbf{w}^T \tau_0 \mathbf{I} \mathbf{w} \\ &+ \frac{1}{2}(\mathbf{w} - \tau_e \boldsymbol{\Sigma}_N \boldsymbol{\Phi}^T \mathbf{t})^T (\tau_0 \mathbf{I} + \tau_e \boldsymbol{\Phi}^T \boldsymbol{\Phi})(\mathbf{w} - \tau_e \boldsymbol{\Sigma}_N \boldsymbol{\Phi}^T \mathbf{t}) \\ &= -\frac{1}{2}(\mathbf{t} - \boldsymbol{\Phi}\boldsymbol{\mu}_N)^T \tau_\epsilon \mathbf{I} (\mathbf{t} - \boldsymbol{\Phi}\boldsymbol{\mu}_N) - \frac{1}{2} \boldsymbol{\mu}_N^T \tau_0 \mathbf{I} \boldsymbol{\mu}_N \quad \text{set } \mathbf{w} = \boldsymbol{\mu}_N \end{aligned}$$

Now let us add the normalization constants:

$$\log p(\mathbf{t}|\mathbf{X}) = -\frac{1}{2}(\mathbf{t} - \boldsymbol{\Phi}\boldsymbol{\mu}_N)^T \tau_\epsilon \mathbf{I} (\mathbf{t} - \boldsymbol{\Phi}\boldsymbol{\mu}_N) - \frac{1}{2} \boldsymbol{\mu}_N^T \tau_0 \mathbf{I} \boldsymbol{\mu}_N + \frac{1}{|r_0|}$$

$$\begin{aligned} \log p(\mathbf{t}|\mathbf{X}) &= -\frac{1}{2}(\mathbf{t} - \boldsymbol{\Phi}\boldsymbol{\mu}_N)^T \tau_\epsilon \mathbf{I} (\mathbf{t} - \boldsymbol{\Phi}\boldsymbol{\mu}_N) - \frac{1}{2} \boldsymbol{\mu}_N^T \tau_0 \mathbf{I} \boldsymbol{\mu}_N \\ &+ \frac{d}{2} \log r_0 + \frac{n}{2} \log r_\epsilon - \frac{n}{2} \log 2\pi + \frac{1}{2} \log |\boldsymbol{\Sigma}_n| \end{aligned}$$

## 6A-2.

Let  $w_{MAX}$  be the (unregularized) maximum likelihood parameters. The lasso path illustrates the current selection of weights  $w_i$  given weight regularization  $\|w_{current}\| = \frac{t}{\|w_{MAX}\|_1}$ ,  $t \in [0, \|w_{MAX}\|_1]$ . For a very thorough and detailed explanation, please refer to page 69 of "Elements of statistical learning".

You can find the online version here:

[https://web.stanford.edu/~hastie/ElemStatLearn//printings/ESLII\\_print10.pdf](https://web.stanford.edu/~hastie/ElemStatLearn//printings/ESLII_print10.pdf)

This means that for low values on the  $x$  axis, the model is heavily regularized. Moving along the  $x$  axis then relaxes the regularization constraint until you get to the unregularized maximum likelihood solution  $\|w_{MAX}\|$ .

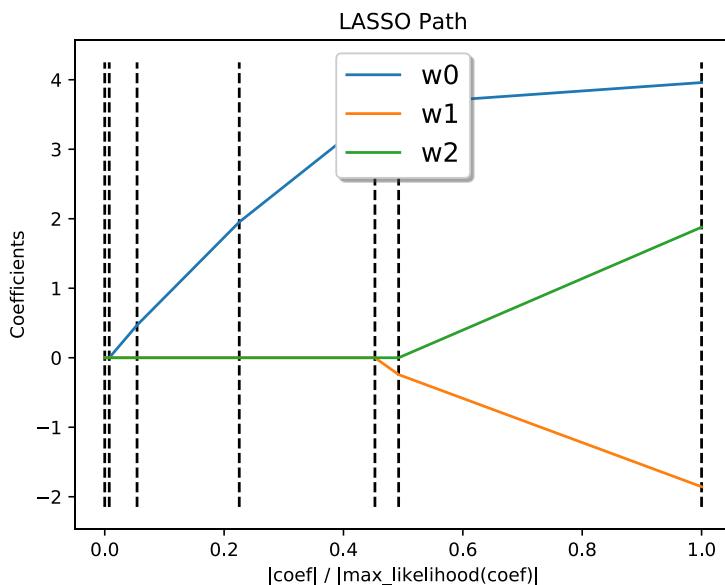
In order to study what happens, we are now using only 3 features. Using a design matrix, we can model each degree  $x^i$  of the regression polynomial  $y(x) = \sum_{i=0}^d w_i x^i$  as a separate dimension for the lasso path. Let us have

$$f(x) = 4 - 2x + 2x^2$$

Given enough samples, the maximum likelihood solution should set our regressor weights as such:

$$w_0 = 4, w_1 = -2, w_2 = 2,$$

This can be seen, if we look at the right end of our lasso path:



The less we regularize our weights (we move towards 1 on the abscissa) the more we obtain the maximum likelihood solution for our weights. Model was created with  $n=3000$  samples and sample interval  $I = (0,2)$ .

Now something should be noted: In the above image, weight  $w_0$  was picked up first, then  $w_1$  and lastly  $w_2$ . This means, in this case, relaxing the lasso translates to gradually picking up model complexity.

Now we are ready for eight features. Our polynomial is

$$g(x) = 2 - 3x + 4x^2 - 5x^3 + 4x^4 - 4x^5 + 6x^6 - 5x^7$$

The order of weight pick up is very much dependent on the sample interval  $I$ . Shown are two pictures of the same function  $g(x)$ , one time with a sample interval of  $I = (0,25)$  and one time with  $I = (-1,1)$ :

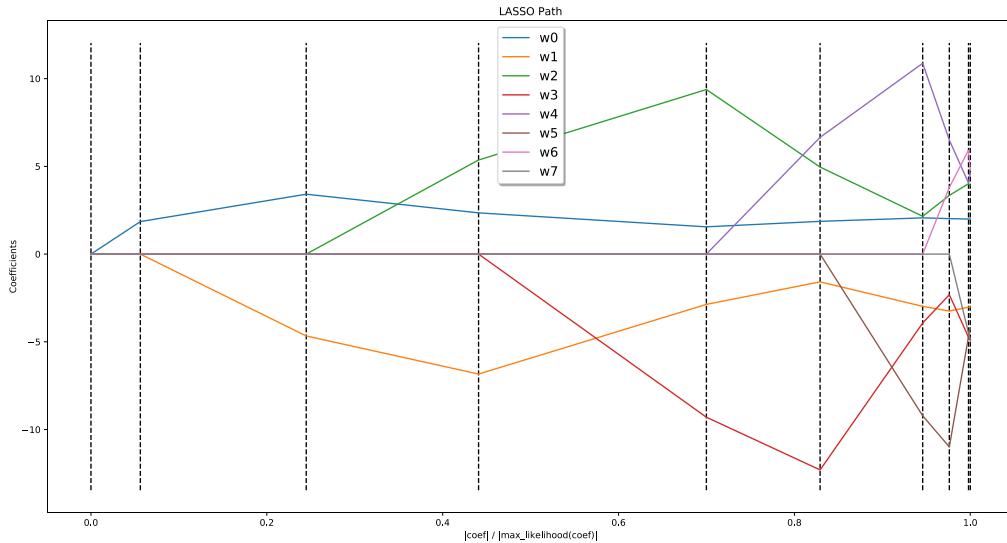


Figure 1  $n=70000, I=(-1,1)$ . The weights are picked up in ascending order of degree.

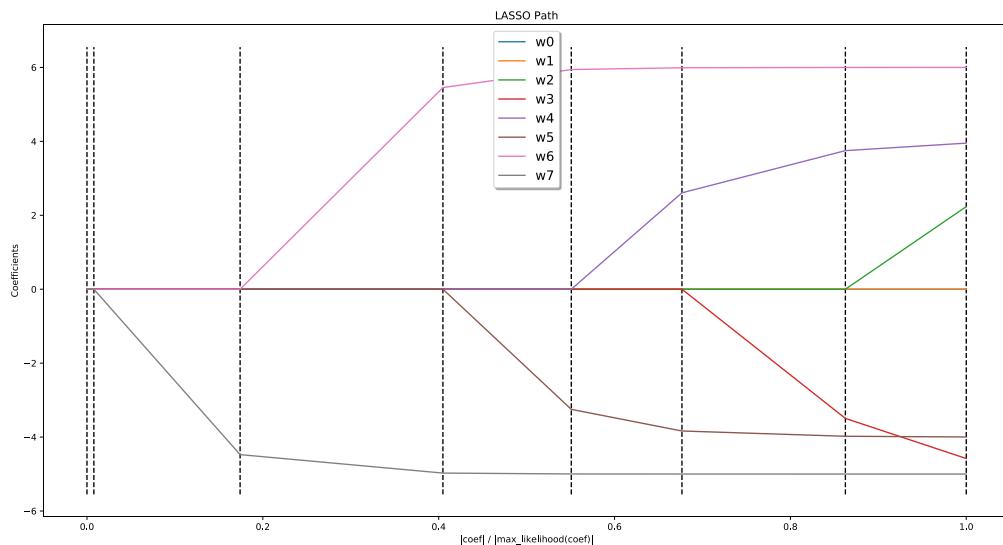


Figure 2  $n=70000, I=(0,25)$ . The weights are mostly picked up in descending order.

Think about why this happens (Tip: With polynomial basis functions, we have  $x^i \geq x^j, i < j, x \in [-1,1]$ . What does this mean for our regularized error?).

If we put the length of our interval between the ascending ( $I=(-1,1)$ ) and descending ( $I=(0,25)$ ) cases, then the following more chaotic weight pick up happens:

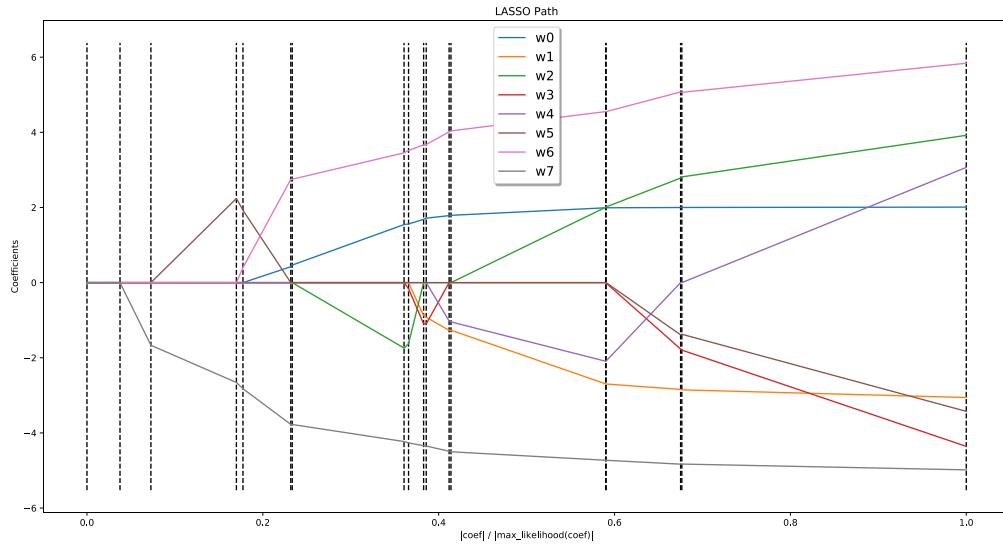


Figure 3  $n=70000, I=(0,2)$  Chaotic pick up of weights

### 6A-3.

The noise  $\epsilon$  is symmetric around  $t$ . So by the same way we derived  $t \sim N(t; \mathbf{w}^T \mathbf{x}, \epsilon)$  when the noise was normal, we can now say

$$t \sim \text{Stud}_t(t; \nu, \mathbf{w}^T \mathbf{x}, \sigma_\epsilon^2)$$

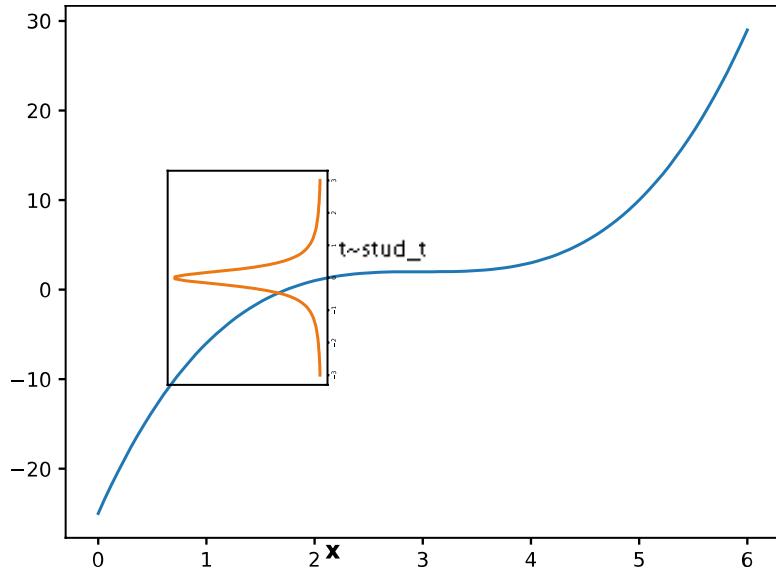
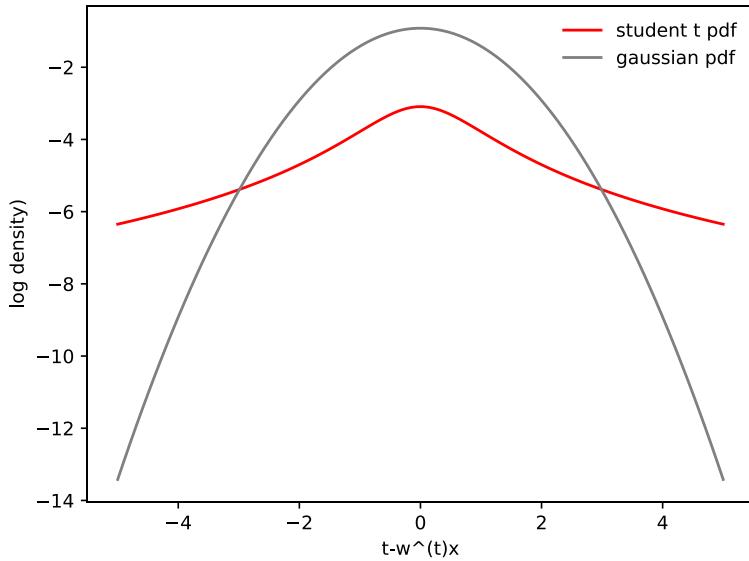


Figure 4 Because the noise is symmetrically distributed around  $t$  and the hidden function  $f(x)$  is non-random ( $p(t|x)=1$  for one value of  $t$ ),  $t$  is identically distributed to its noise with adjusted mean.

This leads to the log likelihood:

$$\begin{aligned}\log p(\mathbf{t}|\mathbf{x}, \nu, \sigma_\epsilon) &= \log \prod_{i=1}^n \text{Stud\_t}(t_i; \nu, \mathbf{w}^T \mathbf{x}_i, \sigma_\epsilon^2) \\ &= n \left[ \log \Gamma\left(\frac{\nu+1}{2}\right) - \log \left( \sqrt{\nu \pi \sigma_\epsilon^2} \Gamma\left(\frac{\nu}{2}\right) \right) \right] \\ &\quad - \frac{(\nu+1)}{2} \sum_{i=1}^n \log \left( 1 + \frac{1}{\nu} \frac{(t_i - \mathbf{w}^T \mathbf{x}_i)^2}{\sigma_\epsilon^2} \right)\end{aligned}$$

Let us set  $d = (t_1 - \mathbf{w}^T \mathbf{x}_1)$ . We now plot the error of one datapoint in comparison to a log normal pdf:



*Figure 5 Comparison of the test error between a log normal pdf and a student t pdf. Image was attained with the parameters  $\sigma_\epsilon = 1, \nu = 1$ .*

For small errors  $\mathbf{w}^T \mathbf{x} \approx t$ , the student t distribution appears to be less forgiving, as less density is concentrated around small errors ( $\log \text{Stud\_t}(t_1; 1, d, 1) < \log N(t_1; d, 1)$ ). Larger errors  $|t - \mathbf{w}^T \mathbf{x}| \gg 1$  are less penalized however, as  $\log N(t_1; d, 1) < \log \text{Stud\_t}(t_1; 1, d, 1)$ .

Generally speaking, this means if we try to maximize our weights  $\mathbf{w}$  with respect to  $e \sim N(0, \sigma^2)$ , then our fit is trying to mitigate large errors while being mostly indifferent to smaller ones.

If we maximize  $\mathbf{w}$  with respect to  $t \sim \text{Stud\_t}(t; \nu, \mathbf{w}^T \mathbf{x}, \sigma_\epsilon^2)$ , we get a more balanced fit.