

Feel free to work with other students, but make sure you write up the homework and code on your own (no copying homework *or* code; no pair programming). Feel free to ask students or instructors for help debugging code or whatever else, though.

**1 (Murphy 12.5 - Deriving the Residual Error for PCA)** It may be helpful to reference section 12.2.2 of Murphy.

(a) Prove that

$$\left\| \mathbf{x}_i - \sum_{j=1}^k z_{ij} \mathbf{v}_j \right\|^2 = \mathbf{x}_i^\top \mathbf{x}_i - \sum_{j=1}^k \mathbf{v}_j^\top \mathbf{x}_i \mathbf{x}_i^\top \mathbf{v}_j.$$

Hint: first consider the case when  $k = 2$ . Use the fact that  $\mathbf{v}_i^\top \mathbf{v}_j$  is 1 if  $i = j$  and 0 otherwise. Recall that  $z_{ij} = \mathbf{x}_i^\top \mathbf{v}_j$ .

(b) Now show that

$$J_k = \frac{1}{n} \sum_{i=1}^n \left( \mathbf{x}_i^\top \mathbf{x}_i - \sum_{j=1}^k \mathbf{v}_j^\top \mathbf{x}_i \mathbf{x}_i^\top \mathbf{v}_j \right) = \frac{1}{n} \sum_{i=1}^n \mathbf{x}_i^\top \mathbf{x}_i - \sum_{j=1}^k \lambda_j.$$

Hint: recall that  $\mathbf{v}_j^\top \Sigma \mathbf{v}_j = \lambda_j \mathbf{v}_j^\top \mathbf{v}_j = \lambda_j$ .

(c) If  $k = d$  there is no truncation, so  $J_d = 0$ . Use this to show that the error from only using  $k < d$  terms is given by

$$J_k = \sum_{j=k+1}^d \lambda_j.$$

Hint: partition the sum  $\sum_{j=1}^d \lambda_j$  into  $\sum_{j=1}^k \lambda_j$  and  $\sum_{j=k+1}^d \lambda_j$ .

a. We know that:

$$\begin{aligned} \left\| \mathbf{x}_i - \sum_{j=1}^k z_{ij} \mathbf{v}_j \right\|^2 &= \left( \mathbf{x}_i - \sum_{j=1}^k z_{ij} \mathbf{v}_j \right)^\top \left( \mathbf{x}_i - \sum_{j=1}^k z_{ij} \mathbf{v}_j \right) \\ &= \mathbf{x}_i^\top \mathbf{x}_i - 2 \sum_{j=1}^k z_{ij} \mathbf{v}_j^\top \mathbf{x}_i + \sum_{j=1}^k z_{ij}^2 \mathbf{v}_j^\top \mathbf{v}_j \end{aligned}$$

Since all  $\mathbf{v}_j$  vectors are orthonormal,  $\mathbf{v}_j^\top \mathbf{v}_j = 1$ , and  $z_{ij} = \mathbf{v}_j^\top \mathbf{x}_i$ , the above results in:

$$\mathbf{x}_i^\top \mathbf{x}_i - 2 \sum_{j=1}^k z_{ij}^2 + \sum_{j=1}^k z_{ij}^2 = \mathbf{x}_i^\top \mathbf{x}_i - \sum_{j=1}^k z_{ij}^2$$

$$= \mathbf{x}_i^T \mathbf{x}_i - \sum_{j=1}^k (\mathbf{x}_i^T \mathbf{v}_j)^T (\mathbf{x}_i^T \mathbf{v}_j) = \mathbf{x}_i^T \mathbf{x}_i - \sum_{j=1}^k \mathbf{v}_j^T \mathbf{x}_i \mathbf{x}_i^T \mathbf{v}_j$$

- b. We know that  $\Sigma$ , the correlation matrix of  $\mathbf{X}$ , can be written as  $\frac{1}{n} \sum_{j=1}^k \mathbf{x}_i \mathbf{x}_i^T$ . Therefore,  $\frac{1}{n} \sum_{j=1}^k \mathbf{v}_j^T \mathbf{x}_i \mathbf{x}_i^T \mathbf{v}_j$  can be written as  $\frac{1}{n} \sum_{j=1}^k \mathbf{v}_j^T \Sigma \mathbf{v}_j$ . Therefore,

$$J_k = \frac{1}{n} \sum_{i=1}^n \mathbf{x}_i^T \mathbf{x}_i - \sum_{j=1}^k \mathbf{v}_j^T \Sigma \mathbf{v}_j$$

. However, since  $\mathbf{v}_j^T \Sigma \mathbf{v}_j = \lambda_j \mathbf{v}_j^T \mathbf{v}_j$ , and  $\mathbf{v}_j^T \mathbf{v}_j = 1$ , the above can be expressed as:

$$J_k = \frac{1}{n} \sum_{i=1}^n \mathbf{x}_i^T \mathbf{x}_i - \sum_{j=1}^k \lambda_j$$

- c. From above, we know that  $J_k = \frac{1}{n} \sum_{i=1}^n \mathbf{x}_i^T \mathbf{x}_i - \sum_{j=1}^k \lambda_j$ , and that  $J_d = 0 = \frac{1}{n} \sum_{i=1}^n \mathbf{x}_i^T \mathbf{x}_i - \sum_{j=1}^d \lambda_j$ . Therefore,

$$J_k - J_d = \sum_{j=1}^d \lambda_j - \sum_{j=1}^k \lambda_j$$

Note that  $J_d = 0$  and  $\sum_{j=1}^d \lambda_j = \sum_{j=1}^k \lambda_j + \sum_{j=k+1}^d \lambda_j$ . The above therefore yields:

$$J_k = \sum_{j=1}^k \lambda_j + \sum_{j=k+1}^d \lambda_j - \sum_{j=1}^k \lambda_j = \sum_{j=k+1}^d \lambda_j$$

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**2 ( $\ell_1$ -Regularization)** Consider the  $\ell_1$  norm of a vector  $\mathbf{x} \in \mathbb{R}^n$ :

$$\|\mathbf{x}\|_1 = \sum_i |\mathbf{x}_i|.$$

Draw the norm-ball  $B_k = \{\mathbf{x} : \|\mathbf{x}\|_1 \leq k\}$  for  $k = 1$ . On the same graph, draw the Euclidean norm-ball  $A_k = \{\mathbf{x} : \|\mathbf{x}\|_2 \leq k\}$  for  $k = 1$  behind the first plot. (Do not need to write any code, draw the graph by hand).

Show that the optimization problem

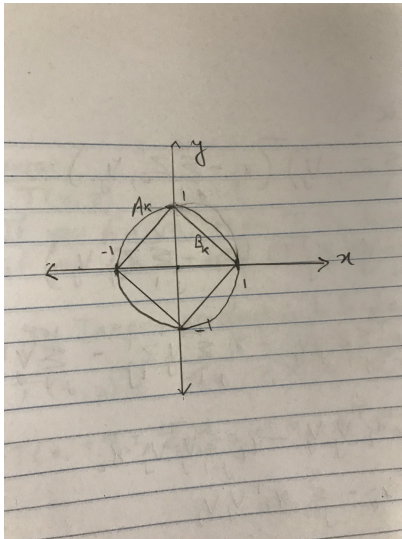
$$\begin{aligned} &\text{minimize: } f(\mathbf{x}) \\ &\text{subj. to: } \|\mathbf{x}\|_p \leq k \end{aligned}$$

is equivalent to

$$\text{minimize: } f(\mathbf{x}) + \lambda \|\mathbf{x}\|_p$$

(hint: create the Lagrangian). With this knowledge, and the plots given above, argue why using  $\ell_1$  regularization (adding a  $\lambda \|\mathbf{x}\|_1$  term to the objective) will give sparser solutions than using  $\ell_2$  regularization for suitably large  $\lambda$ .

The graphs are drawn below:



Minimising  $f(\mathbf{x})$  with constraints  $\|\mathbf{x}\|_p \leq k$  can be solved with Lagrangian Multipliers. Using Lagrangian Multipliers leaves us with the optimisation problem  $L(\mathbf{x}, \lambda) = f(\mathbf{x}) + \lambda(\|\mathbf{x}\|_p - k)$ . In order to create the Lagrangian, we can take the partial derivatives with respect to  $\mathbf{x}$  and  $\lambda$ . We therefore get,

$$\frac{\partial}{\partial \mathbf{x}} L(\mathbf{x}, \lambda) = f'(\mathbf{x}) + \lambda$$

And

$$\frac{\partial}{\partial \lambda} L(\mathbf{x}, \lambda) = \|\mathbf{x}\|_p - k$$

We can see that the terms of the Lagrangian governing the optimal values of  $\mathbf{x}$ , do not depend on  $k$ . This optimisation problem is therefore equivalent to

$$\text{minimise: } f(\mathbf{x}) + \|\mathbf{x}\|_p$$

Based on what was discussed in class, the optimal solution occurs at the point where a contour line of  $f(\mathbf{x})$  meets the constraint. With the  $\ell_1$  norm, we have 'sharper' constraint, which makes it more likely for a contour line of  $f(\mathbf{x})$  to intersect at one of the corners of the norm-ball. With  $\ell_2$  regularisation, a contour of  $f(\mathbf{x})$  is less likely to intersect with the norm-ball along the axes. In fact, there are no 'corners' for the contour to intersect with. Since intersection on the axes, i.e., the corners of the  $\ell_1$  norm ball correspond to sparser solutions (on the axes, at least one dimension is nil), the  $\ell_1$  norm usually yields sparser solutions. Of course, the  $\lambda$  value needs to be suitably large in order to push the parameters values close to 0.

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**Extra Credit (Lasso)** Show that placing an equal zero-mean Laplace prior on each element of the weights  $\theta$  of a model is equivalent to  $\ell_1$  regularization in the Maximum-a-Posteriori estimate

$$\text{maximize: } \mathbb{P}(\theta|\mathcal{D}) = \frac{\mathbb{P}(\mathcal{D}|\theta)\mathbb{P}(\theta)}{\mathbb{P}(\mathcal{D})}.$$

Note the form of the Laplace distribution is

$$\text{Lap}(x|\mu, b) = \frac{1}{2b} \exp\left(-\frac{|x - \mu|}{b}\right)$$

where  $\mu$  is the location parameter and  $b > 0$  controls the variance. Draw (by hand) and compare the density  $\text{Lap}(x|0, 1)$  and the standard normal  $\mathcal{N}(x|0, 1)$  and suggest why this would lead to sparser solutions than a Gaussian prior on each elements of the weights (which correspond to  $\ell_2$  regularization).

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