

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) Consider that  $||\vec{v}||^2 = \vec{v}^\top \vec{v}$ , we can apply the same premise here;

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

(b) Lets begin a key statement that will prove useful,  $\sum \mathbf{x}_i \mathbf{x}_i^\top = \mathbf{\Sigma}$ , we know that the reconstruction error is;

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

(c) We saw in part b that

$$J_d = \sum_{j=1}^d \lambda_j = \frac{1}{n} \left( \sum_{i=1}^n \mathbf{x}_i^\top \mathbf{x}_i \right) - \sum_{j=1}^d \lambda_j$$

We want to find out how much error is introduced for a specific value,  $J_k$ . The expression for this, after partitioning the sum as suggested, in terms of  $d$  will be,

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

This is simply  $J_k = \sum_{j=k+1}^d \lambda_j$  because  $J_d = 0$

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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

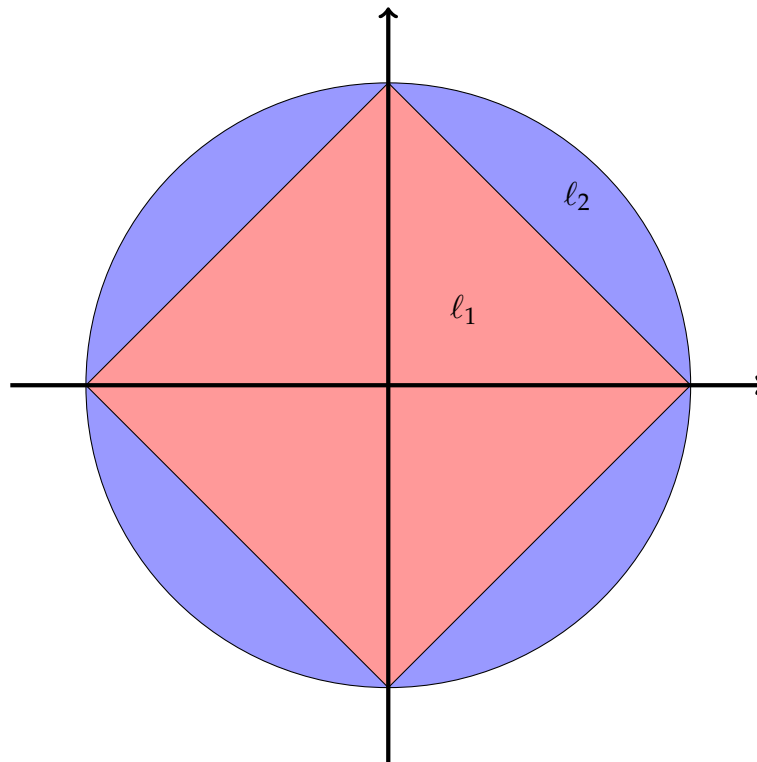
$$\begin{aligned} &\text{minimize: } f(\mathbf{x}) \\ &\text{subj. to: } \|\mathbf{x}\|_p \leq k \end{aligned}$$

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  $\lambda$ .

(a) The desired norm balls



(b) The Lagrange multiplier is defined as  $\mathcal{L}(f(x), g(x)) = f(x) - \lambda g(x)$  Where  $g(x)$  is

our constraint, as such we see that for this problem;

$$\mathcal{L}(f(x), g(x)) = f(x) + \lambda(\|\mathbf{x}\|_p - k)$$

We know that minimizing the Lagrangian is equivalent to minimizing the function  $f(x)$  subject to the constraint and that  $\lambda k$  is not dependent at all upon  $x$ , so we can throw this term away;

$$\text{minimize } \mathcal{L}(f(x), g(x)) = \text{minimize } f(x) + \lambda(\|\mathbf{x}\|_p)$$

Since we have constructed this from a function and constraint given, we can say that the two statements above are equivalent.

- (c) We saw in class that an advantage of the  $\ell_1$  norm is its preference for zeros, and we can think of an optimal solution as residing on a vertex of the norm ball. The  $\ell_1$  norm will have sparser solutions than the  $\ell_2$  norm because it has fewer vertices, and thus fewer optimal solutions.

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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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