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  $\mathbf{v}_i^{\top} \mathbf{\Sigma} \mathbf{v}_j = \lambda_j \mathbf{v}_i^{\top} \mathbf{v}_j = \lambda_j$ .

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

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

Show that the optimization problem

minimize:  $f(\mathbf{x})$  subj. to:  $\|\mathbf{x}\|_p \le k$ 

is equivalent to

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

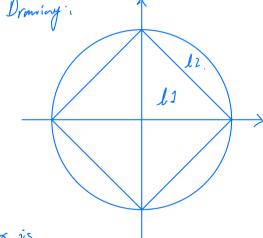
(hint: create the Lagrangian). With this knowledge, and the plots given above, argue why using  $\ell_1$  regularization (adding a  $\lambda \|\mathbf{x}\|_1$  term to the objective) will give sparser solutions than using  $\ell_2$  regularization for suitably large  $\lambda$ .

We know the optimization problem: minimize: f(x) subj. to  $||x||_p \le |x|$ is equivalent to  $\inf_{x} \sup_{\lambda \ge 0} \int_{-\infty}^{\infty} (x, \lambda)^{-1} \inf_{x \ge 0} \sup_{x \ge 0} f(x) + \lambda(||x||_p - |x|).$ 

In its dual we can flip the inf and sup:

sup inf f(x) +  $\lambda(||x||p-|x|) = \sup_{\lambda \neq 0} g(\lambda)$ .

Since the minimizing value of  $f(x) + \lambda(1|x||p-k)$  over x is



equivalent to the minimizing value of  $f(x) + \lambda ||x||p$ , we know that optimizing x will solve minimize:  $f(x) + \lambda ||x||p$  for some suitable value of  $\lambda \geq 0$ . Looking at the alraming, we can consider  $l_1$  regularization as project the autual optimal solution of the problem onto some suitably sized  $l_1$  norm ball. Since the  $l_1$  ball has sharper edges, the probability of landing an our edge and not on the face is infinitely larger than the  $l_2$  ball. This is due to the rotation invariance of the  $l_2$  that clearly aloes it hold for the  $l_1$  ball. Generalizing to higher dimensions, we can see that the  $l_1$  penalty will encourage more weights to be zero compared to the  $l_2$  ball, which is what we want.

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

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

Note the form of the Laplace distribution is

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

where  $\mu$  is the location parameter and b>0 controls the variance. Draw (by hand) and compare the density 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).

We know the Maximum-a-Posteriori problem maximize:  $P(\theta|D) = \frac{P(D|\theta)P(\theta)}{P(D)}$  is equivalent to maximize log  $P(\theta|D) = \frac{P(D|\theta)P(\theta)}{P(D)}$  is equivalent for maximize log  $P(\theta|D) = \log P(D|\theta) + \log P(\theta) - \log P(D)$ .

Since P(D) is a constant not dependent on  $\theta$ , we can drop that term from the problem and flip into a minimization problem, giving animimize:  $-\log P(D|\theta) - \log P(\theta)$ . Given a prior  $\theta : \sim Lap(\theta,b)$ ,  $-\log P(\theta) = -\log T \exp(-\frac{|\theta_0|}{b}) + Z = \frac{1}{b} \sum_{i=0}^{b} |\theta_i| + Z = \lambda ||\theta|| + Z$ .

It following that our original problem is equivalent to minimize:  $-\log P(D|\theta) + \lambda ||\theta||$ , or a regularized maximum likelihood estimate, on desired. Note the plots of the Standard Normal and Laplace Densites.

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We can see that lop(0,1) will place much more mass at x=0. It follows that when we use a loplace prior instead of a Gamssian prior on our weights, (eq. i.i.) our weights will be more 'arranged to be exactly zero, forcing sparslay.