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

(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 can prove this algebraically

$$\left\| x_i - \sum_{j=1}^k z_{ij} v_j \right\|^2 = \left(x_i - \sum_{j=1}^k z_{ij} v_j \right)^T \left(x_i - \sum_{j=1}^k z_{ij} v_j \right)$$

$$= x_i^T x_i - \left(\sum_{j=1}^k z_{ij} v_j\right)^T x_i - x_i^T \left(\sum_{j=1}^k z_{ij} v_j\right) + \left(\sum_{j=1}^k z_{ij} v_j\right)^T \left(\sum_{j=1}^k z_{ij} v_j\right)$$

$$= x_i^T x_i - 2 \sum_{j=1}^k z_{ij} v_j^T x_i + \sum_{j=1}^k \sum_{i=1}^k z_{ij} v_j^T z_{ij} v_j \quad \text{(bringing } x_i^T \text{ into sum)}$$

$$= x_i^T x_i - 2 \sum_{j=1}^k z_{ij} v_j^T x_i + \sum_{j=1}^k v_j^T \left(\sum_{i=1}^k z_{ij} z_{ij} \right) v_j$$

$$= x_i^T x_i - 2 \sum_{j=1}^k z_{ij} v_j^T x_i + \sum_{j=1}^k v_j^T x_i x_i^T v_j \quad \text{(since } v_i^T v_j = 1 \text{ iff } i = j\text{)}$$

$$= x_i^T x_i - 2 \sum_{j=1}^k z_{ij} v_j^T x_i + \sum_{j=1}^k v_j^T x_i x_i^T v_j \quad \text{(since } z_{ij} \in \mathbb{R} \text{)}$$

$$= x_{i}^{T} x_{i} - \sum_{j=1}^{k} v_{j}^{T} x_{i} x_{i}^{T} v_{j},$$

(b) By definition

$$J_{k} = \frac{1}{n} \sum_{i=1}^{n} \left(x_{i}^{T} x_{i} - \sum_{j=1}^{k} v_{j}^{T} x_{i} x_{i}^{T} v_{j} \right)$$

$$= \frac{1}{n} \sum_{i=1}^{n} x_i^T x_i - \sum_{j=1}^{k} v_j^T \left(\frac{1}{n} \sum_{i=1}^{n} x_i x_i^T \right) v_j$$

$$= \frac{1}{n} \sum_{i=1}^{n} x_i^T x_i - \sum_{j=1}^{k} \sum_{i=1}^{n} v_j$$

$$= \frac{1}{n} \sum_{i=1}^{n} x_i^T x_i - \sum_{j=1}^{k} \lambda_j$$

as desired.

(c) Since
$$J_d = 0$$
 we know $\sum_{j=1}^d \lambda_j = \frac{1}{n} \sum_{i=1}^n x_i^T x_i$. Then

$$J_k = \frac{1}{n} \sum_{i=1}^n x_i^T x_i - \sum_{j=1}^d \lambda_j + \sum_{j=k+1}^d \lambda_j = \sum_{j=k+1}^d \lambda_j.$$

3

2 (ℓ_1 -Regularization) Consider the ℓ_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 ℓ_1 regularization (adding a $\lambda \|\mathbf{x}\|_1$ term to the objective) will give sparser solutions than using ℓ_2 regularization for suitably large λ .

Though not included in my solution we can think of the balls, respectively a square and circle on the xy plane for respective L1 and L2 norm. Using these We know the optimization problem

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

is equivalent to

$$\inf_{x} \sup_{\lambda \ge 0} L(x,\lambda) = \inf_{x} \sup_{\lambda \ge 0} f(x) + \lambda(\|x\|_p - k).$$

In its dual we can flip the inf and sup.

$$\sup_{\lambda \ge 0} \inf_{x} f(x) + \lambda(\|x\|_{p} - k) = \sup_{\lambda \ge 0} g(\lambda)$$

Since the minimizing value of $f(x) + \lambda(\|x\|_p - k)$ over x is equivalent to the minimizing value of $f(x) + \lambda \|x\|_p$ (- λk doesn't depend on x), we know the optimizing x will solve

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

for some suitable value of $\lambda \geq 0$. Looking at the plot and this result, we can consider ℓ_1 regularization as projecting the actual optimal solution of your problem onto some suitably sized ℓ_1 norm ball. Since the ℓ_1 ball has sharper edges, the probability of landing on an edge and not on the face (where both elements of the vector are nonzero) is infinitely larger than the ℓ_2 ball. This is due to the rotation invariance of the ℓ_2 that certainly doesn't 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, as desired.

Extra Credit (Lasso) Show that placing an equal zero-mean Laplace prior on each element of the weights θ of a model is equivelent to ℓ_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 μ 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 ℓ_2 regularization).