Exercises

Computational Intelligence Lab

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Problem 1 (PCA Theory):

- 1. (a) $\bar{\mathbf{X}} = \mathbf{X} \mathbf{M}$
 - (b) $\mathbf{\Sigma} = \frac{1}{N} \bar{\mathbf{X}} \bar{\mathbf{X}}^{\top} \in \mathbb{R}^{D \times D}$
 - (c) $\Sigma = \mathbf{U} \mathbf{\Lambda} \mathbf{U}^T$. In the sequel we assume that $\mathbf{\Lambda} = diag(\lambda_1, \dots, \lambda_D)$, where $\lambda_1 \geq \dots \geq \lambda_D \geq 0$. The eigenvalues are positive because Σ is symmetric. Further, the eigenvector matrix \mathbf{U} can be written as $\mathbf{U} = [u_1, \dots, u_D]$, where $u_i \in \mathbb{R}^D$ are unit eigenvectors (i.e. $\|u_i\|_2 = 1$) represented as column vectors.
 - (d) $\bar{\mathbf{Z}}_K = \mathbf{U}_K^{\top} \bar{\mathbf{X}}$. Here, we have \mathbf{U}_K is given by the first K columns of \mathbf{U} , i.e. $\mathbf{U}_K = [u_1, \dots, u_K]$.
 - (e) $\tilde{\mathbf{X}} = \mathbf{U}_K \bar{\mathbf{Z}}_K$
 - (f) We have that $\tilde{\mathbf{X}} = \mathbf{U}_K \mathbf{U}_K^{\top} \bar{\mathbf{X}}$. The reconstruction error is :

$$\mathsf{err} = \frac{1}{N} \sum_{i=1}^N \|\tilde{x}_i - \bar{x}_i\|_2^2 = \|\tilde{\mathbf{X}} - \bar{\mathbf{X}}\|_F^2 = \|(\mathbf{U}_K \mathbf{U}_K^\top - \mathbf{I}_d)\bar{\mathbf{X}}\|_F^2$$

where $||A||_F = \sqrt{\operatorname{trace}(AA^\top)} = \sqrt{\sum_i \sigma_i^2}$ is the Frobenius norm of matrix A and σ_i are its singular values (the same as eigenvalues if A is symmetric). Thus,

$$\begin{split} & \operatorname{err} = \frac{1}{N} \operatorname{trace}((\mathbf{U}_K \mathbf{U}_K^\top - \mathbf{I}_d) \bar{\mathbf{X}} \bar{\mathbf{X}}^\top (\mathbf{U}_K \mathbf{U}_K^\top - \mathbf{I}_d)^\top) \\ & = \operatorname{trace}((\mathbf{U}_K \mathbf{U}_K^\top - \mathbf{I}_d) \mathbf{\Sigma} (\mathbf{U}_K \mathbf{U}_K^\top - \mathbf{I}_d)) \\ & = \operatorname{trace}((\mathbf{U}_K \mathbf{U}_K^\top - \mathbf{I}_d) \mathbf{U} \boldsymbol{\Lambda} \mathbf{U}^\top (\mathbf{U}_K \mathbf{U}_K^\top - \mathbf{I}_d)) \\ & = \operatorname{trace}((\mathbf{U}_K \mathbf{U}_K^\top \mathbf{U} - \mathbf{U}) \boldsymbol{\Lambda} (\mathbf{U}^\top \mathbf{U}_K \mathbf{U}_K^\top - \mathbf{U}^\top)) \\ & = \operatorname{trace}(([\mathbf{U}_K; \mathbf{0}] - \mathbf{U}) \boldsymbol{\Lambda} ([\mathbf{U}_K; \mathbf{0}] - \mathbf{U})^\top) \\ & = \operatorname{trace}(\sum_{i=K+1}^D \lambda_i u_i u_i^\top) \\ & = \sum_{i=K+1}^D \lambda_i \cdot \operatorname{trace}(u_i u_i^\top) \\ & = \sum_{i=K+1}^D \lambda_i \end{split}$$

where we used the fact that $\operatorname{trace}(u_i u_i^{\top}) = ||u_i||_2^2 = 1$.

- 2. (a) Intrinsic dimensionality: high No knee in eigenvalue spectrum
 - (b) No, the approximation error is the sum of the discarded eigenvalues and λ_{100} is still large.
 - (c) D = 100 (no reduction)

3.

1.
$$\begin{bmatrix} 1 & 0 \\ 0 & 2 \end{bmatrix}$$
 Answer: (B)

2.
$$\begin{bmatrix} 1 & -0.5 \\ -0.5 & 1 \end{bmatrix}$$
 Answer: (E)

3.
$$\begin{bmatrix} 1 & 0 \\ 0 & 1 \end{bmatrix}$$
 Answer: (C)

- 4. (a) We would like to decouple the dimensions/measurements in the transformed dataset, i.e. we would like to have uncorrelated dimensions.
 - (b) Consider $\mathbf{Z} = \mathbf{A}^{\top} \mathbf{X}$. Let $\bar{\mathbf{x}}$ be the mean of the dataset \mathbf{X} . We write $\mathbf{M}_{\mathbf{X}} = [\bar{\mathbf{x}},...,\bar{\mathbf{x}}]$, correspondingly, $\mathbf{M}_{\mathbf{Z}} = A^{\top} \mathbf{M}_{\mathbf{X}}$. We can write the covariance matrix of \mathbf{X} as $\mathbf{\Sigma}_{\mathbf{X}} = (\mathbf{X} \mathbf{M}_{\mathbf{X}})(\mathbf{X} \mathbf{M}_{\mathbf{X}})^{\top}$. The covariance of \mathbf{Z} is then given by:

$$\begin{split} \boldsymbol{\Sigma}_{\mathbf{Z}} &= & (\mathbf{Z} - \mathbf{M}_{\mathbf{Z}})(\mathbf{Z} - \mathbf{M}_{\mathbf{Z}})^{\top} \\ &= & (\mathbf{A}^{\top}\mathbf{X} - \mathbf{M}_{\mathbf{Z}})(\mathbf{A}^{\top}\mathbf{X} - \mathbf{M}_{\mathbf{Z}})^{\top} \\ &= & (\mathbf{A}^{\top}\mathbf{X} - \mathbf{A}^{\top}\mathbf{M}_{\mathbf{X}})(\mathbf{A}^{\top}\mathbf{X} - \mathbf{A}^{\top}\mathbf{M}_{\mathbf{X}})^{\top} \\ &= & \mathbf{A}^{\top}(\mathbf{X} - \mathbf{M}_{\mathbf{X}})(\mathbf{A}^{\top}(\mathbf{X} - \mathbf{M}_{\mathbf{X}}))^{\top} \\ &= & \mathbf{A}^{\top}(\mathbf{X} - \mathbf{M}_{\mathbf{X}})(\mathbf{X} - \mathbf{M}_{\mathbf{X}})^{\top}\mathbf{A} \\ &= & \mathbf{A}^{\top}\boldsymbol{\Sigma}_{\mathbf{X}}\mathbf{A} \end{split}$$

(c) If we use
$$\mathbf{A}=\mathbf{U}$$
, we obtain:
$$\begin{aligned} \boldsymbol{\Sigma}_{\mathbf{Z}} &=& \mathbf{A}^{\top}\boldsymbol{\Sigma}_{\mathbf{X}}\mathbf{A}\\ &=& \mathbf{U}^{\top}\boldsymbol{\Sigma}_{\mathbf{X}}\mathbf{U}\\ &=& \mathbf{U}^{\top}\mathbf{U}\boldsymbol{\Lambda}\mathbf{U}^{\top}\mathbf{U}\\ &=& \mathbf{U}^{-1}\mathbf{U}\boldsymbol{\Lambda}\mathbf{U}^{-1}\mathbf{U}\\ &=& \mathbf{I}\boldsymbol{\Lambda}\mathbf{I}\\ &=& \boldsymbol{\Lambda} \end{aligned}$$

We see that the covariance matrix of \mathbf{Z} becomes the diagonal eigenvalue matrix $\boldsymbol{\Lambda}$: Choosing the eigenvectors associated with the highest eigenvalues results in capturing high variances in the transformed dataset.