

CS736 Assignment 2

Adish Shah and Padakanti Akshay Kiran

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1 Problem 1

1.1 d) Initial Pointsets

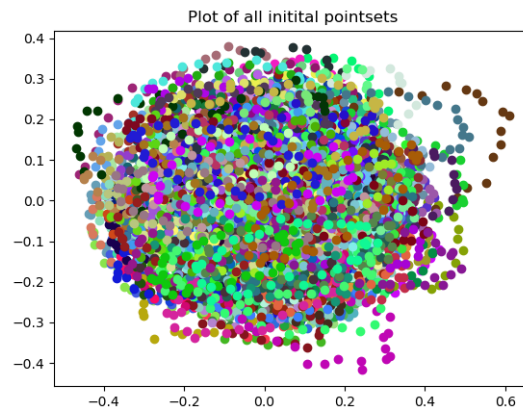


Figure 1: Scatter plot

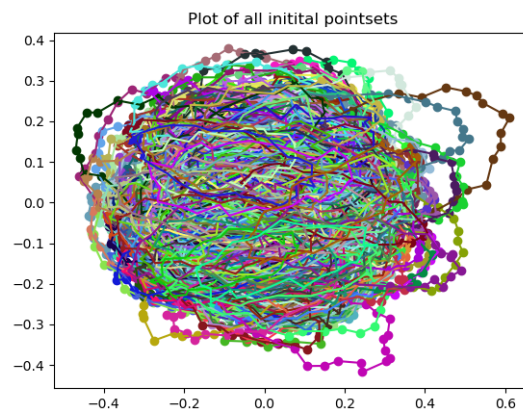


Figure 2: Each pointset represented by polylines

1.2 e)

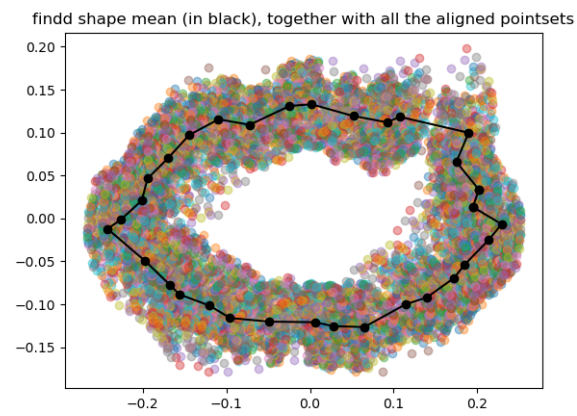


Figure 3: Result based on code11

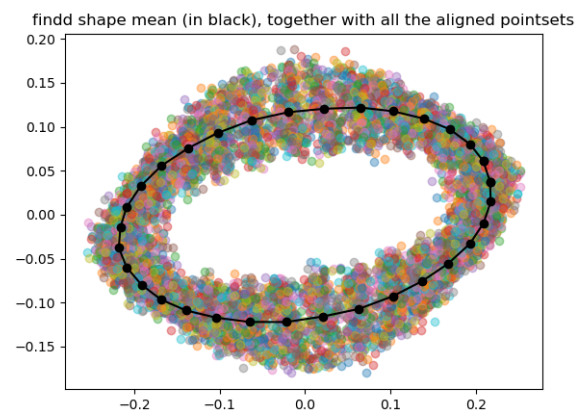


Figure 4: Result based on code22

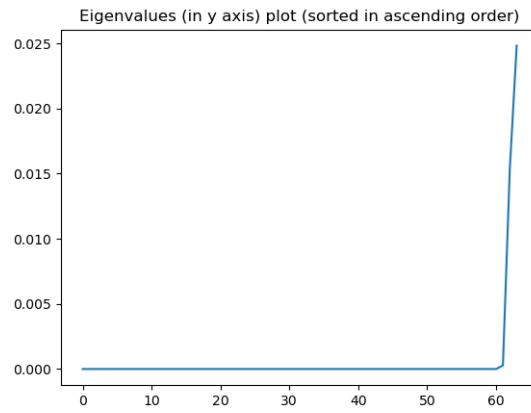
1.3 f)

Figure 5: Result based on code11. First two eigenvalues are significantly greater than other eigenvalues which are close to 0

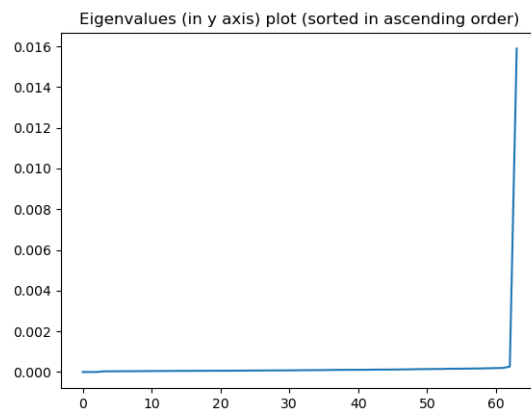


Figure 6: Result based on code22

1.4 g)

1.4.1 Results based on code11

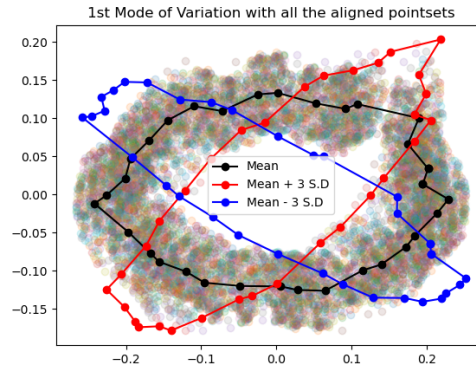


Figure 7: First Mode of variation

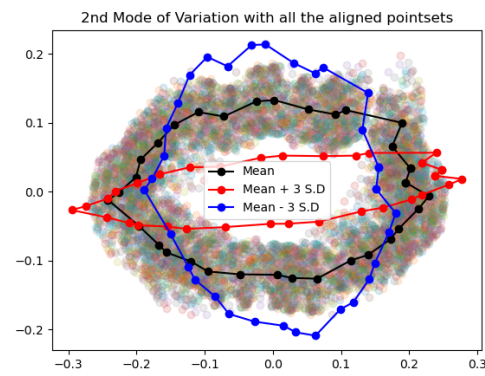


Figure 8: Second Mode of variation

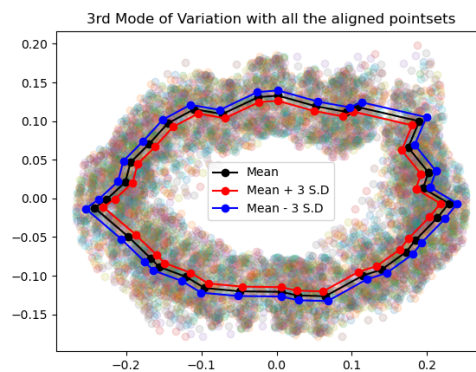


Figure 9: Third mode of variation

1.4.2 Results based on code22

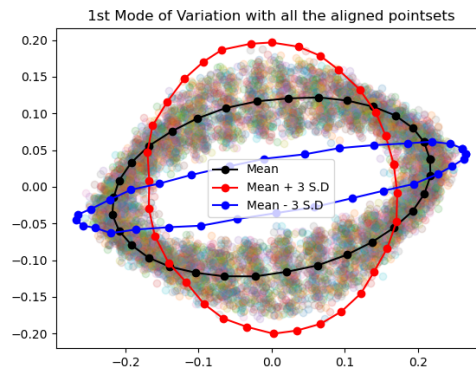


Figure 10: First Mode of variation

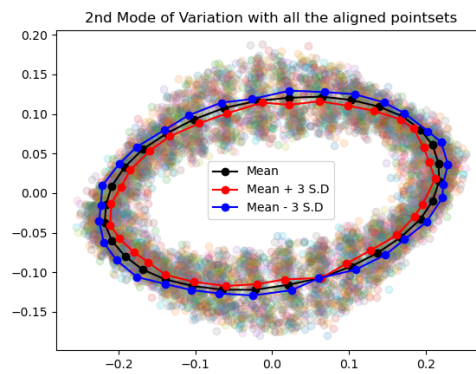


Figure 11: Second Mode of variation

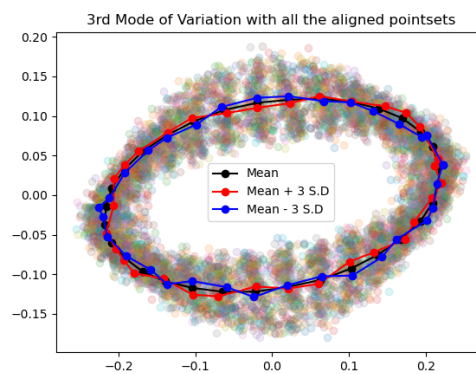


Figure 12: Third mode of variation

2 Problem 2

2.1 d) Initial Pointsets

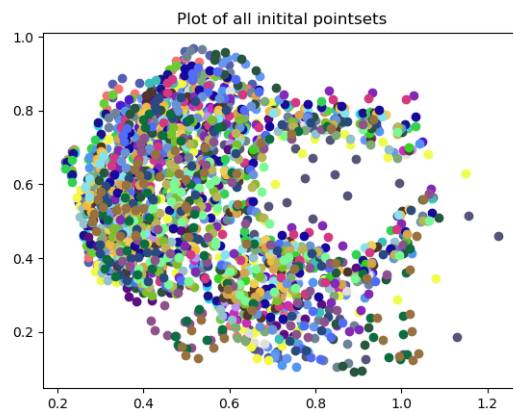


Figure 13: Scatter plot

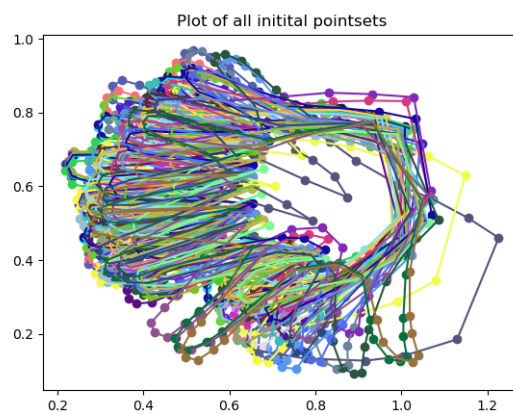


Figure 14: Each pointset represented by polylines

2.2 e)

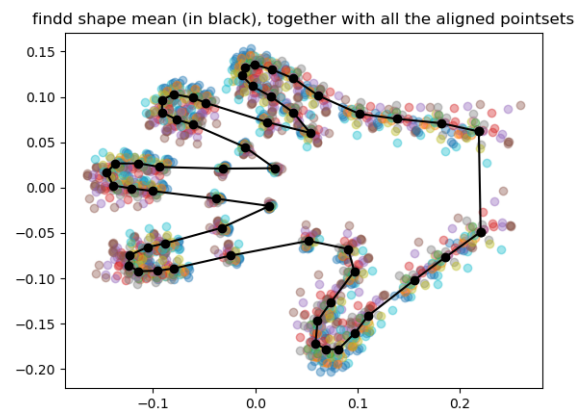


Figure 15: Result based on code11

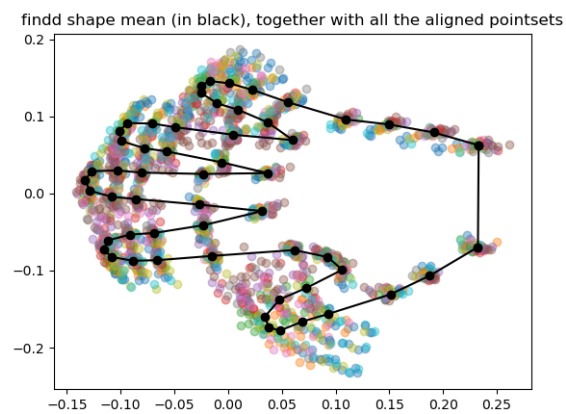


Figure 16: Result based on code22

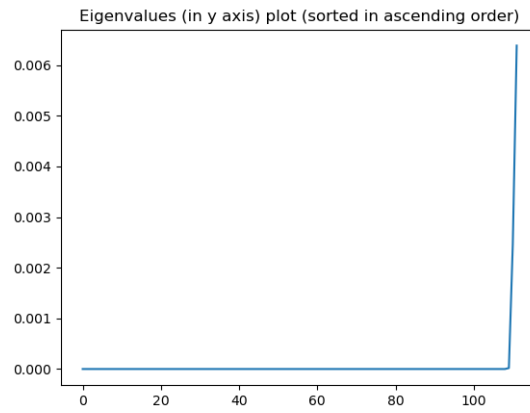
2.3 f)

Figure 17: Result based on code11. First two eigenvalues are significantly greater than other eigenvalues which are close to 0

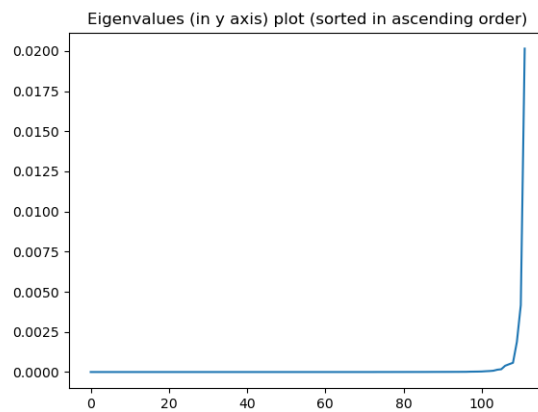


Figure 18: Result based on code22

2.4 g)

2.4.1 Results based on code11

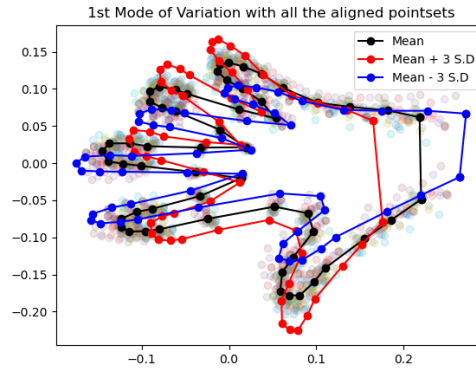


Figure 19: First Mode of variation captures

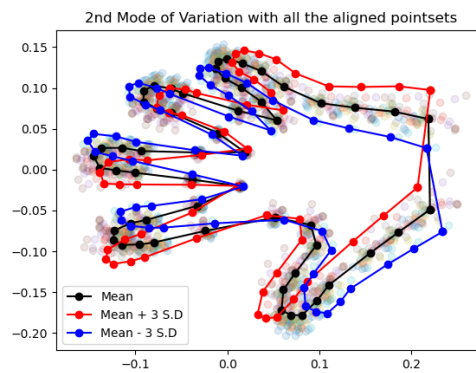


Figure 20: Second Mode of variation

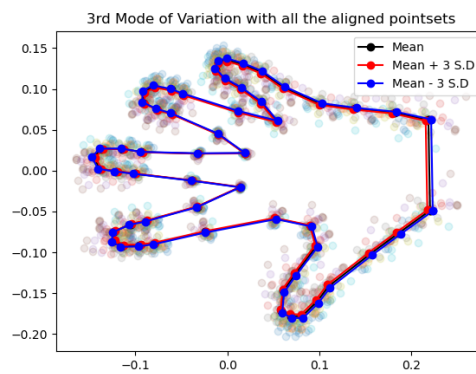


Figure 21: Third mode of variation

2.4.2 Results based on code22

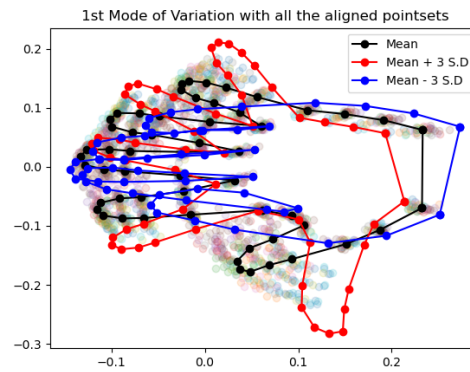


Figure 22: First Mode of variation

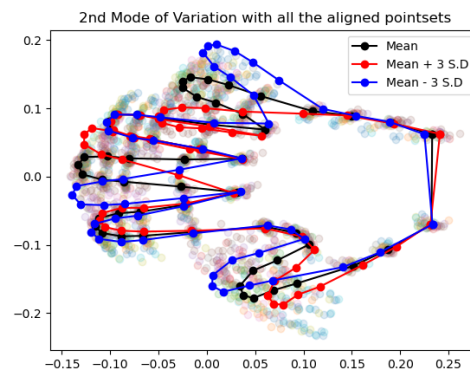


Figure 23: Second Mode of variation

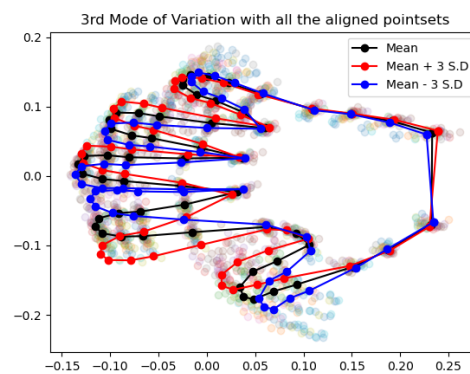


Figure 24: Third mode of variation

3 Problem 3

3.1 Procrustes Distance

The simplest way to define proximity is just the magnitude of the Euclidean norm of the difference between the two shape vectors in \mathbb{R}^{dN} which is the square of the Procrustes distance between the shapes. Procrustes distance between z_1 and z_2 (resulting from aligning z_2 to z_1 by similarity transformations) is

$$\min_{\theta, T, s} \sum_{n=1, \dots, N} \|z_{1n} - sM_{\theta}z_{2n} - T\|$$

where s is the scaling factor, M_{θ} is the rotation matrix of dimensions $d \times d$ and T represents a translation in \mathbb{R}^d .

3.2 K - means algorithm in the Shape space

We apply the k-means algorithm to the points X_1, X_2, \dots, X_n by using the Procrustes distance and Procrustes Mean.

3.3 Objective Function

$$W(\mathcal{C}, [Z_1], \dots, [Z_k]) = \sum_{i=1}^k \sum_{l \in C_i} \text{Procrustes_distance}([X_l], [Z_i])$$

3.4 Algorithm for clustering

1. **Initialisation condition** The starting centroid vector can be assigned in different ways: at random, using prior information, etc. In our application we will choose it at random.
2. Given a vector of shapes $Z = ([Z_1], \dots, [Z_k])$ $[Z_i] \in \sum_3^n$ $i = 1, \dots, k$ we minimize with respect to $\mathcal{C} = (C_1, \dots, C_k)$ assigning each shape $([X_1], \dots, [X_n])$ to the class whose centroid has the minimum Procrustes distance to it.
3. Given \mathcal{C} , we minimize with respect to the vector shapes Z , taking $Z = ([\hat{\mu}_1], \dots, [\hat{\mu}_k])$, and $[\hat{\mu}_i]$ $i = 1, \dots, k$, the Procrustes mean of shapes in C_i . This is calculated using the algorithm mentioned below.
4. Repeat steps 2 and 3 till convergence.
5. **Stopping condition** When the value of the objective function doesn't decrease further, we stop.

3.5 Algorithm to find the Procrustes Means

1. Center the configurations to remove location. Let $X_i, i = 1, \dots, n$ be the configurations. Initially let

$$X_i^P = X_i, i = 1, \dots, n$$

2. Calculate $G = \frac{1}{n} \sum_{i=1}^{n-1} \sum_{j=i+1}^n \|X_i^P - X_j^P\|^2$ For the i th configuration let

$$\bar{X}_{(i)} = \frac{1}{n-1} \sum_{j \neq i} X_j^P$$

We can optimise $\|\bar{X}_{(i)} - X_i^P M\|^2$ over rotations. Set $X_i^P = X_i^P \hat{M}$, where \hat{M} is the optimal rotation matrix. Repeat this for all i . Calculate the new G value. This process is repeated until G can't be reduced further.

3. For the i th configuration calculate $\hat{\beta}_i = \frac{\sum_{k=1}^n \|X_k^P\|^2}{\|X_i^P\|^2}^{1/2} \phi_i$, where ϕ_i is the i th component of the eigenvector ϕ corresponding to the largest eigenvalue of the correlation matrix ϕ of the *vector*(X_i^P). Set $X_i^P = \hat{\beta}_i X_i^P$. Repeat this for all i and calculate the new G value.
4. Repeat steps 2 and 3 until G cannot be reduced further.
- 5.

$$[\hat{\mu}] = \frac{1}{n} \sum_i^n X_i^P$$