

# Skeleton Clustering : A Dimension free Density-Aided Clustering

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# Traditional Clustering Methods

- **k-means clustering:**

- Unable to detect non-convex clusters.
- The center of a non-convex cluster falls outside the cluster itself and may come close to observations from a different cluster.
- In high dimension k-means algorithm may assign all the points to a single cluster.

- **Density Based Clustering:**

- To estimate the underlying PDF and detect clusters based on the PDF.
- The rate of convergence for the density estimates is  $\mathcal{O}_{\mathbb{P}}(n^{-\frac{1}{d+4}})$

- **Hierarchical Clustering:**

- Problem with non-convex clusters persists.
- If any pair of the points in two different clusters lie very close to each other, the two clusters may get merged in this method.

# Skeleton Clustering Framework.

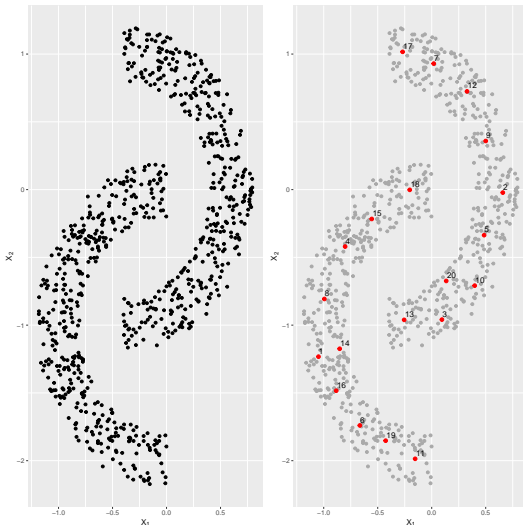
**Input :** Observations  $X_1, X_2, \dots, X_N$ , final number of clusters  $S$ .

- ① **Knot construction** : Perform  $k$ -means clustering with a large number  $k$ ; the centers are the knots.
- ② **Edge construction** : Apply approximate Delaunay triangulation to the knots. Generally we choose  $k = \lfloor \sqrt{n} \rfloor$
- ③ **Edge weights construction** : Add weights to each edge using either Voronoi density, Face density or Tube density similarity measure.
- ④ **Knots segmentation** : Use linkage criterion to segment knots into  $S$  groups based on the edge weights.
- ⑤ **Assignment of labels** : Assign a cluster label to each observation based on which knot group the nearest knot belongs to.

# Knot construction

- Some knots are constructed to give a concise representation of the data structure.
- In practice we use  $k$ -Means to choose  $k = \lfloor \sqrt{n} \rfloor$  knots, where  $n$  is the number of samples.
- Empirically robustness performance with sufficient number of knots.

# Knot Construction

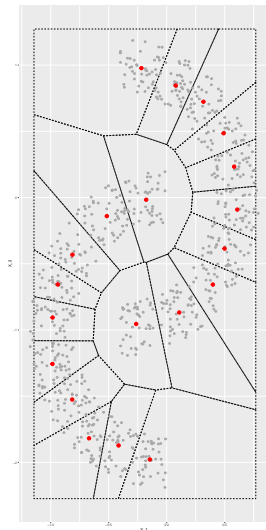


# Edge construction

Let  $c_1, c_2, \dots, c_k$  be the given knots and we use  $\mathcal{C} = \{c_1, c_2, \dots, c_k\}$  to denote their collection of them.

- The Voronoi cell, or Voronoi region,  $\mathbb{C}_j$  associated with a knot  $c_j$  is the set of all points in  $\mathcal{X}$  whose distance to  $c_j$  is the smallest compared to other knots. That is,  $\mathbb{C}_j = \{\mathbf{x} \in \mathcal{X} : d(\mathbf{x}, c_j) \leq d(\mathbf{x}, c_\ell) \forall \ell \neq j\}$  where  $d(\mathbf{x}, \mathbf{y})$  is the usual Euclidean distance.

# Edge Construction

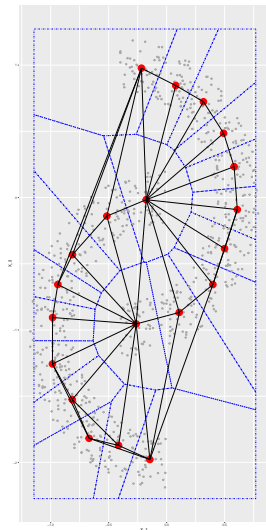


# Edge Construction

- We add an edge between a pair of knots if they are neighbors, with the neighboring condition being that the corresponding Voronoi cells share a common boundary.
- Such resulting graph is the Delaunay Triangulation of the set of knots  $\mathcal{C}$  and we denote it as  $DT(\mathcal{C})$ .
- But in case of high dimensional data, it becomes computationally expensive. Therefore, in practice we approximate the exact Delaunay Triangulation with  $\widehat{DT}(\mathcal{C})$  by examining the 2-nearest knots of the sample data points.



# Edge Construction



# Edge Weight Construction.

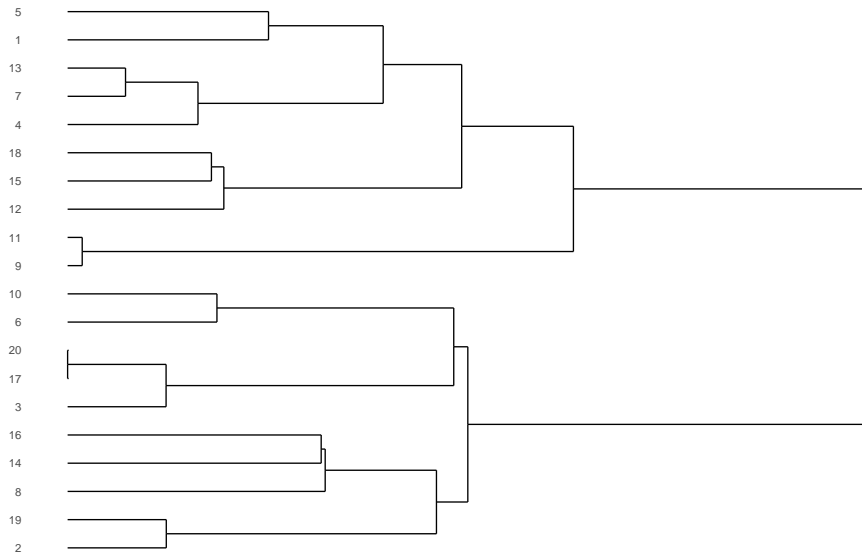
- Given the constructed edges and knots, we assign each edge a weight that represents the similarity between the pair of knots.
- We propose a density aided quantity as edge weights.
- The estimations of the newly proposed similarity measures are reliable even under high-dimensional settings.

# Knot Segmentation.

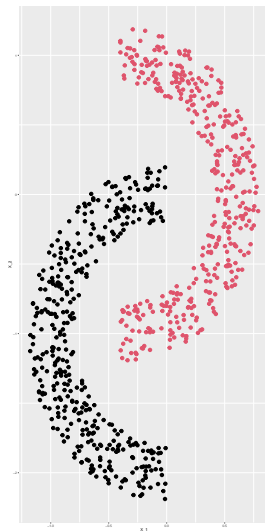
Given the weighted skeleton graph, the next step is to partition the knots into the desired number of final clusters.

- For given similarity measures  $\{s_{ij}\}_{i \neq j}$  where only connected pairs can take nonzero entries and let  $s_{max} = \max_{i \neq j} s_{ij}$ , we define the corresponding distances as  $d_{ij} = 0$  if  $i = j$  and  $d_{ij} = s_{max} - s_{ij}$  otherwise.
- After that, we perform a hierarchical clustering. The choice of the linkage criterion may depend on the underlying geometric structure of the data. In practice, such choice of linkage should be made based on some exploratory understanding of the data structure.

# Knot Segmentation



# Final Clustering Result



# How to choose the final number of Clusters ?

- The number of final clusters  $S$  is an essential parameter for the hierarchical clustering procedure but can be unknown.
- An inferential choice can also be made using the gap statistics, Calinski-Harabasz index etc.

# Some feasible density functions

## Voronoi Density (VD).

- The Voronoi density (VD) measures the similarity between a pair of knots  $(c_j, c_\ell)$  based on the number of observations whose 2-nearest knots are  $c_j$  and  $c_\ell$ .
- Given a metric  $d$  on  $\mathbb{R}^d$ , the 2-Nearest-Neighbor (2-NN) region of a pair of knots  $(c_j, c_\ell)$  is defined as,

•

$$A_{j\ell} = \{\mathbf{x} \in \mathcal{X} : d(\mathbf{x}, c_i) > \max\{d(\mathbf{x}, c_j), d(\mathbf{x}, c_\ell)\} \forall i \neq j, \ell\}$$

- Following the idea of density-based clustering, two knots  $c_j, c_\ell$  belong to the same clusters if they are in a connected high-density region, and we would expect the 2-NN region of  $c_j, c_\ell$  to have a high probability measure. Hence  $\mathbb{P}(A_{j\ell})$  can measure the association between the knots.

# Voronoi Density

- Based on this insight, the Voronoi density measures the edge weight of  $(c_j, c_\ell)$  with,

$$S_{j\ell}^{VD} = \frac{\mathbb{P}(A_{j\ell})}{\|c_j - c_\ell\|}$$

- In practice, we estimate  $S_{j\ell}^{VD}$  by a sample average. Specifically, the numerator  $\mathbb{P}(A_{j\ell})$  is estimated by  $\hat{\mathbb{P}}_n(A_{j\ell}) = \frac{1}{n} \sum_{i=1}^n \mathbf{1}(\mathbf{X}_i \in A_{j\ell})$  and the final estimator for VD is,

$$\hat{S}_{j\ell}^{VD} = \frac{\hat{\mathbb{P}}(A_{j\ell})}{\|c_j - c_\ell\|}$$