

# STATISTICAL DEPTH FUNCTIONS

*In the multivariate and functional setting*

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MASTER'S THESIS

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# ABSTRACT

# DEDICATION

# DECLARATION

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# *Chapter 1*

## INTRODUCTION

### 1.1 Centrality vs Density

### 1.2 Nonparametric procedures

## Chapter 2

# MULTIVARIATE DATA

### 2.1 Depth contours

The following definitions are adapted from Liu et al., [1999](#).

**Definition 2.1.1.** The contour of depth  $t$  is the set  $\{\mathbf{x} \in \mathbb{R}^d : D(\mathbf{x}, F) = t\}$ .

**Definition 2.1.2.** The region enclosed by the contour of depth  $t$  is the set

$$R_F(t) = \{\mathbf{x} \in \mathbb{R}^d : D(\mathbf{x}, F) > t\}. \quad (2.1.1)$$

**Definition 2.1.3.** The  $p$ -th central region is the set

$$C_F(p) = \bigcap_t \{R_F(t) : P_F(R_F(t)) \geq p\}. \quad (2.1.2)$$

**Definition 2.1.4.** The  $p$ -th level contour, or center-outward contour surface, is the set  $Q_F(p) = \partial C_F(p)$ .

### 2.2 Depth-Depth plots

**Definition 2.2.1** (DD plot). Let  $F, G$  be two distributions on  $\mathbb{R}^d$ , and let  $D$  be a depth function. The Depth-Depth plot, also known as the DD plot, of  $F$  and  $G$  is given by

$$\text{DD}(F, G) = \{(D(\mathbf{z}, F), D(\mathbf{z}, G)) : \mathbf{z} \in \mathbb{R}^d\}. \quad (2.2.1)$$

*Remark.* The above definition generalizes naturally to involve more than two distributions on  $\mathbb{R}^d$ .

When the depth function  $D$  only takes values in  $[0, 1]$ , the DD plot is a subset of  $[0, 1]^2$  and hence easily visualized. Clearly when  $F = G$ , the corresponding DD plot is confined to the diagonal  $\{(t, t) : t \in [0, 1]\}$ . However, when  $d \geq 2$  and  $F, G$  are absolutely continuous,  $\text{DD}(F, G)$  has non-zero area (Lebesgue measure) when



$F \neq G$ . Assuming that  $D$  is affine invariant, Liu et al. (1999) propose this area as an affine invariant measure of the discrepancy between  $F$  and  $G$ .

If the distributions  $F, G$  are unknown, we may use data  $\mathcal{D}_F = \{\mathbf{x}_i\}$  and  $\mathcal{D}_G = \{\mathbf{y}_j\}$  where  $\mathbf{x}_1, \dots, \mathbf{x}_n \stackrel{\text{iid}}{\sim} F$  and  $\mathbf{y}_1, \dots, \mathbf{y}_m \stackrel{\text{iid}}{\sim} G$ , then construct empirical distributions  $\hat{F}_n$  and  $\hat{G}_m$ . With this, we may examine the empirical DD plot

$$\text{DD}(\hat{F}_n, \hat{G}_m) = \{(D(\mathbf{z}, \hat{F}_n), D(\mathbf{z}, \hat{G}_m)) : \mathbf{z} \in \mathcal{D}_F \cup \mathcal{D}_G\}. \quad (2.2.2)$$

DD plots can be used as a diagnostic tool to detect differences in location, scale, skewness, and kurtosis between two multivariate distributions (Liu et al., 1999).

1. If the same point  $\mathbf{z}_0$  achieves maximum depths with respect to both distributions  $F$  and  $G$ , this indicates that  $\mathbf{z}_0$  is their common center.
2. Suppose that  $F$  and  $G$  have the same center. If the points in  $\text{DD}(\hat{F}_n, \hat{G}_m)$  arch above the diagonal, i.e. the bulk of points are deeper in  $G$  than in  $F$ , this indicates that  $F$  has a greater spread than  $G$ .

## 2.3 Testing

## 2.4 Classification

The  $k$ -class classification task involves assigning an observation  $\mathbf{x}$  to one of  $k$  populations, described by distributions  $F_i$  for  $1 \leq i \leq k$ . The populations may also be associated with prior probabilities  $\pi_i$ .

**Definition 2.4.1** (Classifier). A classifier is a map  $\hat{\iota}: \mathbb{R}^d \rightarrow \{1, \dots, k\}$ .

**Example 2.4.2** (Bayes classifier). Suppose that the population densities  $f_i$  for each  $1 \leq i \leq k$  are known. The Bayes classifier assigns  $\mathbf{x}$  to the  $\hat{\iota}_B$ -th population where

$$\hat{\iota}_B(\mathbf{x}) = \arg \max_{1 \leq i \leq k} \pi_i f_i(\mathbf{x}). \quad (2.4.1)$$

One way of measuring the performance of a classifier (given the population distributions and their priors) is by measuring its average misclassification rate.

**Definition 2.4.3.** The average misclassification rate of a classifier  $\hat{\iota}$  is given by

$$\Delta(\hat{\iota}) = \sum_{i=1}^k \pi_i P(\hat{\iota}(\mathbf{X}) \neq i \mid \mathbf{X} \sim F_i). \quad (2.4.2)$$

**Proposition 2.4.4.** *The Bayes classifier has the lowest possible average misclassification rate. This is known as the optimal Bayes risk, denoted  $\Delta_B$ .*

The simplest depth based classifier is the maximum depth classifier (Ghosh & Chaudhuri, 2005).

**Example 2.4.5** (Maximum depth classifier). Suppose that the prior probabilities  $\pi_i$  are equal. The maximum depth classifier  $\hat{l}_D$  for a choice of depth function  $D$  is described by

$$\hat{l}_D(\mathbf{x}) = \arg \max_{1 \leq i \leq k} D(\mathbf{x}, F_i). \quad (2.4.3)$$

In practice, instead of having direct access to the population distributions  $F_i$ , we have typically deal with labeled training data

$$\mathcal{D} = \{(\mathbf{x}_{ij}, i)\} \subset \mathbb{R}^d \times \{1, \dots, k\}, \quad (2.4.4)$$

where  $\mathbf{x}_{i1}, \dots, \mathbf{x}_{in_i} \stackrel{\text{iid}}{\sim} F_i$  for each  $1 \leq i \leq k$ . The empirical maximum depth classifier simply replaces the population distributions  $F_i$  with their empirical counterparts  $\hat{F}_i$  determined by  $\mathbf{x}_{i1}, \dots, \mathbf{x}_{in_i}$ . Thus, it is given by

$$\hat{l}_D(\mathbf{x}) = \arg \max_{1 \leq i \leq k} D(\mathbf{x}, \hat{F}_i). \quad (2.4.5)$$

Under certain restrictions, this classifier becomes asymptotically optimal in the following sense.

**Theorem 2.4.6** (Ghosh and Chaudhuri, 2005). *Suppose that the population density functions  $f_i$  are elliptically symmetric, with  $f_i(\mathbf{x}) = g(\mathbf{x} - \boldsymbol{\mu}_i)$  for parameters  $\boldsymbol{\mu}_i$  and a density function  $g$  such that  $g(k\mathbf{x}) \leq g(\mathbf{x})$  for every  $\mathbf{x}$  and  $k > 1$ . Further suppose that the priors on the populations are equal, and the depth function  $D$  is one of HD, SD, MJD, PD. Then,  $\Delta(\hat{l}_D) \rightarrow \Delta_B$  as  $\min\{n_1, \dots, n_k\} \rightarrow \infty$ .*

Note that this result deals with elliptic population densities differing only in location. Relax this assumption, and instead suppose that  $f_i$  are elliptic of the form

$$f_i(\mathbf{x}) = c_i |\Sigma|^{-1/2} h_i((\mathbf{x} - \boldsymbol{\mu}_i)^\top \Sigma^{-1} (\mathbf{x} - \boldsymbol{\mu}_i)) \quad (2.4.6)$$

for strictly decreasing  $h_i$ , and that the depths can be expressed as  $D(\cdot, F_i) = l_i(f_i(\cdot))$  for strictly increasing functions  $l_i$ . It follows that the Bayes decision rule can be reformulated as

$$\pi_i f_i(\mathbf{x}) > \pi_j f_j(\mathbf{x}) \iff D(\mathbf{x}, F_i) > r_{ij}(D(\mathbf{x}, F_j)) \quad (2.4.7)$$

for some real increasing function  $r_{ij}$ . Using this observation, the DD classifier (Li et al., 2012) picks separating functions  $r_{ij}$  which best classify the training data  $\mathcal{D}$ .

**Definition 2.4.7.** The empirical misclassification rate of a classifier  $\hat{l}$ , with respect to data  $\mathcal{D}$ , is given by

$$\hat{\Delta}(\hat{l}) = \sum_{i=1}^k \frac{\pi_i}{n_i} \sum_{j=1}^{n_i} \mathbf{1}(\hat{l}(\mathbf{x}_{ij}) \neq i). \quad (2.4.8)$$

**Definition 2.4.8** (DD classifier). Suppose that  $k = 2$ , that  $D$  is a depth function, and that  $r: [0, 1] \rightarrow [0, 1]$  is an increasing function. The DD classifier  $\hat{l}_{D,r}$  is given by

$$\hat{l}_{D,r}(\mathbf{x}) = \begin{cases} 1, & \text{if } D(\mathbf{x}, F_2) \leq r(D(\mathbf{x}, F_1)), \\ 2, & \text{if } D(\mathbf{x}, F_2) > r(D(\mathbf{x}, F_1)). \end{cases} \quad (2.4.9)$$

The empirical DD classifier  $\hat{l}_{D,\hat{r},\mathcal{D}}$  replaces  $F_i$  by their empirical counterparts  $\hat{F}_i$ . Here, the separating curve  $\hat{r}$  is chosen from a family  $\Gamma$  so as to minimize the empirical misclassification rate, i.e.

$$\hat{r} = \arg \min_{r' \in \Gamma} \hat{\Delta}(\hat{l}_{D,r'}). \quad (2.4.10)$$

*Remark.* The maximum depth classifier  $\hat{l}_D$  is simply the DD classifier  $\hat{l}_{D,\text{id}}$ , where  $\text{id}(x) = x$ .

Li et al. (2012) show that under certain restrictions, the empirical DD classifier is asymptotically equivalent to the Bayes rule. We give one such instance below.

**Lemma 2.4.9.** *Suppose that the following conditions hold.*

1.  $\Gamma$  is the class of polynomial functions on  $[0, 1]$ .
2. The depth functions  $D(\cdot, F_i)$  are continuous.
3. As  $\min\{n_1, n_2\} \rightarrow \infty$ , we have for each  $i \in \mathbb{N}_2$ ,

$$\sup_{\mathbf{z} \in \mathbb{R}^d} |D(\mathbf{z}, \hat{F}_i) - D(\mathbf{z}, F_i)| \xrightarrow{a.s.} 0. \quad (2.4.11)$$

4. For each  $i \in \mathbb{N}_2$ , the distributions  $F_i$  are elliptical and satisfy for all  $\delta \in \mathbb{R}$

$$P(D(\mathbf{Z}, F_i) = \delta \mid \mathbf{Z} \sim F_i) = 0. \quad (2.4.12)$$

Then,  $\Delta(\hat{l}_{D,\hat{r}}) \rightarrow \Delta_B$  as  $\min\{n_1, n_2\} \rightarrow \infty$ .

In all the depth based classifiers we have seen so far, the classification rule depends on the observation  $\mathbf{x}$  only through the depths  $D(\mathbf{x}, F_i)$ . Thus, we are motivated to define the following transformation from  $\mathbb{R}^d$  to a depth feature space.

**Definition 2.4.10.** The depth feature vector  $\mathbf{x}^D$  of an observation  $\mathbf{x}$ , with respect to the population distributions  $F_i$  and a choice of depth function  $D$ , is defined as

$$\mathbf{x}^D = (D(\mathbf{x}, F_1), \dots, D(\mathbf{x}, F_k)). \quad (2.4.13)$$

*Remark.* The graph

$$\text{DD}(F_1, \dots, F_k) = \{\mathbf{x}^D : \mathbf{x} \in \mathbb{R}^d\} \quad (2.4.14)$$

is the analogue of the **DD plot**, with  $k$  distributions.

Assuming that the depth function  $D$  only takes values in  $[0, 1]$ , the map  $\mathbf{x} \mapsto \mathbf{x}^D$  takes values in  $[0, 1]^k$ , regardless of the dimensionality of the original vector  $\mathbf{x}$ . With this, the maximum depth classification rule can be expressed as

$$\hat{\iota}_D(\mathbf{x}) = i \iff \mathbf{x}^D \in R_i^D = \{\mathbf{y} \in [0, 1]^k : y_i = \max_j y_j\}. \quad (2.4.15)$$

Indeed, any partition of the unit cube  $[0, 1]^k$  into  $k$  decision regions  $R_i^D$  gives rise to a depth based classifier. The DD classifier achieves this by using an increasing separating function  $r$  to partition  $[0, 1]^2$ . Furthermore,  $r \in \Gamma$  is chosen so as to best separate the training data  $\mathcal{D}$  transformed into the depth feature space. However, we can in principle use the transformed training data

$$\mathcal{D}^D = \{(\mathbf{x}_{ij}^D, i)\} \subset [0, 1]^k \times \{1, \dots, k\} \quad (2.4.16)$$

along with any multivariate classification algorithm (LDA, QDA,  $k$ NN, GLM, etc) to devise suitable decision regions. This is the basis of the  $DD^G$  classifier (Cuesta-Albertos et al., 2017).

## 2.5 Clustering

## 2.6 Outlier detection

## *Chapter 3*

# FUNCTIONAL DATA

### 3.1 Classification

### 3.2 Clustering

### 3.3 Outlier detection

### 3.4 Partially Observed Functional Data

## *Chapter 4*

# LOCAL DEPTH FUNCTIONS

## 4.1 Regression using Local Depth Regions

## *Chapter 5*

## CONCLUSION

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