# Summary Part 3: Applications of Concentration Inequalities

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## 1 Applications

#### 1.1 U-Statistics

• Example (*U-Statistics*) [Wainwright, 2019] Let  $g: \mathbb{R}^2 \to \mathbb{R}$  be a *symmetric function* of its arguments. Given an i.i.d. sequence  $\{X_k, k \geq 1\}$ , of random variables, the quantity

$$U := \frac{1}{\binom{n}{2}} \sum_{j < k} g(X_j, X_k) \tag{1}$$

is known as a **pairwise U-statistic**. For instance, if g(s,t) = |s-t|, then U is an unbiased estimator of the mean absolute pairwise deviation  $\mathbb{E}[|X_1 - X_2|]$ . Note that, while U is **not** a sum of independent random variables, the dependence is relatively weak, and this fact can be revealed by a martingale analysis.

If g is bounded (say  $||g||_{\infty} \leq b$ ), then the Bounded Difference Inequality can be used to establish the concentration of U around its mean. Viewing U as a function  $f(x) = f(x_1, \ldots, x_n)$ , for any given coordinate k, we have

$$\left| f(x) - f(x^{(-k)}) \right| \le \frac{1}{\binom{n}{2}} \sum_{j \neq k} \left| g(x_j, x_k) - g(x_j, x_k') \right| 
\le \frac{(n-1)2b}{\binom{n}{2}} = \frac{4b}{n},$$

so that the bounded differences property holds with parameter  $L_k = \frac{4b}{n}$  in each coordinate. Thus, we conclude that

$$\mathbb{P}\left\{|U - \mathbb{E}\left[U\right]| \ge t\right\} \le 2\exp\left(-\frac{n\,t^2}{8b^2}\right),\,$$

This tail inequality implies that U is a consistent estimate of  $\mathbb{E}[U]$ , and also yields finite sample bounds on its quality as an estimator. Similar techniques can be used to obtain tail bounds on U-statistics of higher order, involving sums over k-tuples of variables.

#### 1.2 Jackknife Estimation and Boostrapping

#### 1.3 Kernel-Density Estimation

• Example (Kernel Density Estimation)

Let  $Z_1, \ldots, Z_n$  be i.i.d. samples drawn according to some (unknown) density  $\phi$  on the real line. The density is estimated by the kernel estimate

$$\phi_n(z) = \frac{1}{n h_n} \sum_{i=1}^n K\left(\frac{z - Z_i}{h_n}\right),$$

where  $h_n > 0$  is a *smoothing parameter*, and K is a nonnegative function with  $\int K(z) = 1$ . The performance of the estimate is typically measured by **the**  $L_1$  **error**:

$$X(n) := f(Z_1, ..., Z_n) = \int |\phi(z) - \phi_n(z)| dz.$$

It is easy to see that

$$\left| f(z_1, \dots, z_n) - f_i(z_1, \dots, z_{i-1}, z_i', z_{i+1}, \dots, z_n) \right| \le \frac{1}{nh_n} \int \left| K\left(\frac{z - z_i}{h_n}\right) - K\left(\frac{z - z_i'}{h_n}\right) \right| dz$$

$$\le \frac{2}{n},$$

so without further work we obtain

$$\operatorname{Var}(X(n)) \le \frac{1}{n}$$

It is known that for every  $\phi$ ,  $\sqrt{n}\mathbb{E}[X(n)] \to \infty$ , which implies, by Chebyshev's inequality, that for every  $\epsilon > 0$ 

$$\mathbb{P}\left\{\left|\frac{X(n)}{\mathbb{E}\left[X(n)\right]} - 1\right| > \epsilon\right\} = \mathbb{P}\left\{\left|X(n) - \mathbb{E}\left[X(n)\right]\right| > \epsilon \mathbb{E}\left[X(n)\right]\right\} \le \frac{\operatorname{Var}(X(n))}{\epsilon^2(\mathbb{E}\left[X(n)\right])^2} \to 0$$

as  $n \to \infty$ . That is,  $\frac{X(n)}{\mathbb{E}[X(n)]} \to 1$  in probability, or in other words, X(n) is relatively stable. This means that the random  $L_1$ -error essentially behaves like its expected value.

By bounded difference inequality, we have

$$\mathbb{P}\left\{|X(n) - \mathbb{E}\left[X(n)\right]| \ge t\right\} \le 2\exp\left(-\frac{nt^2}{2}\right) \quad \blacksquare$$

#### 1.4 Random Graph

• Example (Clique Number in Erdös-Rényi Random Graphs) [Wainwright, 2019] Let  $\mathcal{G} = (\mathcal{V}, \mathcal{E})$  be an undirected graph, where  $\mathcal{V} = \{1, \ldots, d\}$  is the vertex set and  $\mathcal{E} = \{(i, j), i, j \in \mathcal{V}\}$  is the undirected edge set. A graph clique C is a subset of vertices such that  $(i, j) \in \mathcal{E}$  for all  $i, j \in C$ . The clique number  $C(\mathcal{G})$  of the graph is the cardinality of the largest clique. Note that  $C(\mathcal{G}) \in [1, d]$ . When the edges  $\mathcal{E}$  of the graph are drawn according to some random process, then the clique number  $C(\mathcal{G})$  is a random variable, and we can study its concentration around its mean  $\mathbb{E}[C(\mathcal{G})]$ .

<u>The Erdös-Rényi ensemble</u> of random graphs is one of the most well-studied models: it is defined by a parameter  $p \in (0,1)$  that specifies the probability with which each edge (i,j) is included in the graph, independently across all  $\binom{d}{2}$  edges. More formally, for each i < j, let us introduce a **Bernoulli edge-indicator** variable  $X_{i,j}$  with parameter p, where  $X_{i,j} = 1$  means that edge (i,j) is included in the graph, and  $X_{i,j} = 0$  means that it is not included.

Note that the  $\binom{d}{2}$ -dimensional random vector  $Z := \{X_{i,j}\}_{i < j}$  specifies the edge set; thus, we may view the clique number  $C(\mathcal{G})$  as a function  $Z \to f(Z)$ . Based on definition in Section ??, we see that f(Z) is a **configuration function** with property of "being in a clique".

Let Z' denote a vector in which a *single coordinate* of Z has been changed, and let  $\mathcal{G}'$  and  $\mathcal{G}$  be the associated graphs. It is easy to see that  $C(\mathcal{G}')$  can differ from  $C(\mathcal{G})$  by at most 1, so that

$$\left| f(Z) - f(Z') \right| \le 1,$$

Thus, the function  $C(\mathcal{G}) = f(Z)$  satisfies the bounded difference property in each coordinate with parameter L = 1, so that

$$\mathbb{P}\left\{\frac{1}{n}\left|C(\mathcal{G}) - \mathbb{E}\left[C(\mathcal{G})\right]\right| \ge \delta\right\} \le 2\exp\left(-2n\delta^2\right).$$

Consequently, we see that the clique number of an  $Erd\ddot{o}s$ - $R\acute{e}nyi\ random\ graph$  is  $very\ sharply\ concentrated\ around\ its\ expectation.$ 

#### 1.5 Minimum Weight Spanning Tree

#### 1.6 Rademacher Complexity

• Example (Rademacher Complexity) [Wainwright, 2019] Let  $\{\epsilon_k\}_{k=1}^n$  be an i.i.d. sequence of Rademacher variables (i.e., taking the values  $\{-1, +1\}$  equiprobably). Given a collection of vectors  $\mathcal{A} \subset \mathbb{R}^n$ , define the random variable

$$Z := \sup_{a \in \mathcal{A}} \sum_{k=1}^{n} \epsilon_k a_k = \sup_{a \in \mathcal{A}} \langle a, \epsilon \rangle.$$
 (2)

The random variable Z measures the size of A in a certain sense, and its expectation

$$\mathfrak{R}(\mathcal{A}) := \mathbb{E}\left[Z(\mathcal{A})\right] \tag{3}$$

is known as the Rademacher complexity of the set A.

Let us now show how the bounded difference inequality can be used to establish that Z(A) is **sub-Gaussian**. Viewing Z(A) as a function  $(\epsilon_1, \ldots, \epsilon_n) \to f(\epsilon_1, \ldots, \epsilon_n)$ , we need to bound the maximum change when coordinate k is changed. Given two Rademacher vectors  $\epsilon, \epsilon' \in \{-1, +1\}^n$ , recall our definition of the modified vector  $\epsilon^{(-k)}$ . Since

$$f(\epsilon^{(-k)}) \ge \langle a, \epsilon^{(-k)} \rangle$$
, for any  $a \in \mathcal{A}$ ,

we have

$$\langle a, \epsilon \rangle - f(\epsilon^{(-k)}) \le \langle a, \epsilon - \epsilon^{(-k)} \rangle = a_k(\epsilon_k - \epsilon'_k) \le 2 |a_k|.$$

Taking the supremum over A on both sides, we obtain the inequality

$$f(\epsilon) - f(\epsilon^{(-k)}) \le 2 \sup_{a \in \mathcal{A}} |a_k|.$$

Since the same argument applies with the roles of  $\epsilon$  and  $\epsilon^{(-k)}$  reversed, we conclude that f satisfies the bounded difference inequality in coordinate k with parameter  $L_k := 2 \sup_{a \in \mathcal{A}} |a_k|$ .

Consequently, the bounded difference inequality implies that the random variable  $Z(\mathcal{A})$  is sub-Gaussian with parameter at most  $2\sqrt{\sum_{k=1}^n\sup_{a\in\mathcal{A}}a_k^2}$ . This sub-Gaussian parameter can be reduced to the (potentially much) smaller quantity  $\sqrt{\sup_{a\in\mathcal{A}}\sum_{k=1}^na_k^2}$  using alternative techniques.

#### 1.7 Dimensionality Reduction

# 2 Self-Bounding Functions

#### 2.1 Definitions, Variance Bounds and Concentration

• Another simple property which is satisfied for many important examples is the so-called self-bounding property.

#### Definition (Self-Bounding Property)

A nonnegative function  $f: \mathcal{X}^n \to [0, \infty)$  has the <u>self-bounding property</u> if there exist functions  $f_i: \mathcal{X}^{n-1} \to \mathbb{R}$  such that for all  $z_1, \ldots, z_n \in \mathcal{X}$  and all  $i = 1, \ldots, n$ ,

$$0 \le f(z_1, \dots, z_n) - f_i(z_1, \dots, z_{i-1}, z_{i+1}, \dots, z_n) \le 1$$
(4)

and also

$$\sum_{i=1}^{n} \left( f(z_1, \dots, z_n) - f_i(z_1, \dots, z_{i-1}, z_{i+1}, \dots, z_n) \right) \le f(z_1, \dots, z_n).$$
 (5)

• Remark Clearly if f has the self-bounding property,

$$\sum_{i=1}^{n} \left( f(z_1, \dots, z_n) - f_i(z_1, \dots, z_{i-1}, z_{i+1}, \dots, z_n) \right)^2 \le f(z_1, \dots, z_n)$$
 (6)

Taking expectation on both sides, we have the following inequality

• Corollary 2.1 [Boucheron et al., 2013]
If f has the self-bounding property, then

$$Var(f(Z)) < \mathbb{E}[f(Z)].$$

• Remark (*Relative Stability*) [Boucheron et al., 2013] A sequence of nonnegative random variables  $(Z_n)_{n\in\mathbb{N}}$  is said to be *relatively stable* if

$$\frac{Z_n}{\mathbb{E}\left[Z_n\right]} \stackrel{\mathbb{P}}{\to} 1.$$

This property guarantees that the random fluctuations of  $Z_n$  around its expectation are of negligible size when compared to the expectation, and therefore most information about the size of  $Z_n$  is given by  $\mathbb{E}[Z_n]$ .

Bounding the variance of  $Z_n$  by its expected value implies, in many cases, the relative stability of  $(Z_n)_{n\in\mathbb{N}}$ . If  $Z_n$  has the self-bounding property, then, by Chebyshev's inequality, for all  $\epsilon > 0$ ,

$$\mathbb{P}\left\{ \left| \frac{Z_n}{\mathbb{E}\left[Z_n\right]} - 1 \right| > \epsilon \right\} \le \frac{\operatorname{Var}(Z_n)}{\epsilon^2 (\mathbb{E}\left[Z_n\right])^2} \le \frac{1}{\epsilon^2 \mathbb{E}\left[Z_n\right]}.$$

Thus, for relative stability, it suffices to have  $\mathbb{E}[Z_n] \to \infty$ .

#### 2.2 Configuration Function

• An important class of functions satisfying the self-bounding property consists of the so-called configuration functions.

#### **Definition** (Configuration Function)

Assume that we have a property  $\Pi$  defined over the union of finite products of a set  $\mathcal{X}$ , that is, a sequence of sets

$$\Pi_1 \subset \mathcal{X}, \ \Pi_2 \subset \mathcal{X} \times \mathcal{X}, \ \dots, \ \Pi_n \subset \mathcal{X}^n.$$

We say that  $(z_1, \ldots, z_m) \in \mathcal{X}^m$  satisfies the property  $\Pi$  if  $(z_1, \ldots, z_m) \in \Pi_m$ .

We assume that  $\Pi$  is  $\underline{hereditary}$  in the sense that if  $(z_1, \ldots, z_m)$  satisfies  $\Pi$  then so does any  $\underline{sub\text{-}sequence}$   $\{z_{i_1}, \ldots, z_{i_k}\}$  of  $(z_1, \ldots, z_m)$ .

The function f that maps any vector  $z = (z_1, \ldots, z_n)$  to **the size** of a **largest sub-sequence** satisfying  $\Pi$  is **the configuration function** associated with property  $\Pi$ .

• Corollary 2.2 [Boucheron et al., 2013] Let f be a configuration function, and let  $X = f(Z_1, ..., Z_n)$ , where  $Z_1, ..., Z_n$  are independent random variables. Then

$$Var(f(Z)) \leq \mathbb{E}[f(Z)].$$

#### 2.3 VC-Dimension and Growth Function

• Example (VC Dimension)

Let  $\mathcal{H}$  be an arbitrary collection of subsets of  $\mathcal{X}$ , and let  $x = (x_1, \ldots, x_n)$  be a vector of n points of  $\mathcal{X}$ . Define the **trace** of  $\mathcal{H}$  on x by

$$\operatorname{tr}(x) = \{A \cap \{x_1, \dots, x_n\} : A \in \mathcal{H}\}.$$

The shatter coefficient, (or Vapnik-Chervonenkis growth function) of  $\mathcal{H}$  in x is  $\tau_{\mathcal{H}}(x) = |\operatorname{tr}(x)|$ , the size of the trace.  $\tau_{\mathcal{H}}(x)$  is the number of different subsets of the n-point set  $\{x_1,\ldots,x_n\}$  generated by intersecting it with elements of  $\mathcal{H}$ . A subset  $\{x_{i_1},\ldots,x_{i_k}\}$  of  $\{x_1,\ldots,x_n\}$  is said to be **shattered** if  $2^k = T(x_{i_1},\ldots,x_{i_k})$ .

The VC dimension D(x) of  $\mathcal{H}$  (with respect to x) is the cardinality k of the largest shattered subset of x. From the definition it is obvious that f(x) = D(x) is a **configuration function** (associated with the property of "shatteredness") and therefore if  $X_1, \ldots, X_n$  are independent random variables, then

$$Var(D(X)) \le \mathbb{E}[D(X)].$$

- 2.4 Longest Increasing Subsequence
- 2.5 Weakly Self-Bounding Functions
- 3 Random Matrices
- 3.1 Definitions
- 3.2 Concentration Inequalities of Random Vectors
- 3.3 Concentration of Norm of Gaussian Vectors
- 3.4 Spectral Distribution of Hermitian Matrix: Semi-Circular Law
- 3.5 Largest Eigenvalue of Hermitian Random Matrix
- 4 Empirical Process
- 4.1 Definition
- 4.2 Tail Bounds for Empirical Processes
- 4.3 Uniform Law of Large Numbers
- 4.4 Suprema of Empirical Processes
- 4.5 Covering Number, Packing Number and Metric Entropy
- 4.6 Chaining
- 4.7 VC-Dimension
- 4.8 Variance Bounds

# References

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