The Foundation of Big Data: Experiments, Formulation, and Applications

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1 INTRODUCTION

The central theme of this talk \cite{1} is to promote the non-asymptotic statistical viewpoint in the context of massive datasets. The classical viewpoint breaks down when the data size becomes large.

2 Non-Asymptotic Viewpoint

In \cite{2}, the author’s group has shown that the classical GLRT is not optimal for large data lengths. This discovery has triggered a series of books \cite{3, 4, 5}.

Very recently\cite{2}, the work of Polyanskiy, Poor and Verdu (2010) \cite{6} has come to our attention. In finite-blocklength coding, Polyanskiy, Poor and Verdu proves the following result

\[
R(\varepsilon, n) = C - \sqrt{\frac{V}{n}} Q^{-1}(\varepsilon) + O\left(\frac{1}{\sqrt{n}}\right),
\]

where \(\varepsilon\) is the error probability, \(n\) is the coding length, \(V\) is the dispersion and \(Q(x)\) is the standard error function. This result says that Shannon’s result on channel capacity \(C\) is only the asymptotic term. His result is not optimal for finite blocklengths. As a matter of fact, the second order term \(\sqrt{\frac{V}{n}} Q^{-1}(\varepsilon)\) is explicitly given in \cite{1}.

The spirits of two above works are identical. The classical results are optimal only in the asymptotic limit \(n \to \infty\). In the age of Big Data, the size of \(n\) is large, but FINITE! The large but finite \(n\) asks for a new paradigm shift in analytical tools.

3 Work at TTU Since 2010

High scalability is the most important requirement for big data analysis. Size is the only thing that matters. It is natural to model massive datasets by using large random matrices. In the asymptotic regimes, the sizes of random matrices are assumed to approach the infinity. For example, for a random matrix \(X\) of size \(m \times n\), we assume the asymptotic regime:

\[m \to \infty, n \to \infty \text{ but } m/n \to c \in (0, \infty).\]

The work \cite{3} give a survey of the work in this direction and study the connection with cognitive radio network.

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4 Data Collection and Modeling

4.1 Data Collection

Figure 1 shows a 80-node network made of software-defined radios (USRP). The potential research topics include: time-varying topology, UAVs, random arrays, anti-jamming, network security and privacy of data.

The central message here is to show the big data problem arose from the natural engineering environment. Google downloads all webpages from the Internet, and saves them for data mining. Big Data at least traces back to its practice. In the same spirit, can we save and makes sense of "all the data" in wireless networks? The answer is a big YES. TTU has saved all the radio waveforms (modulated waveforms) and attempted to make sense of these massive datasets. Real-time, streaming data of massive size is the hallmark of this work.

Figure 2 shows the real deployment of TTU’s cognitive radio network. We have 80-node USRPs. Massive MIMO can be studied in this testbed. Figure 3 illustrates waveforms of 80 soft-defined radios. One asks a basic question: How do we represent these massive datasets? In the next subsection, we propose to use large random matrices for this purpose.

Fig. 1: A Large-Scale Advanced Radio Network Testbed.

The non-asymptotic regime is defined as: $m$ and $n$ are large, but finite. In [4], we use concentration inequalities as the mathematical foundation. The book goes to study compressed sensing, low rank matrix recovery, estimation, detection and database-friendly computing. Large random matrices can unify many big data problems and form the foundation for big data analysis. In [5], we build upon this key observation. There we go on to study non-Hermitian random matrices using free probability theory. The three books promote the framework of using large random matrices (Hermitian or non-Hermitian) to represent massive datasets.
Fig. 2: The photo of the 80-node testbed at TTU.

Fig. 3: Waveforms of soft-defined radios. The above represents 40 time slots of data recorded at one USPR receiver; the bottom is zoomed in for each time slot. The sample rate is 20 Msp.
4 Data Collection and Modeling

4.2 Large Random Matrices

In [7,8], we model the massive datasets using large random matrices.

4.3 Empirical Spectrum Density

Figure 4 shows empirical spectral density: noise only. Any deviation from "noise only" case will indicate the presence of "some signal". We observe a bulk cluster that is detached from the noise cluster. Note the noise cluster follows the Marchenko-Pasture law, as shown in Figure 4.

4.4 Product of Non-Hermitian Random Matrices

Figure 6 shows the product of non-Hermitian random matrices: noise only. The product of non-Hermitian random matrices can be studied by the new free probability theory. The ring law is so universal for a basic building block, a rectangular random matrix.

Figure 7 shows the product of non-Hermitian random matrices: signal plus noise. Compared with the previous case of "Noise only", correlation of the signal reduces the inner radius of the ring law. We also observed this using data in other fields.

Figure 8 the product of large random Ginibre matrices. The product of $k$ large random Ginibre matrices has closed form. Gaussian is a special case of Ginibre matrix.
4.5 5G Massive MIMO: Euclidean Random Matrices

In the next generation (5G) wireless system, massive MIMO uses hundreds (even thousands of) antennas that co-located or distributed in two dimensions (2D) or three dimensions (3D).

Here we are interested in the massive datasets (random fields distributed in 3D). Modeling the random fields (or the channel) is a challenge. The idea here is to study the free-space propagation and thus the closed-form solution can be obtained [5]. Figure 10 shows massive MIMO modeled using Euclidean random matrices.
4 Data Collection and Modeling

The unit circles assume the noise only. We need a theory for predicting the signal.

Experiments agree with (closed-form) theoretical predictions.

\[
\rho^{(2)}(x) = \frac{2^{1/3}}{12\pi} \sqrt{3 \left( 27 + 3\sqrt{81 - 12x} \right)} \left( 27 + 3\sqrt{81 - 12x} \right)^{2/3} - 6x^{2/3},
\]

\[x \in [0, \frac{27}{4}].\]

Fig. 7: Product of Non-Hermitian Random Matrices: Signal plus Noise

Fig. 8: Product of Large Random Ginibre Matrices
The Green’s function between an arbitrary transmitter located at $r_i$ and an arbitrary receiver located at $r_j$ is expressed as

$$\left(\nabla^2 + k_0^2 + i\varepsilon_0\right) G(r_i, r_j) = -\frac{4\pi}{k_0} \delta(r_i, r_j).$$

where $k_0 = 2\pi/\lambda_0$ with $\lambda_0$ being the wavelength in free space, and $\varepsilon_0$ is the electric constant. In free space, we obtain the closed-form expression

$$G(r_i, r_j) = \exp \left(i k_0 |r_i - r_j|\right) \frac{|r_i - r_j|}{|r_i - r_j|}. \tag{2}$$

The channel propagation between an arbitrary transmitter located at $r_i$ and an arbitrary receiver located at $r_j$ is modeled by matrix entries

$$A_{ij} = f(r_i, r_j), i, j = 1, \ldots, N. \tag{3}$$

Here $N = 10,000$ is considered. It turns out that (3) can be represented in terms of the product of large random matrices

$$A = HTH^H, \tag{4}$$

leads to the use of free probability theory. In free space, we have

$$A_{ij} = (1 - \delta_{ij}) \exp \left(i k_0 |r_i - r_j|\right) \frac{|r_i - r_j|}{|r_i - r_j|}. \tag{5}$$

The fundamental property of the large random matrix $A$ is that $A$ is non-Hermitian and thus its eigenvalues are complex in general, as shown in Figure [10]. Two free parameters determine the distribution of eigenvalues in the complex plane. One is the signal-to-noise ratio $\gamma$, and the other is $\rho \lambda_0^3$, and $\rho$ is the density of users (number of users per meter).
A Non-Asymptotic Theory of Detection

In [2], we discover that the larger the length of the data, the bigger the gain (compared with the classical estimator-correlator test). It is commonly believed that the classical estimator-correlator test is optimal. To understand this effect, we are motivated for these three works [3–5].

Let us give a tutorial introduction of this problem. We follow the classic Van Tree (1968) [9] for the classical treatment. For (column) random vectors

\[ x = [X_1, ..., X_n]^T, \quad y = [Y_1, ..., Y_n]^T, \quad s = [S_1, ..., S_n]^T, \]

we define their covariance matrices as

\[ R_x = \mathbb{E} [xx^T]; \quad R_y = \mathbb{E} [yy^T]; \quad R_s = \mathbb{E} [ss^T], \]

where \( \mathbb{E} \) is the mathematical expectation. In practice, we only access \( N \) independent copies of random vectors so form the random matrices

\[ X = [x_1, ..., x_N]; \quad Y = [y_1, ..., y_N]; \quad S = [s_1, ..., s_N]. \]

We are, in particular, interested in the following regime: \( N \) and \( n \) are large but finite! Now we are ready to compare three formulations

\[
\begin{align*}
\mathcal{H}_0: y &= x, \\
\mathcal{H}_1: y &= s + x, \\
\mathcal{H}_0: R_y &= R_x \\
\mathcal{H}_1: R_y &= R_s + R_x \\
\mathcal{H}_0: Y &= X \\
\mathcal{H}_1: Y &= S + X,
\end{align*}
\]

where the first formulation follows from Van Tree (1968) [9], and the second one is for quantum detection [3], and the third one is used in [2]. Formulation II is the limit form for

\[ n \text{ is fixed, but } N \rightarrow \infty, \]

Massive MIMO: Euclidean Random Matrices

\[ A_{ij} = f(r_i, r_j), \quad i, j = 1, ..., N. \]

\[ A = HTH^H \]

\[ (\nabla^2 + k_0^2 + i\varepsilon)G(r_i, r_j) = -\frac{4\pi}{k_0}\delta(r_i - r_j), \]

\[ G(r_i, r_j) = \frac{\exp(\frac{i k_0 |r_i - r_j|}{k_0})}{k_0 |r_i - r_j|}. \]

\[ A_j = (1 - \delta_j)\frac{\exp(\frac{i k_0 |r_i - r_j|}{k_0})}{k_0 |r_i - r_j|}. \]

Fig. 10: Massive MIMO: Euclidean Random Matrices
which, however, is not true in general. We are often interested in the non-asymptotic regime:

\[ N \to \infty, n \to \infty \text{ but } N/n \to c \in (0, \infty). \quad (7) \]

Consider Formulation I first, for simplicity we assume that \( x \) a zero-mean, Gaussian random vector. A vector \( z \) is a Gaussian random vector when its components \( z_1, z_2, \ldots, z_n \) are jointly Gaussian random variables. We further assume that

\[ \mathbb{E}[y] = \mathbb{E}[s] = m. \]

Of course \( m \) is a deterministic vector. In practice, we need to estimate \( m \). Thus we have

\[ R_y = \mathbb{E}[(y - m)(y - m)^T]. \]

The probability density of \( y \) is

\[ p_y = \frac{1}{(2\pi)^{n/2}(\det R_y)^{1/2}} \exp\left[-\frac{1}{2} (y - m)^T R_y^{-1} (y - m)\right]. \]

The likelihood ratio test follows easily:

\[ \Lambda = \frac{1}{(2\pi)^{n/2}(\det R_y)^{1/2}} \exp\left[-\frac{1}{2} y^T R_y^{-1} y\right] < \eta, \quad \mathcal{H}_1 \]

Taking logarithms, we obtain

\[ \frac{1}{2} (y - m)^T R_y^{-1} (y - m) - \frac{1}{2} y^T R_y^{-1} y > \ln \Lambda + \frac{1}{2} \ln \det R_y - \frac{1}{2} \ln \det R_x \triangleq \gamma^*. \quad (8) \]

Consider the special case:

\[ R_y = R_x = R. \quad (9) \]
Equation (8) is equivalent to
\[
   l(y) = m^T R y > \gamma^*.
\]
The quantity on the left is a scalar Gaussian random variable, for it was obtained by a linear transformation of jointly Gaussian random variables. The distance between the means on the two hypotheses is defined as
\[
   d^2 \triangleq \frac{[E(l|H_1) - E(l|H_0)]^2}{\text{Var}(l|H_1)}.
\]
After some manipulation, we obtain
\[
   d^2 = m^T R m.
\]
When for white Gaussian noise
\[
   R = \sigma^2 I_n,
\]
we have
\[
   d^2 = \frac{m^T m}{\sigma^2} = \|m\|^2/\sigma^2
\]
where \(\|\cdot\|\) is the Euclidean norm of a vector. \(\|m\|^2\) can be viewed as the signal power.

Now let us turn our attention Formulation III in (5). We can always write
\[
   \mathcal{H}_0 : \frac{1}{N} Y^H Y = \frac{1}{N} X^H X = \frac{1}{N} \sum_{i=1}^N x_i^H x_i,
\]
\[
   \mathcal{H}_1 : \frac{1}{N} Y^H Y = \frac{1}{N} S^H S + \frac{1}{N} X^H X = \frac{1}{N} \sum_{i=1}^N s_i^H s_i + \frac{1}{N} \sum_{i=1}^N x_i^H x_i,
\]
where \(x_1, \ldots, x_N\) are \(N\) independent copies of the random vector \(x\) and \(s_1, \ldots, s_N\) are \(N\) independent copies of the random vector \(s\).

In this modern non-asymptotic view of (7), the derivation of Formulation I is not justified. The central argument is the expectation operation \(E[\cdot]\) in analysis, which requires (6), indirectly through the central limit theorem [10].

**Theorem 5.1 (Central Limit Theorem):** Let \((x_i)_{i \geq 1}\) be i.i.d. \(\mathbb{R}^n\)–valued random variables. Let the (vector) \(\mu = E[x_i]\), and let \(\Sigma\) denote the covariance matrix: \(\Sigma = (\Sigma_{k,l})_{1 \leq k,l \leq n}\), where \(\Sigma_{k,l} = \text{Cov}(X_i^k, X_i^l)\), where \(X_i^k\) is the \(k\)–th component of the \(\mathbb{R}^n\)–valued random variable \(x_i\). Then
\[
   \lim_{N \to 0} \frac{s_N - N\mu}{\sqrt{N}} = z,
\]
where \(s_N = \sum_{i=1}^N x_i\) and the distribution of \(z\) is zero-mean Gaussian \(\mathcal{N}(0, \Sigma)\) and where the convergence is in distribution.

The central limit theorem is not valid for non-asymptotic regime (7).
Consider an example for white Gaussian noise, \((12)\) cannot be justified. We consider the hollow Wishart matrix

\[
W_n = \frac{1}{N} \sum_{i=1}^{N} x_i^H x_i - \sigma^2 I_n. \tag{15}
\]

For massive data, the noise is aggregated. The effect of noise aggregation cannot be ignored, as shown in Figure \(12\). Understanding noise aggregation is central to big data problems. Without loss of generality, Formulation III \((14)\) can be rewritten as

\[
\mathcal{H}_0 : W_n, \quad \mathcal{H}_1 : \frac{1}{N} \sum_{i=1}^{N} s_i^H s_i + W_n \tag{16}
\]

where \(W_n\) is the \(n \times n\) random matrix. Our intuition is that the trace (sometimes called average) of \(W_n\) tends to vanish when \(n\) is large enough. The central task is to find the non-asymptotic relation as a function of \(n\).

Using concentration inequalities, it is shown in \([4]\) that the scalar random variable

\[
Z = \text{Tr} \left( W_n \right) \tag{17}
\]

behaves very nicely. We obtain \([4]\) p.221

\[
\mathbb{E} [Z] = 0, \quad \text{Var} [Z] \leq \frac{c}{n^2}, \tag{18}
\]

where \(c\) is a constant. We guess that the convergence rate \(1/n^2\) may be optimal. Our algorithm claims \(\mathcal{H}_1\) when \(Z\) is above zero.
6 The Mathematical Foundation of Big Data

In the next subsection, we take material from the NSF workshop [11].

6.1 Overview of Big Data

- Big engineering data has unique characteristics: more disciplined and regulated.
- Emerging engineering systems with big-data opportunities: smart grids, sensor nets, transportation, telemedicine, aerospace, testing, safety, nuclear, design blueprints and more.
- Rethink of data collection and storage to facilitate big data processing and inference tasks.
- How do we trade-off complexity for accuracy in massive decentralized signal and data analysis tasks?
- How can efficient signal and data analysis algorithms be developed for big, unstructured or loosely structured data?
- What are the basic principles and useful methodologies to scale inference and learning algorithms and trade off the computational resources (e.g., time, space and energy) according to the needs of engineering practice (e.g. robustness vs. efficiency, real-time)?

Big Data processing and analysis is summarized below

- Integration of very heterogeneous data: correlation mining in massive database;
- Data at vastly different scales and noise levels; Mixture of continuous and categorical variables.
- Reliable and robust quantitative models: Uncertainty quantitation; Adaptive to drift over time.
- High throughput real-time processing: Smart adaptive sampling and compression; Distributed or parallel processing architectures.
- Interactive user interfaces: Human-in-the-loop processing;
- Visualization and dimensionality reduction.

Some signal processing challenges include

- Heterogeneous data integration: Ranking signals for human-aided selection of relevant variables; Fusing graphs, tensors, and sequence data; Active visualization: dimensionality reduction.
- Flexible low complexity modeling and computation: Scalable signal processing: distributed algorithms and implementation; Smart sampling: feedback controlled signal search and acquisition.
- Reliable robust models for anomaly detection and classification: Parsimonious signal processing: Sparse correlation graphical models; Decomposable signal processing: factored models and algorithms
Large random matrices for Big Data. Random matrices play a central role in statistics in the context of multivariate
data. The continued growth of big data has given rise to high dimensional statistics analysis. Convex analysis,
Riemannian geometry and combinatorics are relevant. Random matrix theory (RMT) has emerged as a particularly
useful framework for many theoretical questions associated with the analysis of high-dimensional multivariate data.
RMT affects the modern statistical thinking in two ways. Most of the mathematical treatments of RMT have focused
on matrices with high degree of independence in the entries, which one may refer to as “unstructured” random
matrices. Recall that about 75% of big data are unstructured. In high-dimensional statistics, we are primarily
interested in problems where there are lower dimensional structures buried under random noise.

Traditional statistical theory, particularly in multivariate analysis, did not contemplate the demands of high dimen-
sionality in data analysis. Classical multivariate analysis textbooks consider: (1) Only asymptotically optimum, not
optimum for any size data; (2) dimension of the dataset, \( p \) is fixed small constant or at least negligible compared
with the sample size \( n \). Modern datasets have: (1) Their dimensions \( p \) can be proportionally large compared with the
sample size \( n \). (2) Examples: sensor network data, wireless network data, smart grid data, financial data, consumer
data, manufacturing data, multimedia data, etc. (3) The likelihood ratio test (LRT), the classical statistical inference,
must be revisited.

In this next subsection, we draw material heavily from [12].

6.2 Probability in High Dimension

“High dimension” refers to the presence of many distinct but interacting random variables, including

- Large random structure: random matrices, random graphs,
- Statistics and machine learning: estimation, prediction and model selection for high-dimensional data.
- Randomized algorithms in computer science.
- Random codes in information theory.
- Statistical physics: Gibbs measure, percolation, spin glasses,
- Random combinatorial structures: longest increasing subsequence, spanning trees, travelling salesman prob-
  lem, . . .
- Probability in Banach spaces: probabilistic limit theorems for Banach valued random variables, empirical
  processes, local theory of Banach spaces, geometric functional analysis, convex geometry.
- Mixing times and other phenomena in high-dimensional Markov chains.

Non-asymptotic probabilistic inequalities are especially relevant when the number of observables is large. They are
suitable for large complex sets of data with possibly huge collections of models at different scales. This framework
allows the collection of models together with their dimensions \( n \) to vary freely, letting the dimensions \( n \) be possibly
of the same order of magnitude as the number of observations \( p \). We are not concerned with limit theorems (as in
many classical probabilistic results), but rather with explicit estimates that are either dimension-free, or that capture
precisely the dependence of the problem on the relevant dimensional parameters. In asymptotic results, one must
take all these parameters to the limit in a fixed relation to one another, while the *non-asymptotic* viewpoint allows to express the interrelation between the different parameters in a much more precise way.

High-dimensional problems typically involve interactions between a large number of degrees of freedom whose aggregate contributions to the phenomenon of interest must be accounted for in the mathematical analysis; the explicit nature of non-asymptotic estimates makes them particularly well suited to be used as basic ingredients of the analysis, even if the ultimate result of interest is asymptotic in nature.

In Section 5 we give an introduction to the non-asymptotic theory of random matrix-valued hypothesis detection. In 2010, we were lucky to discover that concentration inequalities were indeed the probabilistic tools.

Three basic principles formulated informally as “theorems” are pointed out below. Of course, we need to make precise some terms. But rigor is not concern here.

**Theorem 6.1 (Concentration):** If \( X_1, \ldots, X_n \) are independent (or weakly dependent) random variables, then the random variable \( f(X_1, \ldots, X_n) \) is “close” to its mean \( \mathbb{E}[f(X_1, \ldots, X_n)] \), provided that the function \( f(x_1, \ldots, x_n) \) is not too “sensitive” to any of the coordinates \( x_i, i = 1, \ldots, n \).

This theorem is valid for very complicate functions such as eigenvalues of large random matrices. Sometimes the expectation of the function \( \mathbb{E}[f(X_1, \ldots, X_n)] \) may be unknown. Even in this scenario, concentration inequalities can still be used to find the fluctuations of this function. When the Central Limit Theorem is invalid, the next best thing you can do is to resort to concentration inequalities.

**Theorem 6.2 (Supremum):** If the random process \( \{X_t\}_{t \in T} \) is “sufficiently continuous”, then the magnitude of the supremum \( \sup_{t \in T} X_t \) is controlled by the “complexity” of the index set \( T \).

A family of random variables indexed by a set \( T \) that is frequently high- or infinite-dimensional. We often involve a large number of interdependent degrees of freedom; there is a need to obtain *simultaneous control* over many random variables. The investigation of suprema is in fact surprisingly general. Four functions are defined below in this framework. The first one is the operator norm of a matrix \( A \),

\[
\|A\| = \sup_{u,v \in B_2} \langle u, Av \rangle,
\]

where \( \langle \cdot \rangle \) represents the inner product of two vectors, \( B_2 \) is the Euclidean unit ball in \( \mathbb{R}^n \).

The second function is related to the norms of random vectors. Let \( X \) be a random vector in \( \mathbb{R}^n \), and let \( \|\cdot\|_B \) be any norm on \( \mathbb{R}^n \), where \( B \) denotes the unit ball of \( \|\cdot\|_B \). The duality theory of Banach spaces implies that we can write

\[
\|X\|_B = \sup_{t \in B^o} \langle t, X \rangle,
\]

where \( B^o \) denotes the dual ball. The supremum of the random processes

\[
\{X_t = \langle t, X \rangle\}_{t \in B^o}
\]

arises naturally in probability in Banach spaces.
The third function is related to empirical risk minimization. The problem is to estimate how close the empirical risk minimizer is to the optimal parameter as a function of the number of samples \( n \), the dimension of the parameter space \( \Theta \), the dimension of the state space \( X \), etc. Many problems in statistics and machine learning may be formulated as the problem of computing

\[
\arg \min_{\theta \in \Theta} \mathbb{E} [l(\theta, X)]
\]
given only observed “data” consisting of i.i.d. samples \( X_1, \ldots, X_n \sim X \). That is to, without knowledge of the law of \( X \). Here \( lF \) is a given loss function, and \( \Theta \) is a given parameter space, which depends on the problem at hand. Formally we have

\[
\sup_{\theta \in \Theta} \frac{1}{n} \sum_{i=1}^{n} \{l(\theta, X_i) - \mathbb{E}[l(\theta, X_i)]\},
\]

which is the supremum of a random process.

The fourth function is related a convex function \( f: \mathbb{R}^n \to \mathbb{R} \)

\[
f(x) = \sup_{y \in \mathbb{R}^n} \{\langle y, x \rangle - f^*(y)\},
\]

where \( f^* \) denotes the convex conjugate of \( f \).

Now we are ready to talk about the third principle: universality. The high-dimensional phenomenon is, in a sense, robust to the precise details of the model ingredients.

**Theorem 6.3 (Universality):** If \( X_1, \ldots, X_n \) are independent (or weakly dependent) random variables, then the expectation \( \mathbb{E} [f(X_1, \ldots, X_n)] \) is “insensitive” to the distribution of when the function \( f \) is “sufficiently smooth”.

Let \( X_1, X_2, \ldots \) be i.i.d. random variables with finite variance. The law of large numbers says

\[
\frac{1}{\sqrt{n}} \sum_{i=1}^{n} \{X_i - \mathbb{E}X_i\} \to 0 \text{ as } n \to \infty.
\]

We do not only know that the fluctuations are of order \( n^{-1/2} \) (as is captured by the concentration phenomenon), but we have much more precise information as well: by the central limit theorem, we have a precise description of the distribution of the fluctuations, as

\[
\frac{1}{\sqrt{n}} \sum_{i=1}^{n} \{X_i - \mathbb{E}X_i\} \approx \text{Gaussian}
\]

when \( n \) is large. A different way of saying this property is that

\[
\frac{1}{\sqrt{n}} \sum_{i=1}^{n} \{X_i - \mathbb{E}X_i\} \approx \frac{1}{\sqrt{n}} \sum_{i=1}^{n} \{G_i - \mathbb{E}G_i\},
\]

where \( G_k \) are independent Gaussian random variables with the same mean and variance of \( X_k \) (here \( \approx \) denotes closeness of the distributions).

As in the case of concentration, it turns out that this phenomenon is not restricted to linear functions of independent random variables, but is in fact a manifestation of a more general principle.
6.3 Talagrand’s Concentration of Measure

6.4 Free Entropy

7 Conclusion

Big data problems are ubiquitous. The basic questions: How do we represent the massive datasets? How do we make sense of those data? We promote the non-asymptotic statistical viewpoint, that is built upon concentration inequalities. Data-driven system design is enabled. Large Random matrices serve as basic building blocks. We only know little about the Big Data paradigm. One wonder if new Data Science is emerging for today’s and future’s challenges.

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