Robust estimations from distribution structures: V. Non-asymptotic

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Due to the complexity of order statistics, the finite sample bias of robust statistics is generally not analytically solvable. While the Monte Carlo method can provide approximate solutions, its convergence rate is typically very slow, making the computational cost to achieve the desired accuracy unaffordable for ordinary users. In this paper, we propose an approach analogous to the Fourier transformation to decompose the finite sample structure of the uniform distribution. By obtaining a set of sequences that are simultaneously consistent with a parametric distribution for the first four sample moments, we can approximate the finite sample behavior of robust estimators with significantly reduced computational costs. This article reveals the underlying structure of randomness and presents a novel approach to integrate two or more assumptions. 1 2 3 4 5 6 7 8 9 10 11 12 13

finite sample bias | order statistics | variance reduction | Monte Carlo study | uniform distribution

1 **I**₂ **I**_a n the early nineteenth century, Bessel deduced the unbiased sample variance and found it has a correction term of $\frac{n}{n-1}$. 1 n the early nineteenth century, Bessel deduced the unbiased ³ Later, Cramér [\(1\)](#page-3-0) in his classic textbook *Mathematical Meth-*⁴ *ods of Statistics* deduced unbiased sample central moments ⁵ with a linear time complexity. However, apart from the mean and central moments, the finite sample behavior of nearly all ⁷ other estimators depends on the underlying distribution and lacks a simple non-parametric correction term. For example, ⁹ the simplest robust estimator, the median, exhibits a highly 10 complex finite sample behavior. If *n* is odd, $E[median_n] =$ $\int_{-\infty}^{\infty} \left(\frac{n+1}{2}\right) \left(\frac{n}{2} - \frac{1}{2}\right) F(x)^{\frac{n}{2} - \frac{1}{2}} \left[1 - F(x)\right]^{\frac{n}{2} - \frac{1}{2}} f(x) x dx$ [\(2\)](#page-3-1), 12 where $F(x)$ and $f(x)$ represent the cumulative distribution ¹³ function (cdf) and probability density function (pdf) of ¹⁴ the assumed distribution, respectively. For the exponential ¹⁵ distribution, the above equation is analytically solvable, yield- $2^{-n-1}(n+1)\left(\frac{n}{2}\right)\left(H_n-H_{\frac{n-1}{2}}\right)$ $\Gamma(\frac{n+1}{2})\sqrt{\pi}$

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E[median_n] = \frac{2}{\sqrt{(\frac{n}{2}+1)}}
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,
\nwhere H_n denotes the *n*th Harmonic number, Γ represents
\nthe gamma function, and λ stands for the scale parameter of
\nthe exponential distribution. However, for distributions with
\nmore complex pdfs, such equations are generally unsolvable.
\n24 Another widely used exact finite sample bias correction is
\n25 the factor for unbiased standard deviation in the Gaussian
\n26 distinct distribution, which can be deduced using Cochran's theorem
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\n28 (3). For more complex estimators, writing their exact
\n29 finite-sample distribution formulas becomes challenging. In
\n2013, Nagatsuka, Kawakami, Kamakura, and Yamamoto
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\n2913, Nagatsuka, Kawakami, Kamakura, and Yamamoto
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simulation is often too high to be processed on a typical 33 PC. For example, for median absolute deviation, Croux and 34 Rousseeuw (1992) provided correction factors with a precision ³⁵ of three decimal places for $n \leq 9$ using 200,000 pseudorandom ³⁶ Gaussian sample (5) . Hayes (2014) reported correction factors 37 for $n \leq 100$ using 1 million pseudorandom samples for each ∞ value of *n* to ensure the accuracy to four decimal places (6) . 39 Recently, Akinshin (2022) [\(7\)](#page-3-6) presented correction factors ⁴⁰ for $n \leq 3000$ using 0.2-1 billion pseudorandom Gaussian 41 samples. His result suggest that, for the median absolute 42 deviation, finite sample bias correction is required to ensure 43 a precision of three decimal places when the sample size is ⁴⁴ smaller than 2000. This highlights the importance of finite $\frac{45}{45}$ sample bias correction. However, since different correction 46 factors are required for different parametric assumptions, the ⁴⁷ computational cost of addressing all possible cases in the real ⁴⁸ world becomes significant, especially for complex models. 49

In addition to computational challenges, there exists an 50 inherent difficulty in dealing with randomness. The theory 51 of probability provides a framework for modeling and under- ⁵² standing random phenomena. However, the practical imple- ⁵³ mentation of these models can be challenging, as discussed, $\frac{54}{4}$ and their complexity greatly hinders our comprehension. The 55 quality of randomness can significantly impact the validity of $\frac{56}{60}$ simulation results, and a deeper understanding of randomness 57 may offer a more effective and cost-efficient solution. The 58 purpose of this brief report is to demonstrate that the finite 59 sample structure of uniform random variables can be decom- 60 posed using a few well-designed sequences with high accuracy. ⁶¹ Furthermore, we show that the computational cost of estimating finite sample bias from a Monte Carlo study can be 63

Significance Statement

Most contemporary statistics theories focus on asymptotic analysis due to its tractability and simplicity. Non-asymptotic statistics are crucial when dealing with small or moderate sample sizes, which is often the case in practice. In situations where analytical results are difficult or impossible to obtain, Monte Carlo studies serve as a powerful tool for addressing non-asymptotic behavior. However, these studies can be computationally expensive, particularly when high precision is required or when the statistical model demands significant computational time. Here, we propose calibrated Monte Carlo study that aims to approximate the randomness structures using a small set of sequences. This approach sheds light on understanding the general structure of randomness.

Fig. 1. The frequency histograms of pseudo-random sequences on the interval [0,1] with size 80.

⁶⁴ dramatically improved by obtaining a set of sequences that ⁶⁵ are simultaneously consistent with a parametric distribution

⁶⁷ **Decomposing the finite sample structure of uniform** ⁶⁸ **distribution**

⁶⁹ Any continuous distribution can be linked to the uniform dis- π tribution on the interval [0, 1] through its quantile function. ⁷¹ This fundamental concept in Monte Carlo study implies that ⁷² understanding the finite sample structure of uniform random ⁷³ variables can be leveraged to understand the finite sample ⁷⁴ structure of any other continuous random variable through the τ ⁵ quantile transform. The Glivenko–Cantelli theorem $(8, 9)$ $(8, 9)$ $(8, 9)$ en-⁷⁶ sures the almost-sure convergence of the empirical distribution ⁷⁷ function to the true distribution function. However, the indi-⁷⁸ vidual empirical distribution often deviates significantly from ⁷⁹ the asymptotic distribution even when the sample size is not ⁸⁰ small (Figure [1,](#page-1-0) sample size is 80), which cause finite sample 81 biases of common estimators. Let μ , μ_2 , \ldots , μ_k denote the first ⁸² **k** central moments of a probability distribution. According to $\frac{1}{83}$ the unbiased sample central moment (1) , the expected value of the sample central moment, $m_{\mathbf{k}} = \frac{1}{n} \sum_{\mathbf{k}=1}^{n} (x_{\mathbf{k}} - \bar{x})^{\mathbf{k}}$, can 85 be deduced, denoted as $E[m_k]$. Let $S = \{sequence[i]|i \in \mathbb{N}\}\$ ⁸⁶ be a set of number sequences ranging from 0 to 1, where ⁸⁷ *sequence*[*i*] represents the *i*th sequence in the set, and N is 88 the set of natural numbers, with $i \leq N$. Transform every ⁸⁹ number in *S* using the quantile function of a parametric dis-⁹⁰ tribution, *PD*. The transformed sequences can be denoted as ⁹¹ *SPD*. Denote the set of the **k**th sample central moments 92 for these transformed sequences as $M_{\mathbf{k}} = \{m_{\mathbf{k},i} | i \in \mathbb{N}\}.$ 93 *S* is consistent with *PD* for all m_k when $k \leq k$, if and ⁹⁴ only if the following system of linear equations is consistent, $\int m_{1,1}w_1 + \ldots + m_{1,i}w_i + \ldots + m_{1,N}w_N = E[m_1]$ \int $\overline{\mathcal{L}}$ *. . .* $m_{\mathbf{k},1}w_1 + \ldots + m_{\mathbf{k},i}w_i + \ldots + m_{\mathbf{k},N}w_N = E[m_{\mathbf{k}}]$, where *. . .* $m_{k,1}w_1 + \ldots + m_{k,i}w_i + \ldots + m_{k,N}w_N = E[m_k]$ $w_1 + \ldots + w_i + \ldots + w_N = 1$

 $w_1, \ldots, w_i, \ldots, w_N$ are the unknowns of the system, with

 y^2 *N* $\geq k+1$. $w_1, \ldots, w_i, \ldots, w_N$ can be determined using a typi-

⁹⁸ cal constraint optimization algorithm. The Monte Carlo study

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can be seen as a special case when $w_1 = \ldots = w_i = \ldots = w_N$, so and the sequences in *S* are all random number sequences. The 100 strong law of large numbers (proven by Kolmogorov in 1933) 101 (10) ensures that in this case, when the number of sequences 102 $N \to \infty$ or when the sample size $n \to \infty$, the above system of 103 linear equations is always consistent.

Low-discrepancy sequences are commonly used as a re- ¹⁰⁵ placement of uniformly distributed random numbers to re- ¹⁰⁶ duce computational cost. When considering a sequence to 107 approximate the structure of uniform random variables, the ¹⁰⁸ most natural choice is the arithmetic sequence, denoted as 109 ${x_i}_{i=1}^n = {\frac{i}{n+1}}_{i=1}^n$. However, the arithmetic central moments estimated from the arithmetic sequence for the Gaus- 111 sian distribution differ significantly from their expected values 112 (Figure $2A$). The arithmetic sequence lacks the variability 113 of true random samples which produce additional biases for ¹¹⁴ even order moments. The beta distribution is defined on the ¹¹⁵ interval $(0, 1)$ in terms of two shape parameters, denoted by $_{116}$ *α* and *β*. When $α = β$, the beta distribution is symmetric. 117 To better replicate the features of uniform random variables, 118 we introduced beta distributions with a variety of parameters. 119 The arithmetic sequences were transformed by the quantile 120 functions of these beta distributions to form beta-sequences, 121 resulting in sequences that are U-shape ($\alpha = \beta = 0.547$), leftskewed ($\alpha = 46.761$, $\beta = 20.108$), right-skewed ($\alpha = 20.108$, 123 *β* = 46*.761*), monotonic decreasing (α = 0*.478*, *β* = 38*.53*), 124 monotonic increasing ($\alpha = 38.53$, $\beta = 0.478$), their left-skewed 125 self-mixtures ($\alpha = \beta = 0.369$, $\alpha = \beta = 18.933$), their rightskewed self-mixtures ($\alpha = \beta = 0.369$, $\alpha = \beta = 18.933$), 127 their left-skewed mixture with the arithmetic sequence $(\alpha = 128)$ $\beta = 0.328$, their right-skewed mixture with the arithmetic 129 sequence $(\alpha = \beta = 0.328)$ (Figure [2B](#page-2-0)). Besides beta sequences 130 with a U-shape, other sequences are paired so an additional 131 constraint is set to ensure equal weight for each pair. Besides 132 these 9 sequences and arithmetic sequences, a pseudo-random 133 sequence is introduced to further approximate the structure 134 and avoid inconsistent scenarios. Finally, a complement se- ¹³⁵ quence is introduced which if combining all the sequences with 136 corresponding weights, the overall sequence is nearly uniform. ¹³⁷

⁶⁶ for sample central moments.

Fig. 2. A. The first four sample central moments for the Gaussian distribution are plotted over a sample size ranging from 2 to 100. The red lines represent the expected values, while the blue lines depict the values estimated from the arithmetic sequences. B. The histograms of different beta sequences, their self-mixtures, and mixtures with arithmetic sequences.

Fig. 3. The first plot shows the weights assigned to different sequences as the sample size increases. The second plot depicts the sample standard deviations estimated from designed and arithmetic sequences and compares them to the true values. The designed sequences were repeated 10 times to reduce the variation due to the random sequences.

¹³⁸ **Results**

 The most surprising result in this article is that, by carefully selecting/designing sequences in *S*, even when *N* and *n* are very small, e.g., less than 20, the above system of linear equations can still be consistent, while the weight assigns to the random and complement sequences are extremely small $144 \left(< 0.01 \text{ on average} \right)$. Using just 12 sequences, when $n = 10$, the constraint optimization algorithm can assign weights to all these sequences with errors less than 10^{-10} . This means that technically, these sequences are consistent with the Gaussian distribution for the first four moments. More importantly, the findings suggest that when the sample size is small, the beta sequence with a U-shape accounts for approximately 50-60% of the finite sample properties of uniform random variables, while arithmetic, monotonic beta, beta-beta mixed, skewed beta distributions each contribute about 2-10% (Figure [3\)](#page-3-11). As the sample size grows, as expected, the weight of the arithmetic sequence increases and dominants while the weights of other sequences gradually decrease. However, the beta sequence with a U-shape still still holds about 10% weight even when the sample size is 100 (Figure [3\)](#page-3-11).

 The obtained weights can be used to estimate the finite sample behaviour of other related estimators, such as the stan- dard deviation and median absolute deviation for the Gaussian distribution. We found that by using the 12 well-designed sequences, the performance is much better than the arithmetic sequence (Figure [3\)](#page-3-11). To further increase precision, we adopted a stochastic method. We pseudo-randomly generated twelve sequences and evaluated their efficacy in approximating the finite sample structure of uniform random variables by solving the above system of linear equations for the first four moments. Sequences that met the predetermined accuracy threshold (er-170 ror less than 10^{-5}) were retained, while those that did not meet the requirement were discarded in favor of a new set. Upon identifying twenty qualified sets, these sets were ap- plied to assess the finite sample biases in other estimators for the Gaussian distribution. The outcomes indicate that using merely fifty sets of sequences, totaling 600 sequences, which can be executed on a standard PC in a negligible amount of time, achieves a precision of approximately 0.005 for the stan-dard deviation and median absolute deviation. In contrast, attaining the same level of precision using classic Monte Carlo ¹⁷⁹ methods would require roughly 0.1 million pseudo-random 180 samples. 181

Data and Software Availability 182

All data are included in the brief report and SI Dataset S1. 183 All codes have been deposited in [GitHub.](https://github.com/tubanlee/REDS_Nonasymptotic)

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