

Homework 2: Random Numbers and Multidimensional Monte-Carlo Integration

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Assignment 2

Random Numbers and Multidimensional Monte-Carlo Integration

2.1 Pseudo Random

1. The term "random number" is largely a misnomer when referring to numbers generated by machines. Machine algorithm is intrinsically deterministic and it is for this reason that numbers generated in such a manner are termed pseudo random numbers. Pseudo random number generators (PRNG) cannot be "truly" random, however, they may have the appearance of being so over a specific period. PRNG's can be assessed by their ability to mask their period of repetition when producing quantities of PRN's that are comparable to the generators period. In this section, the discussion of PRNG's will include examples from the Power-Residue method, `rand48()` and `random.org` distributions. Below, Figures 2.1 and 2.2 each provide "*PRNvis*" plots respectively.

Figure 2.1: 15,000 Pseudo Random Numbers using `rand48()`

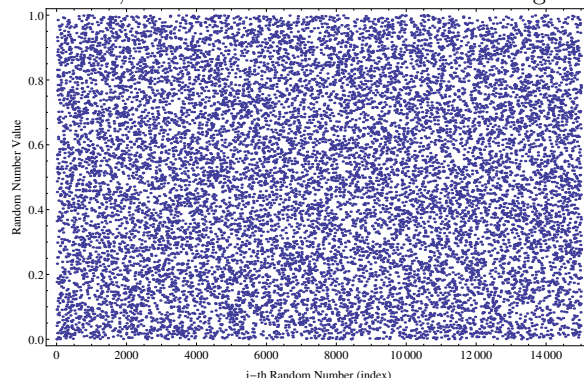
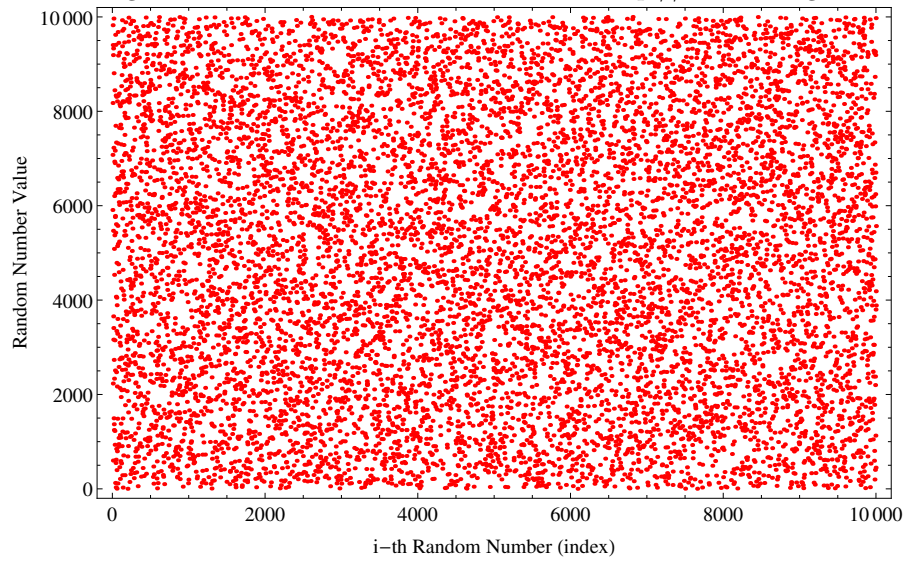


Figure 2.2: 10,000 Random Numbers from <http://random.org>

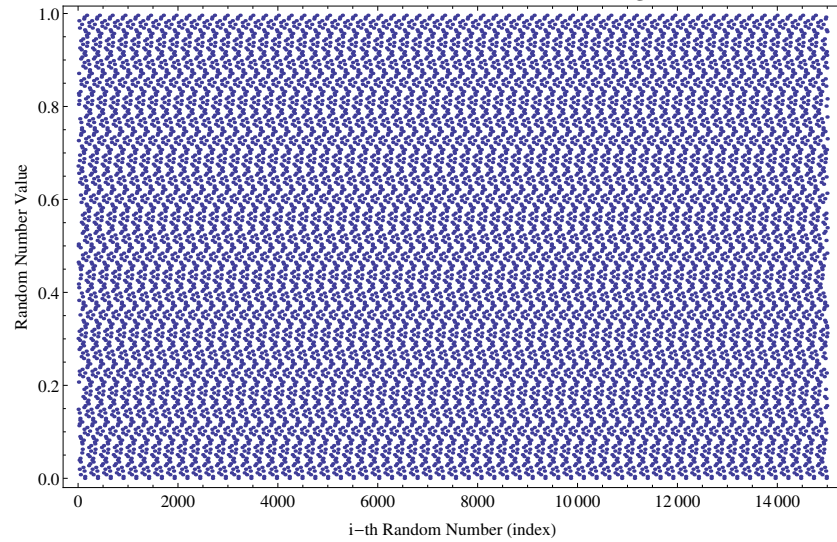


Qualitative Analysis:

The usefulness of a PRNG may be determined by assessing the uniformity and non-correlation of its plot(s). By "uniformity", it is meant that the plot of PRN's is uniformly distributed throughout its plot. By "non-correlation", it is meant that there appears no pattern or correlation between the PRN's. These criteria may be assessed qualitatively by the user visually inspecting the PRNG's plots for uniformity and non-correlation. Notice then that the `drand48()` and `random.org` plots in Figures 2.1 and 2.2 both pass the visual uniformity and non-correlation tests.

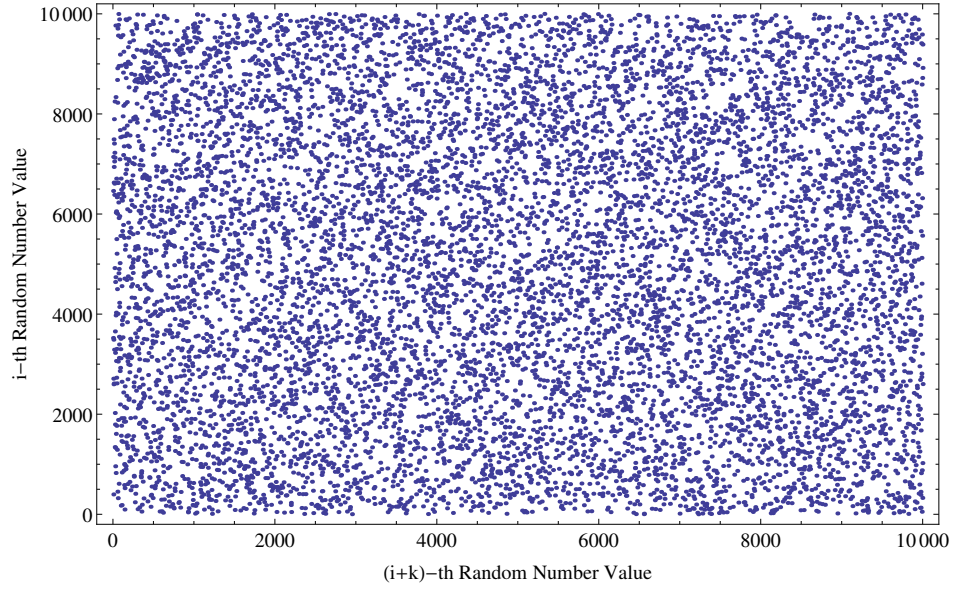
To contrast the successful tests above, Figure 2.3 depicts the Power-Residue PRNG failing the visual inspection. Notice that while the Power-Residue method appears well distributed, the correlation between its PRN's are visually simple to identify.

Figure 2.3: 15,000 Shameful Pseudo Random Numbers using Power-Residue method



Another method of qualitatively assessing non-correlation is to plot " x_i vs x_{i+k} " where x_i is the i^{th} random number and k is a "small" value which shifts the index. This plot may allow us to visually detect correlation that was not previously evident in the plot of " i vs x_i ". Figure 2.4 represents what the " x_i vs x_{i+k} " plots look like for both the drand48() and random.org distributions.

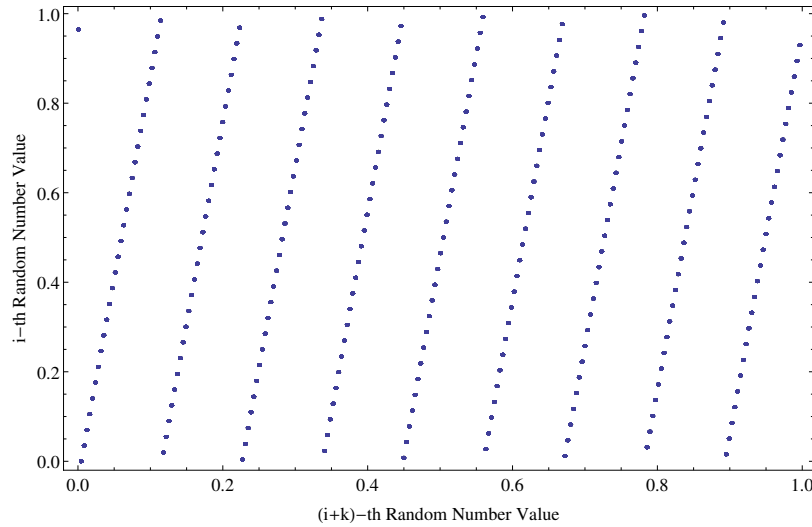
Figure 2.4: Depiction of " x_i vs x_{i+k} " plots for both drand48() and random.org



Notice that the " x_i vs x_{i+k} " plot appears identical to the original " i vs x_i " plots. Because of this, we gain confidence in each PRNG's ability to produce, "random", uniform distributions.

To contrast this this result, Figure 2.5 shows the Power-Residue PRNG's " x_i vs x_{i+k} " plot. We've already determined that the Power-Residue method fails the visual correlation criterion in Figure 2.3" and so the pattern evident in Figure 2.5 is expected.

Figure 2.5: Shameful " x_i vs x_{i+k} " plot for Power Residue Method



Histograms are another method by which the uniformity of a distribution may be visualized. This is done by dividing the sample range into intervals known as bins. For example, we used `drand48()` to calculate 15,000 PRN's in the range of $[0, 1]$. If we divide this range into 10 intervals, each bin in the x-axis of the histogram would be equal to 0.1. The distributions PRN's are each placed into their respective bin, whereby a histogram displaying the distributions uniformity is produced. If each bin has an equal number of PRN's, the top of the bins will all align and the distribution is considered uniform. The more offset between bins, the less uniform the distribution.

Figures 2.6 and 2.7 provide histograms of `drand48()`'s and `random.org`'s distributions. Notice the unevenness in `drand48()`'s histogram as opposed to `random.org`'s. These histograms suggest that `random.org`'s distribution is absolutely uniform whereas `drand48()`'s is close, but not perfect.

Figure 2.6: `drand48()` 10-Bin Histogram

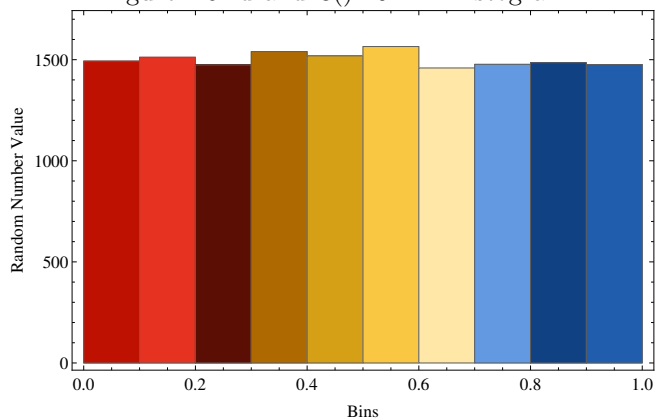
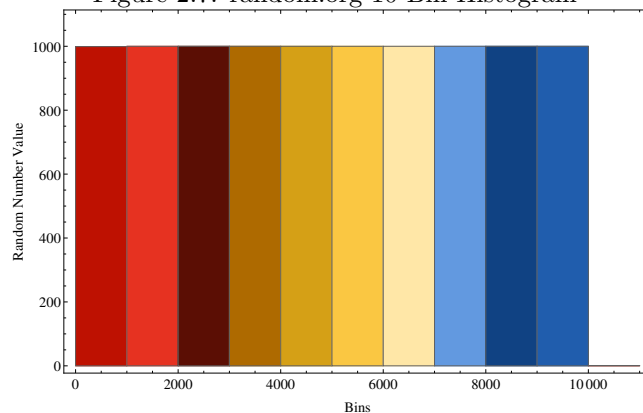


Figure 2.7: `random.org` 10-Bin Histogram



Quantitative Analysis:

The above methodologies take advantage of our ability to qualitatively assess uniformity and non-correlation. While these tests are effective and an ease to implement, there exist numerous quantitative measures as discussed below:

The k^{th} -moment:

The k^{th} moment test of a PRN distribution addresses uniformity. If a PRNG produces uniformly distributed PRN's, then the average of all of the PRN's raised to a sufficiently small k should approximate $\frac{1}{k+1}$. That is:

$$\frac{1}{N} \sum_{i=1}^n x_i^k \approx \frac{1}{k+1} \quad (2.1)$$

Note: If the standard deviation (average distance from the average point) varies as $\frac{1}{N}$, then the distribution may be considered "random".

Below are the k^{th} moment calculations for the rand48() and random.org distributions:

$$\begin{aligned} \cdot \quad \quad \quad rand48() : \quad & \frac{1}{1.5 \times 10^4} \sum_{i=1}^{1.5 \times 10^4} x_i^4 \approx \frac{1}{4+1} \\ \cdot \quad \quad \quad & 0.198068074284546 \approx 0.2 \end{aligned}$$

Relative Error between Calculated and Theoretical k^{th} moments = 0.0097538471

$$\begin{aligned} \cdot \quad \quad \quad random.org \quad & \frac{1}{1 \times 10^4} \sum_{i=1}^{1 \times 10^4} x_i^{4 \times 10^{-5}} \approx \frac{1}{4 \times 10^{-5} + 1} \\ \cdot \quad \quad \quad & 1.0003284904 \approx 0.9999600016 \end{aligned}$$

Relative Error between Calculated and Theoretical k^{th} moments = 0.0003685036

Notice here the significant decrease in error that random.org's distribution displays over rand48() –Nearly 2 orders of magnitude less error! These results depict the strong non-correlation advantage of the random.org distribution.

Auto Correlation (Near-neighbor Correlation):

Auto correlation tests non-correlation between PRN's within a distribution. If a PRNG produces uniformly distributed PRN's, then the average of all $x_i x_{i+k}$ PRN's should approximate $\frac{1}{4}$. More specifically, the formula for Auto Correlation is as follows:

$$\frac{1}{N} \sum_{i=1}^n x_i x_{i+k} \approx \frac{1}{4} \quad (2.2)$$

Note: The standard deviation (average distance from the average point) varying as $\frac{1}{N}$ ensures that the distribution is "random".

Below are Auto Correlation calculations for the rand48() and random.org distributions:

$$\begin{aligned} \cdot \quad & rand48() : \quad \frac{1}{1.5 \times 10^4} \sum_{i=1}^{1.5 \times 10^4} x_i x_{i+1} \approx \frac{1}{4} \\ \cdot \quad & \quad \quad \quad 0.248318844894805 \approx 0.25 \\ \cdot \quad & \quad \quad \quad \text{Relative Error between Calculated and Theoretical Auto Correlations} = 0.006724620420 \end{aligned}$$

$$\begin{aligned} \cdot \quad & random.org : \quad \frac{1}{1 \times 10^4} \sum_{i=1}^{1 \times 10^4} x_i x_{i+0.00004} \approx \frac{1}{4} \\ \cdot \quad & \quad \quad \quad 0.248989071076 \approx 0.25 \\ \cdot \quad & \quad \quad \quad \text{Relative Error between Calculated and Theoretical Auto Correlations} = 0.004043715696 \end{aligned}$$

Notice here the decrease in error that random.org's distribution displays over drand48() – While the difference in error is not as significant as in the k^{th} moment calculations, the results still display a non-correlation advantage in random.org's distribution.

2.2 Multidimensional Monte-Carlo Integration

1. The Monte-Carlo integration method is a technique used for numerically approximating the area under an n-dimensional function. The method is implemented much alike the mean value theorem. The two methods are similar in that they both use the limits of integration $(b - a)$ as the width in the area calculation. Where the methods differ is in Monte-Carlo's application of random numbers to determine the height in the area calculation. The height is determined by taking N evenly distributed random numbers, inputting them into the given function, and averaging their resulting values.

The formula for a multidimensional Monte Carlo integral is as follows:

$$\int_a^b \int_c^d \int_e^f f(x_1, x_2, x_3) dx_1 dx_2 dx_3 \approx (b - a)(d - c)(f - e) \frac{1}{N} \sum_{i=0}^N f(x_i) \quad (2.3)$$

Here, " $f(x_1, x_2, x_3)$ " is a surface in \mathbb{R}^3 , the variables in " $(b - a)(d - c)(f - e)$ " correlate directly to the limits of integration and " N " represents the number of iterations/random numbers used to determine a mean value for $f(x_1, x_2, x_3)$.

As an example, consider the following 5D integral:

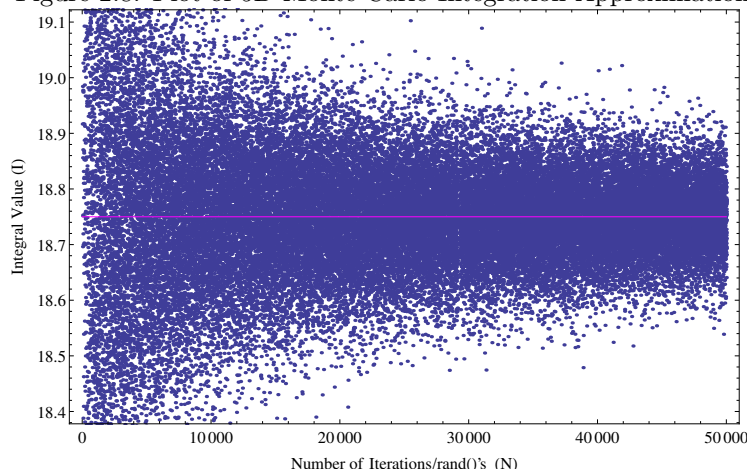
$$I = \int_0^1 \int_0^1 \int_0^1 \int_0^1 \int_0^1 (x_1 + x_2 + x_3 + x_4 + x_5)^3 dx_1 dx_2 dx_3 dx_4 dx_5 = 18.75 \quad (2.4)$$

Using Monte-Carlo integration, we chose $N = 5 \times 10^4$ random numbers to approximate $\int_0^1 \int_0^1 \int_0^1 \int_0^1 \int_0^1 f(x_1 \dots x_5)^3 dx_1 \dots dx_5$ as:

$$\approx (1 - 0)(1 - 0)(1 - 0)(1 - 0)(1 - 0) \left(\frac{1}{5 \times 10^4} \right) \sum_{i=0}^{5 \times 10^4} f(x_{i1} + x_{i2} + x_{i3} + x_{i4} + x_{i5})^3 \approx 18.75296845$$

The relative error between the Approximate and Calculated values of the integral is $1.583175277 \times 10^{-4}$.

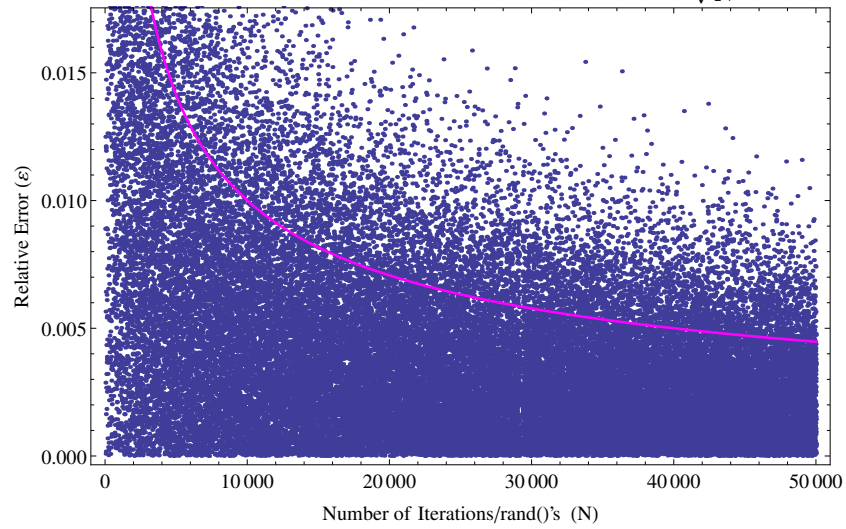
Figure 2.8: Plot of 5D Monte-Carlo Integration Approximation



Monte-Carlo Error:

Error in the Monte-Carlo method decreases, independent of dimension, as $\frac{1}{\sqrt{N}}$. This is evident Figure 2.9 below:

Figure 2.9: Error in 5D Monte-Carlo Integration(blue) and $\frac{1}{\sqrt{N}}$ (magenta)



Notice how the plot of Relative-Error vs. Iterations(N) behaves as $\frac{1}{\sqrt{N}}$.

Finally, the variance of function is defined as the standard deviation squared, that is, σ^2 .

Here, $\sigma^2 = \langle f^2 \rangle - \langle f \rangle^2 = 0.061207254801$.