Modeling Random Telegraph Noise as a Randomness Source and Its Application in True Random Number Generation

Xiaoming Chen, Member, IEEE, Lin Wang, Boxun Li, Student Member, IEEE, Yu Wang, Senior Member, IEEE, Xin Li, Senior Member, IEEE, Yongpan Liu, Member, IEEE, and Huazhong Yang, Senior Member, IEEE

Abstract—The random telegraph noise (RTN) is becoming more serious in advanced technologies. Due to the unpredictability of the physical phenomenon, RTN is a good randomness source for true random number generators (TRNG). In this paper, we build fundamental randomness models for TRNGs based on single trap- and multiple traps-induced RTN. We theoretically derive the autocorrelation coefficient, bias, and hit rate for RTN-based TRNGs. Two representative RTN-based TRNG schemes are simulated to verify the proposed randomness models. An oscillator-based TRNG is also studied based on the theoretical randomness model of multiple traps-induced RTN. We also provide basic guidelines for designing RTN-based TRNGs.

Index Terms—Random telegraph noise, true random number generator, randomness modeling

I. INTRODUCTION

The random telegraph noise (RTN) is a growing reliability issue in advanced integrated circuit (IC) technologies. RTN causes random fluctuations in electrical parameters such as the threshold voltage ($V_{th}$) and the source-drain current ($I_{ds}$). Recent studies have shown that at the 22nm node, the RTN-induced $V_{th}$ fluctuation can be larger than 70mV so RTN becomes a major noise source [1], [2]. RTN significantly affects the reliability of modern ICs. However, on the other hand, since the physical phenomenon of RTN is unpredictable, RTN potentially provides an excellent randomness source for creating true random number generators (TRNG). TRNG is an essential foundation in lots of cryptographic algorithms. Using pseudo random numbers will cause a big vulnerability because of the predictability. On the contrary, if the randomness source is unpredictable, like RTN, the generated random numbers will also be unpredictable. This is why TRNGs are essentially demanded for security.

A TRNG is typically composed of three major modules: entropy source, harvester, and post-processing. The entropy source provides raw random signals which are extracted from unpredictable physical phenomena. Raw signals are usually analog and biased. The harvester converts the raw signals to a digital bit stream. The post-processing is used to reduce bias to balance the probabilities of zeros and ones in the output.

A wide range of physical phenomena can be adopted to act as the entropy source, such as device noises, clock jitter, metastability, chaos, and etc. It is claimed that some conventional entropy sources cannot offer high randomness, such as the clock jitter [3]. On the other hand, reliability mechanisms-based TRNGs are claimed to provide higher randomness [4], [5]. As a growing reliability mechanism, RTN offers much larger random fluctuations than well-known device noises, providing an excellent entropy source for TRNGs. In this paper, we will investigate the methodology of utilizing RTN as an entropy source in TRNGs. The purpose is to build fundamental randomness models for RTN and give basic guidelines to RTN-based TRNG design. To achieve this goal, we build systematic models to evaluate the randomness of RTN in theory. We also give basic methodologies to eliminate the autocorrelation and ensure high randomness for TRNGs based on both single trap- and multiple traps-induced RTN. Advantages of RTN-based TRNGs are emphasized by comparisons with conventional noise- and clock jitter-based TRNGs.

The rest of this paper is organized as follows. We review the related work and present our motivation in Section II. Statistical modeling and simulation methodology of RTN are introduced in Section III. In Sections IV and V, we present randomness models for single trap- and multiple traps-induced RTN, respectively. In Section VI, we study an oscillator-based TRNG using the proposed models. In Section VII, we compare RTN-based TRNGs with conventional noise- and clock jitter-based TRNGs. Finally, Section VIII concludes the paper.

II. RELATED WORK AND MOTIVATION

In this section, we first briefly review some related work, and then present the motivation of this work, followed by a
summarized description of the proposed models.

A. Randomness Modeling of Noise

Kirton et al. gave a fundamental introduction on the physical origin and statistical characteristics of RTN based on the trap switching theory in [6]. Several studies built theoretical models for the jitter and phase noise in ring oscillators (RO) by modeling the white noise and $1/f$ noise [7]–[9]. White noise was theoretically modeled to generate random numbers in [10].

For RTN, although many studies have developed statistical models based on measured data (e.g., [11]–[15]), they focus on modeling the physical phenomenon of RTN. Currently, there is no research that gives a systematic study on randomness modeling of RTN.

B. Existing TRNG Designs

According to the entropy source, there are several types of TRNGs. Entropy sources adopted by popular TRNGs include noises [16], [17], clock jitter in free-running ROs [18], [19], metastability [20], and chaos [21]. The breakdown mechanism of metal-oxide-semiconductor field-effect transistors (MOSFET) has also been studied to generate random numbers [4]. The only two RTN-based TRNGs are proposed in [5], [22].

Almost all of these publications only provide implementations without any theory base on the randomness, naturally raising a question: is the randomness of such implementations really high enough? For example, the jitter-to-mean period ratio is at the magnitude of $10^{-4}$ in free-running ROs [18], such that the RO period must be very long to make a big jitter to create TRNGs. It is claimed that conventional noise- and metastability-based TRNGs cannot provide high randomness, due to the low magnitude of randomness or mismatch of devices [3]. In addition, whether a hard-to-describe chaotic system really behaves in a physically random fashion is unclear [3].

C. Motivation

As can be seen from the above, most of the existing TRNG designs lack for a theoretical derivation of the randomness. Utilizing RTN as a randomness source in TRNGs has two prominent advantages. As a growing reliability issue, RTN offers significantly large random fluctuations in advanced technologies, so that the fluctuations can be easily extracted and converted to random bits. In addition, more bits can be generated from each sampling due to the large fluctuation magnitude. We will further analyze the advantages of RTN-based TRNGs in Section VII. The physical phenomenon of RTN has been well modeled, but the randomness of RTN has never been systematically studied. This paper will build fundamental randomness models for RTN to provide a theoretical foundation for designing RTN-based TRNGs. We will focus on analyzing the autocorrelation coefficient, bias, and bit rate of RTN-based TRNGs. Among them, making a zero bias is the minimum requirement for random bits. However, only reducing the bias cannot guarantee the randomness. For example, a sequence $\{1, 0, 1, 0, 1, 0, 1, 0, \cdots\}$ has no bias but it is not random at all. Therefore, the autocorrelation which has a large impact on the randomness is also analyzed. A near-zero autocorrelation indicates that the next bit is difficult to predict when the previous bit is known. Bit rate is an important performance metric. Although there are other metrics or methods which can also be adopted to evaluate the randomness, such as the information entropy and the test suite provided by the National Institute of Standards and Technology (NIST) [23], they are high-level scores without tight connections with the physical characteristics of random bits. We choose autocorrelation, bias, and bit rate because they have clear physical meanings and crucial influences on the randomness and performance of TRNGs.

D. Key Points of Proposed Models

For TRNGs based on single trap-induced RTN, we will show how to select the sampling frequency to ensure a small autocorrelation. The maximum sampling frequency is constrained by the time constants of the trap. The bias can be eliminated by post-processing like the von Neumann corrector [24]. For TRNGs based on multiple traps-induced RTN, the sampling frequency can be close to the switching frequency of the fastest trap. We will show how to design a bit truncation scheme to eliminate the bias and the high autocorrelation.

III. STATISTICAL MODELING AND SIMULATION METHODOLOGY OF RTN

In this section, we introduce existing statistical models and our simulation methodology of RTN.

A. Statistical Modeling of RTN

1) Physics of RTN: RTN can be explained by the random capture/emission process of charge carriers caused by oxide traps [25], as shown in Fig. 1. A trap in the oxide can occasionally capture a charge carrier from the channel, and the captured carrier can be emitted back to the channel after a period of time. The duration time of the captured and emitted states are denoted as $\tau_e$ (time before emission) and $\tau_r$ (time before capture), respectively, as marked in Fig. 2a. In the time
domain, $V_{th}$ shows a binary fluctuation caused by a single trap. In the frequency domain, the power spectrum density (PSD) of the RTN-induced $V_{th}$ fluctuation shows a Lorentzian shaped spectrum with a slope of $\frac{1}{\tau^2}$ [6], as shown in Fig. 2b.

2) Time Constants: The switching process of an individual trap obeys a Poisson process [6]. As a result, the duration time of the captured and emitted states follow exponential distributions with mean values $\bar{\tau}_c$ and $\bar{\tau}_e$, respectively [12]:

$$f(\tau_c) = \frac{1}{\bar{\tau}_c} e^{-\frac{\tau_c}{\bar{\tau}_c}}, f(\tau_e) = \frac{1}{\bar{\tau}_e} e^{-\frac{\tau_e}{\bar{\tau}_e}}.$$  \hspace{1cm} (1)

$\bar{\tau}_c$ and $\bar{\tau}_e$ are called the capture and emission time constants, respectively. They have a wide range from microsecond to millisecond when sampling a number of traps [11], [12]. Time constants of numerous traps can be approximated modeled by uniform distributions in the logarithmic scale [6], [12], [26]:

$$\log_{10}(\bar{\tau}_c) \sim U(A_c, B_c), \log_{10}(\bar{\tau}_e) \sim U(A_e, B_e).$$  \hspace{1cm} (2)

where $U(A, B)$ denotes the uniform distribution in the interval $(A, B)$. Eq. (2) is a statistical model for a number of traps. For each individual trap, its $\bar{\tau}_c$ and $\bar{\tau}_e$ are strongly correlated [12]. Although an analytical relation between $\bar{\tau}_c$ and $\bar{\tau}_e$ has been given in [6], [11], it relies on some low-level parameters which are difficult to obtain and model. For convenience, the model can be simplified to:

$$\bar{\tau}_e = 10^m, m \sim U \left( \bar{m} - \frac{\sigma_m}{2}, \bar{m} + \frac{\sigma_m}{2} \right),$$  \hspace{1cm} (3)

where $\bar{m}$ and $\sigma_m$ are fitting parameters. $\bar{m}$ is linear to the bias voltage $V_{gs}$ which indicates that $\bar{\tau}_c$ and $\bar{\tau}_e$ are approximately exponential to $V_{gs}$ [11]. Using $\sigma_m = 2$ can generally fit the silicon data presented in [12]. The randomness of $m$ denotes the statistical characteristics of trap positions and energies.

The above model is applicable in the case of a constant bias condition. Actually, $\bar{\tau}_c$ and $\bar{\tau}_e$ both depend on the bias condition. When a transistor undergoes periodically alternate two states (ST1 and ST2) with a fixed duty cycle, the periodic behavior of RTN can be described by an equivalent stationary RTN with two equivalent time constants [13]:

$$\frac{1}{\bar{\tau}_c} = \frac{\alpha}{\bar{\tau}_c^{(eq)}(ST1)} + \frac{1 - \alpha}{\bar{\tau}_c^{(eq)}(ST2)}, \frac{1}{\bar{\tau}_e} = \frac{\alpha}{\bar{\tau}_e^{(eq)}(ST1)} + \frac{1 - \alpha}{\bar{\tau}_e^{(eq)}(ST2)},$$  \hspace{1cm} (4)

where $\alpha$ is the duty cycle. $\bar{\tau}_c^{(eq)}(ST1)$ and $\bar{\tau}_c^{(eq)}(ST2)$ are fixed if $\alpha$ is fixed. Consequently, the model with fixed time constants can also be used in the case of a cyclostationary state.

3) Number of Traps: For numerous transistors, the number of detectable traps in each transistor statistically obeys a Poisson distribution [14]:

$$P(N_t = k) = \frac{\langle N \rangle^k e^{-\langle N \rangle}}{k!},$$  \hspace{1cm} (5)

where $N_t$ is the number of detectable traps in a transistor $t$, and $\langle N \rangle$ is the mean value of the Poisson distribution. pMOSFETs have more traps than nMOSFETs [14]. $\langle N \rangle$ increases with the shrinking of the feature size [27].

4) $V_{th}$ Amplitude: The RTN-induced $V_{th}$ fluctuation of numerous traps statistically obey an exponential distribution [15]:

$$f(\Delta V_{th,i}) = \frac{1}{\langle \Delta V_{th} \rangle e^{-\frac{\Delta V_{th,i}}{\langle \Delta V_{th} \rangle}}},$$  \hspace{1cm} (6)

where $\Delta V_{th,i}$ is the $V_{th}$ fluctuation caused by a trap $i$, and $\langle \Delta V_{th} \rangle$ is the mean value of the exponential distribution. $\langle \Delta V_{th} \rangle$ increases with the shrinking of the feature size [15]. In this paper, we set $\langle \Delta V_{th} \rangle$ according to the 22nm and 32nm silicon data presented in [2], [28]. The total $V_{th}$ fluctuation of a transistor $t$ is the superposition of the effects of all the individual traps in the transistor [29]:

$$\Delta V_{th} = \sum_{i=1}^{N_t} (\Delta V_{th,i} \times S_i),$$  \hspace{1cm} (7)

where $S_i \in \{0, 1\}$ indicates the state of the $i$th trap in the transistor (0 for the emitted state and 1 for the captured state).

B. Simulation Methodology

Since currently there is no available circuit simulator which can natively support RTN, we build an in-house simulator for RTN estimation. Our simulator is a standard simulation program with integrated circuit emphasis (SPICE) [30]-based tool with the BSIM4 [31] model integrated. We follow the approach proposed in [29] to integrate the RTN models presented in Section III-A into the simulator. The simulation flow is shown in Fig. 3, where shaded blocks are RTN-related.

Before transient simulation, the RTN profile is given or randomly generated using the following three steps.

- The number of traps in each transistor $N_t$ is given, or generated by Eq. (5) where $\langle N \rangle$ is given;
- The $V_{th}$ fluctuation of each trap $\Delta V_{th,i}$ is given, or generated by Eq. (6) where $\langle \Delta V_{th} \rangle$ is given;
- The capture time constant of each trap $\bar{\tau}_c,i$ is given, or generated by Eq. (2) where the range is given. The emission time constant of each trap $\bar{\tau}_e,i$ is given, or generated by Eq. (3) where $\bar{m}$ is given.

Note that each trap has its own time constants and $V_{th}$ amplitude [29]. As shown in Fig. 3, each trap $i$ keeps a parameter $t_{next,i}$ which indicates when it will change its state $S_i$. During transient simulation, once the time point $t$ reaches $t_{next,i}$, trap $i$ changes its state $S_i$, and then $t_{next,i}$ is
updated by the duration time of the next state, which obeys an exponential distribution. In simulation, the duration time of the next state is randomly generated by the following approach [29]:

$$\tau_e = -\ln(RND_{(0,1)}) \times \bar{\tau}_e, \quad \tau_c = -\ln(RND_{(0,1)}) \times \bar{\tau}_c,$$ (8)

where $RND_{(0,1)}$ denotes a uniformly distributed random number in the interval $(0, 1)$. The RTN-induced $\Delta V_{th}$ of each transistor is calculated by Eq. (7) and added to the original $V_{th}$ of each transistor, and then the BSIM4 model evaluation is performed based on the updated $V_{th}$. To avoid missing any state change during transient simulation, the time point is always not larger than the minimum value of all the $t_{\text{next,i}}$'s.

The accuracy of our simulator have been verified by comparing waveforms with commercial tools. Here we verify the simulated PSD of single trap-induced RTN. In this test, we use the 22nm high-performance predictive technology model (PTM) [32] in our netlists. A circuit with a single nMOSFET (the width is 50nm) which has a single trap-induced RTN effect is simulated. The nMOSFET is stressed by a fixed bias condition of $V_{gs} = 0.6V$ and drives a load resistance of 10kΩ. We use a fixed RTN profile in this test: $\bar{\tau}_e = \bar{\tau}_c = 0.5ms$, $\Delta V_{th} = 20mV$, and $N_t = 1$. Fig. 4 shows the simulated PSD of $I_{ds}$, and the theoretical Lorentzian PSD which is given by [6]

$$S(f) = \frac{4(\Delta I_{ds})^2}{\tau_e + \tau_c} \frac{1}{1 + (2\pi f \tau_0)^2},$$ (9)

where

$$\frac{1}{\tau_0} = \frac{1}{\tau_e} + \frac{1}{\tau_c}.$$ (10)

$\Delta I_{ds} \approx 2.1\mu A$ is obtained from the simulated $I_{ds}$ waveform. Fig. 4 proves that the simulated PSD is well consistent with the theoretical Lorentzian spectrum.

IV. MODELING SINGLE TRAP-INDUCED RTN

In this section, we first derive a theoretical randomness model for single trap-induced RTN, and then analyze the performance of a representative TRNG scheme based on single trap-induced RTN.

Since the fluctuation caused by single trap-induced RTN has only two discrete values, the simplest way to generate random numbers is to convert the fluctuation into binary bits by a periodically sampled comparator, just like the approach proposed in [5]. Fig. 5 shows a representative scheme for this method. An amplifier may be used if the RTN-induced fluctuation is not large enough. Actually the implementation can be flexible and the binary bits can be converted from other parameters which are affected by RTN. Other implementations based on the same theory are equivalent to the representative scheme.

A. Randomness Modeling

1) Autocorrelation: Let $X$ be the state of a trap. $X = 0/1$ indicates the emitted/captured state. The probabilities of the two states in a stationary state are given by

$$P(X=1) = P_1 = \frac{\bar{\tau}_e}{\bar{\tau}_e + \bar{\tau}_c}, \quad P(X=0) = P_0 = \frac{\bar{\tau}_c}{\bar{\tau}_e + \bar{\tau}_c}.$$ (11)

Considering two time points $s$ and $s + t$, the transition probabilities of a trap, which are also the prediction probability $P(X_{s+t}|X_s)$, are given by [33]

$$P(X_{s+t}=1|X_s=1) = P_{11} = P_1 + P_0 e^{-\frac{t}{\tau_0}},$$

$$P(X_{s+t}=0|X_s=1) = P_{10} = P_0 - P_0 e^{-\frac{t}{\tau_0}},$$

$$P(X_{s+t}=1|X_s=0) = P_{01} = P_0 + P_1 e^{-\frac{t}{\tau_0}},$$

$$P(X_{s+t}=0|X_s=0) = P_{00} = P_1 - P_1 e^{-\frac{t}{\tau_0}},$$ (12)

where $P_{ij}$ means the probability of ending at state $j$ after an elapsed time $t$, when starting from state $i$. In practice, $t$ is the sampling period $t_s = \frac{1}{f_s}$, where $f_s$ is the sampling frequency. According to Eq. (12), when $t_s$ is long enough (e.g., $t_s > 3\tau_0$), the prediction probabilities are close to the stationary state probabilities $P_0$ and $P_1$, which means that knowing the previous state does not provide useful information for predicting the next state.

Let $\Delta A$ be the RTN-induced fluctuation (e.g., $\Delta V_{th}$). Without loss of generality, we assume that $\Delta A$ is zero-mean. The autocorrelation function of the RTN-induced fluctuation is expressed as

$$C(s, s + t_s) = E(X_s X_{s+t_s}) = (\Delta A)^2 \left( P_{01}^2 P_{11} - P_{00}^2 P_{10} - P_{01}^2 P_{10} + P_{01}^2 P_{01} + P_{11}^2 P_{00} \right)$$

$$= (\Delta A)^2 \frac{\tau_0}{\bar{\tau}_e + \bar{\tau}_c} e^{-\frac{t_s}{\tau_0}}.$$ (13)

The autocorrelation function only depends on the duration time $t_s$ so it can be written as $C(t_s)$. The first-order autocorrelation coefficient is given by

$$\rho(t_s) = \frac{C(t_s)}{C(0)} = e^{-\frac{t_s}{\tau_0}}.$$ (14)

Increasing $t_s$ can decrease the autocorrelation, which is consistent with the prediction probabilities shown in Eq. (12). To ensure high randomness, $t_s$ should be high enough. For example, if $t_s \geq 3\tau_0$, the autocorrelation coefficient is less than
The autocorrelation may be partly eliminated by applying some post-processing methods, so the sampling frequency can be higher. According to Eq. (11), $P_0 = P_1$ if and only if $\bar{\tau}_e = \bar{\tau}_e$, which is impossible in actual devices. As a result, post-processing is always required. In this paper, we will take the popular von Neumann corrector [24] as an example to derive the autocorrelation, bias, and bit rate. The von Neumann corrector outputs “0” or “1” if two successive input bits are “01” or “10”, but discards “00” and “11”. After applying the von Neumann corrector, the autocorrelation coefficient can be approximated by

$$\rho_{eN}(t_s) \approx \frac{1}{e^{-\frac{2}{\tau_0}} + 4 + \frac{3}{2P_0P_1 - 1}}, t_s > 1.5\tau_0. \tag{15}$$

The derivation is complicated so we put it in Appendix A. The autocorrelation after the von Neumann corrector is applied is less than the original value given by Eq. (14). For example, if $f_s = \frac{1}{\tau_0}$, Eq. (15) gives an autocorrelation of about 2.5%.

Till now, by deriving the autocorrelation coefficient, we have proved that when using a proper sampling frequency, the autocorrelation can be quite small such that the next bit cannot be predicted when the previous bit is known. This proves the randomness of single trap-induced RTN. Eq. (15) can be used to estimate the proper sampling frequency when the von Neumann corrector is used.

2) Bias: Let the probabilities of ones and zeros are $0.5 + b$ and $0.5 - b$, respectively, where $b$ is the bias, i.e.,

$$b = \frac{\bar{\tau}_e - \bar{\tau}_e}{2(\bar{\tau}_e + \bar{\tau}_e)}. \tag{16}$$

The bias after the von Neumann corrector is applied is expressed as

$$b_{eN} = P(\text{output “1” has output}) - 0.5 = \frac{P_1P_0}{P_1P_10 + P_0P_01} - 0.5 \equiv 0. \tag{17}$$

As can be seen, the bias is completely eliminated, regardless of the sampling frequency and the original bias.

3) Bit Rate: For the von Neumann corrector, the ratio of the output bit rate to the sampling frequency equals half of the probability of observing “10” or “01”, which is given by

$$R_{eN}/f_s = \frac{1}{2} \left(P_0P_10 + P_1P_01\right) = \frac{\tau_0}{\bar{\tau}_e + \bar{\tau}_e} \left(1 - e^{-\frac{t_s}{\tau_0}}\right)$$

$$\approx \frac{1}{4} - b^2 \left(1 - \rho(t_s)\right) < \frac{1}{4}, \tag{18}$$

where $R_{eN}$ is the bit rate of the von Neumann corrector. $R_{eN}$ depends on both the original bias and autocorrelation. In any case, the bit rate is less than $\frac{1}{4}$ of the sampling frequency. $\bar{\tau}_e = \bar{\tau}_e$ (i.e., $b = 0$) yields the maximum bit rate. Applying the 1st-order Taylor expansion to Eq. (18) yields

$$R_{eN} < \frac{1}{4f_s} \left(1 - e^{-\frac{t_s}{\tau_0}}\right) \approx \frac{1}{4} \cdot \frac{t_s}{4\tau_0} = \frac{1}{4\tau_0}. \tag{19}$$

Eq. (19) gives the maximum ideal bit rate that the von Neumann corrector can achieve. It is achieved only when $\bar{\tau}_e = \bar{\tau}_e$ and $f_s$ is high enough. However, using a high $f_s$ leads to a high autocorrelation, so the maximum ideal bit rate cannot be achieved in practice.

### B. Randomness Source Selection

As mentioned in Section III-A, RTN has a big uncertainty and the time constants have a wide range. It is difficult to make a specific transistor behave as what we expect. A feasible solution is to select an adequate transistor from a large transistor array [5]. We should select a transistor with exactly one observable trap, small time constants, and significant $\Delta V_{th}$. To derive the probability of finding an adequate transistor, we first need the probability density function (PDF) of $\tau_0$.

According to Eq. (2), the PDF of $\tau_0$ is expressed as

$$pdf(\tau_0) = \frac{1}{\tau_0 \ln \frac{\tau_{0,\max}}{\tau_{0,\min}}}, \tau_{0,\min} \leq \tau_0 \leq \tau_{0,\max}, \tag{20}$$

where $\tau_{0,\max}$ and $\tau_{0,\min}$ are the maximum and minimum values of $\tau_0$, respectively. Although $\bar{\tau}_e$ is generated according to Eq. (3) with a small randomness on $m$, for simplicity in deriving the model, here we assume that $m$ is fixed so $\bar{\tau}_e$ is linear with $\bar{\tau}_e$. Consequently, $\tau_0$ is also linear with $\bar{\tau}_e$.

$$\tau_0 = \frac{10^m}{10^m + 1} \bar{\tau}_e = \beta \bar{\tau}_e. \tag{21}$$

Then the PDF of $\tau_0$ is expressed as

$$pdf(\tau_0) = \frac{1}{\tau_0 \ln \frac{\tau_{0,\max}}{\tau_{0,\min}}}, \tau_{0,\min} \leq \tau_0 \leq \tau_{0,\max}, \tag{22}$$

where $\tau_{0,\max} = \beta \tau_{c,\max}$ and $\tau_{0,\min} = \beta \tau_{c,\min}$.

Assuming that we have $N$ transistors in total, the probability of finding at least one transistor, such that it has exactly one trap, $\tau_0 \in [\tau_{0,\min}, \delta \tau_{0,\min}]$, and $\Delta V_{th} \geq \gamma (\Delta V_{th})$, is expressed as

$$P = \left(1 - \left(1 - \frac{\delta \tau_{0,\min}}{\tau_{0,\min}} \langle N_t \rangle e^{-\langle N_t \rangle} e^{-\gamma} \right)^N \right)^{-1} \tag{23}$$

$N$ can be solved from Eq. (23) when $P$ is given. For example, if $\delta = 2$, $\frac{\tau_{0,\max}}{\tau_{0,\min}} = 10^4$, $\gamma = 2$, and $\langle N_t \rangle = 0.8$, then using $N \geq 1884$ can ensure a probability of larger than 0.999.

### C. Numerical Results

To verify the proposed randomness model of single trap-induced RTN, the TRNG scheme shown in Fig. 5 is simulated using the 22nm PTM [32]. The nMOSFET is affected by a single trap. We use a fixed RTN profile in this test: $N_t = 1$, $\tau_0 = 10\mu$s, and $\Delta V_{th} = 40$mV. We will analyze the performance under different $\bar{\tau}_e$ and $\tau_0$ (Eq. (10) is always satisfied). We adopt the concept of the approximate entropy (ApEn) [34] to evaluate the generated random numbers with the von Neumann corrector applied. All the results reported in this subsection are the mean values of 5 runs.

Fig. 6 shows the simulated ApEn, under different $\frac{\tau_e}{\tau_0}$ and $f_s$. Note that the maximum ideal ApEn is $\ln(2) \approx 0.69315$. As can be seen, ApEn is higher than 0.69 only when $f_s = 20$kHz and 45kHz. Increasing $f_s$ greatly decreases ApEn. ApEn almost keeps constant when $\tau_0$ varies, revealing that the randomness is mainly determined by $\tau_0$ and $f_s$. The simulated
The generated random numbers are used to evaluate the value of \( \pi \) by the Monte Carlo (MC) method, as shown in Fig. 8.

**Fig. 8: Evaluated \( \pi \) by the Monte Carlo method.**

The bit rate is always less than \( \frac{1}{22r_0} \). The bit rate decreases significantly when \( \frac{e}{r_e} \) is far away from 1.0.

**D. Summary**

In this section, we have analyzed the autocorrelation coefficient, bias, and bit rate of a representative TRNG scheme based on single trap-induced RTN. As predicted by the proposed model and verified by the numerical results, using a too high sampling frequency greatly decreases the randomness of the output bits. In practice, the maximum sampling frequency is approximately \( \frac{1}{2\bar{r}} \), and the corresponding output bit rate is about \( \frac{0.4}{\bar{r}_c} \). Selecting a trap with almost equal \( \bar{r}_c \) and \( \bar{r}_e \) maximizes the output bit rate. A large transistor array should be constructed to ensure that we can always find an adequate transistor to act as the randomness source.

**V. MODELING MULTIPLE TRAPS-INDUCED RTN**

In this section, we first derive a theoretical randomness model for multiple traps-induced RTN, and then analyze the performance of a representative TRNG scheme based on multiple traps-induced RTN.

A simple idea to generate random numbers from multiple traps-induced RTN is to combine several individual circuits shown in Fig. 5 in parallel such that multiple bits can be generated in one sampling. However, these are two practical problems. First, the time constants of many traps have a wide range so it is difficult to synchronize all the circuits using a unified sampling frequency. Second, the number of observable traps in each transistor is random, so some transistors do not show any fluctuation and some may show more-than-two-level fluctuations. In the single trap case, the two problems do not appear because we can select an adequate transistor from a large transistor array. Consequently, the effects of all the individual traps should be combined together to act as a single randomness source. Statistical laws and stochastic process theories will ensure that the overall RTN effect of numerous traps obeys a certain statistical rule.

**Fig. 10: Representative scheme for generating random numbers from multiple traps-induced RTN.**

The bit rates are well consistent with the predictions. As predicted by Eq. (18), the bit rate is always less than \( \frac{1}{22r_0} \). The bit rate decreases significantly when \( \frac{e}{r_e} \) is far away from 1.0.

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**Fig. 6: ApEn of random numbers generated from single trap-induced RTN.**

**Fig. 7: Simulated and theoretical autocorrelation coefficients.**

**Fig. 9: Output bit rates of the von Neumann corrector.**

**Fig. 10: Representative scheme for generating random numbers from multiple traps-induced RTN.**
only two discrete levels, the superposed fluctuation will have many discrete levels so it will look like a continuous signal. The superposed fluctuation is amplified and converted to digital words. The converted digital words have bias and autocorrelation, so the von Neumann corrector may also be applied, leading to a low bit rate. Considering the fact that in the converted digital words, high-order bits change slow and low-order bits change fast, low-order bits tend to be more random. This is a special feature of TRNGs based on multiple traps-induced RTN. In this section, we will investigate that by truncating a few high-order bits from the digital words, the remaining bits will have near-zero autocorrelation and bias. We will also show that the bit truncation scheme has a higher bit rate than the von Neumann corrector for TRNGs based on multiple traps-induced RTN.

A. Superposition of Multiple Lorentzian PSDs

We will first derive the PSD caused by multiple traps-induced RTN based on the statistical RTN model in Section III-A. The PSD of multiple traps-induced RTN is a superposition of PSDs of all the individual traps. By substituting Eq. (21) into Eq. (9), the Lorentzian PSD can be rewritten into the following form, with two random variables (\(A_\Delta\) and \(\tau_0\)):

\[
S(f) = 4\beta(1 - \beta)(\Delta A)^2 \frac{\tau_0}{1 + (2\pi f \tau_0)^2}.
\]  
(24)

Assuming that there are \(N\) traps in total, the superposition of all the individual PSDs is expressed as

\[
S_N(f) = \sum_{i=1}^{N} 4\beta(1 - \beta)(\Delta A_i)^2 \frac{\tau_{0,i}}{1 + (2\pi f \tau_{0,i})^2},
\]  
(25)

where \(\Delta A_i\) is the RTN-induced fluctuation. Note that each trap has its own amplitude and time constants which is mentioned in Section III-B. Actually, \(N\) is also a random variable. Since the summation of multiple independent Poisson distributions is still a Poisson distribution, \(N\) also obeys a Poisson distribution. For a large \(N\), the summation in Eq. (25) can be converted into an integral:

\[
S_N(f) = 4\beta(1 - \beta) \int_0^\infty (\Delta A)^2 pdf(\Delta A) d\Delta A \int_{\tau_{0,\min}}^{\tau_{0,\max}} \frac{\tau_0}{1 + (2\pi f \tau_0)^2} pdf(\tau_0) d\tau_0,
\]  
(26)

where \(pdf(\Delta A)\) depends on the implementation. The integral of \(\int_0^\infty (\Delta A)^2 pdf(\Delta A) d\Delta A\) is always a constant (denoted as \(A\)) regardless of the detailed pdf(\(\Delta A\)). Consequently, Eq. (26) can be converted into a closed form

\[
S_N(f) = \frac{4\beta(1 - \beta)A}{\ln \frac{\tau_{0,\max}}{\tau_{0,\min}}} \int_{\tau_{0,\min}}^{\tau_{0,\max}} \frac{1}{1 + (2\pi f \tau_0)^2} d\tau_0
\]

\[
= \frac{2\beta(1 - \beta)A}{\pi f \ln \frac{\tau_{0,\max}}{\tau_{0,\min}}} \left[\arctan(2\pi f \tau_{0,\max}) - \arctan(2\pi f \tau_{0,\min})\right].
\]  
(27)

Eq. (27) can be approximated by applying the 1st-order Taylor expansion to \(\arctan\) according to the value of \(f\) [35]:

\[
S_N(f) \approx \begin{cases} 
\frac{\beta(1 - \beta)A}{\pi^2 f^2 \ln \frac{\tau_{0,\max}}{\tau_{0,\min}}} \left(\frac{1}{\tau_{0,\min}} - \frac{1}{\tau_{0,\max}}\right), & f \gg \frac{1}{2\pi \tau_{0,\min}} \\
\frac{\beta(1 - \beta)A}{f \ln \frac{\tau_{0,\max}}{\tau_{0,\min}}}, & \frac{1}{2\pi \tau_{0,\min}} < f < \frac{1}{2\pi \tau_{0,\max}} \\
\frac{4\beta(1 - \beta)A(\tau_{0,\max} - \tau_{0,\min})}{\ln \frac{\tau_{0,\max}}{\tau_{0,\min}}}, & f \ll \frac{1}{2\pi \tau_{0,\max}}.
\end{cases}
\]  
(28)

Considering that \(N\) is also a random variable, the superposed PSD has exactly the same form as Eq. (28) with the only difference on the amplitude \(A\), since Eq. (28) is independent with \(N\). For convenience, we still use Eq. (28) to express the superposed PSD. The superposed PSD shows three different shaped spectrums (i.e., white, \(\frac{1}{f}\), and \(\frac{1}{f^2}\)). Among them, the \(\frac{1}{f}\) spectrum occupies a wide frequency range if \(\tau_{0,\max} \gg \tau_{0,\min}\).

B. Randomness Modeling

1) Autocorrelation: The autocorrelation function can be calculated from the PSD based on the Wiener-Khinchin theorem [36], i.e.,

\[
C(t) = \int_0^\infty S_N(f) \cos(2\pi f t) df.
\]  
(29)

By applying the autocorrelation function of the band-limited \(\frac{1}{f}\) spectrum [37] and ignoring the \(\frac{1}{f^2}\) part which is very small, the autocorrelation function is approximated by

\[
C(t) \approx A\beta(1 - \beta) \left(\frac{2}{\pi} + \ln \frac{\tau_{0,\max}}{\tau_{0,\min}}\right), \quad \tau_{0,\min} \leq t \leq \tau_{0,\max}.
\]  
(30)

We also have

\[
C(0) = \int S_N(f) df = \beta(1 - \beta)A \left(\frac{4}{\pi \ln \frac{\tau_{0,\max}}{\tau_{0,\min}}} + 1\right).
\]  
(31)

Then the autocorrelation coefficient is expressed as

\[
\rho(t_s) = \frac{C(t_s)}{C(0)} \approx \frac{1}{2} + \frac{\ln \sqrt{\frac{\tau_{0,\max}}{\tau_{0,\min}}}}{\frac{1}{2} + \ln \frac{\tau_{0,\max}}{\tau_{0,\min}}}, \quad \tau_{0,\min} \leq t_s \leq \tau_{0,\max}.
\]  
(32)

When \(t_s\) is close to \(\tau_{0,\min}\), the autocorrelation is close to 1.0 (if \(\tau_{0,\max} \gg \tau_{0,\min}\). Although Eq. (32) is not applicable for \(t_s < \tau_{0,\min}\), it is no doubt that the autocorrelation is approximately 1.0 in this case, because the sampling is faster than the fastest trap and two successive samplings tend to get identical values. To make the autocorrelation small, \(t_s\) should be close to \(\tau_{0,\max}\), which means that the sampling frequency should match the slowest trap, leading to a very low bit rate. We will show that, after truncating a few high-order bits from the converted digital words, the autocorrelation of the remaining bits can be significantly eliminated and the bias is also close to zero, such that \(t_s\) is not required to be close to \(\tau_{0,\max}\).
Let \( U, D \) and \( T \) be the amplified fluctuation, the converted digital words, and the words after truncation, respectively. According to the central limit theorem, the overall RTN effect caused by numerous traps can be modeled by a Gaussian process. As a result, the fluctuation \( U \) follows a normal distribution. Let \( \phi_U(\mu_U, \sigma_U; x) \) be the PDF of \( U \), where \( \mu_U \) and \( \sigma_U \) are the mean value and the standard deviation, respectively. Amplification does not affect the autocorrelation, so the autocorrelation coefficient of \( U \) equals \( \rho(U) \) which is given by Eq. (32). \( U \) is converted to digital words via an analog-to-digital converter (ADC) which encodes \( n \) bits (i.e., the output range is from 0 to \( 2^n - 1 \)). The ratio of \( \frac{2^n}{n} \) can affect \( \mu_U \) and \( \sigma_U \). However, for a theoretical analysis, we assume that the full-scale range of the ADC matches \( \mu_U \pm 3\sigma_U \). Actually, this can be achieved by adjusting the amplifier. We ignore the fluctuation out of the \( \mu_U \pm 3\sigma_U \) range since its probability is negligible (i.e., 0.0027). The autocorrelation coefficient of \( D \) is expressed as \( \rho(D) \) to eliminate the autocorrelation of \( D, n-k \) high-order bits of \( D \) equals \( s \) and \( k \) low-order bits are kept. In what follows, we will derive the autocorrelation coefficient, bias, and bit rate of \( T \).

First, we have the following probability for \( D \):

\[
P(D = i) = P(iQ - 3\sigma_U \leq X \leq (i + 1)Q - 3\sigma_U) = \int_{iQ - 3\sigma_U}^{(i+1)Q - 3\sigma_U} \phi_U(\mu_U, \sigma_U; x) \, dx, 0 \leq i \leq 2^n - 1,
\]

(33)

where \( Q = \frac{6\sigma_U}{\pi} \) is the quantized level of the ADC. Since \( T = i \) if and only if the digital value expressed by the \( k \) low-order bits of \( D \) equals \( i \), the probability of observing \( T = i \) is expressed as

\[
P(T = i) = \sum_{z=0}^{2^n-k-1} P(D = i + z2^k), 0 \leq i \leq 2^k - 1.
\]

(34)

Since the integral of a normal distribution PDF has no analytical solution, Eq. (34) can be only estimated by numerical methods. We have found that if \( n-k \geq 2 \), all the \( P(T = i) \)'s are almost equal, i.e.,

\[
P(T = i) \approx \frac{1}{2^k}, 0 \leq i \leq 2^k - 1, k \leq n - 2.
\]

(35)

The maximum relative error of this approximation is less than 0.65%. An explanation of Eq. (35) is provided in Appendix B.

Considering two successive sampling time points \( s \) and \( s + t_s \), the joint probability of observing \( D_s = i \) and \( D_{s+t_s} = j \) is given by

\[
P(D_s = i \cap D_{s+t_s} = j) = \int_{iQ-3\sigma_U}^{(i+1)Q-3\sigma_U} \int_{jQ-3\sigma_U}^{(j+1)Q-3\sigma_U} \phi_2(\mu_U, \sigma_U, \rho_U(t_s); x, y) \, dx \, dy,
\]

(36)

where \( \phi_2(\mu_U, \sigma_U, \rho_U(t_s); x, y) \) is the joint PDF of a bivariate normal distribution with an autocorrelation coefficient \( \rho_U(t_s) \) which is given by Eq. (32). The joint probability of observing \( T_s = i \) and \( T_{s+t_s} = j \) is given by

\[
P(T_s = i \cap T_{s+t_s} = j) = \sum_{z_1=0}^{2^n-k-1} \sum_{z_2=0}^{2^n-k-1} P(D_s = i + z_12^k \cap D_{s+t_s} = j + z_22^k), 0 \leq i, j \leq 2^k - 1.
\]

(37)

The autocorrelation coefficient of \( T \) is expressed as

\[
\rho_T(t_s) = \frac{\sum_{i=0}^{2^k-1} \sum_{j=0}^{2^k-1} ij P(T_s = i \cap T_{s+t_s} = j) - \mu_T^2}{\sigma_T^2},
\]

(38)

where \( \mu_T \) and \( \sigma_T \) are the mean value and standard deviation of \( T \), which can be easily calculated based on Eq. (35).

Since Eq. (36) involves a double integral of the joint PDF of a bivariate normal distribution, Eq. (38) has no closed form. Eq. (38) can be estimated by numerical methods, and then \( k \) can be selected such that \( \rho_T(t_s) \) is small enough. We have derived a heuristic and effective method to directly calculate the optimal \( k \) such that \( \rho_T(t_s) \) is negligible. The derivation is complicated so we put it in Appendix C. The optimal \( k \) is given by the following closed form

\[
k = \frac{n - \log_2\left( \frac{6}{\pi\epsilon} \right) \left( 2\epsilon \sqrt{1 - \rho_T^2(t_s)} \right) \ln(2\epsilon \sqrt{1 - \rho_T^2(t_s)})}{6},
\]

(39)

where \( \epsilon \) is a near-zero threshold to control the accuracy. We use \( \epsilon = 10^{-6} \) in this paper. According to the explanation in Appendix C, if \( k \) is selected based on Eq. (39), \( \rho_T(t_s) \approx 0 \).

2) Bias: The bias of \( T \) is expressed as

\[
b_T = \frac{1}{k} \sum_{i=0}^{2^k-1} \text{ones}(i) P(T = i) - \frac{1}{2},
\]

(40)

where \( \text{ones}(i) \) is the number of ones in the binary representation of integer \( i \). Based on that the maximum relative error of the approximation of Eq. (35) is less than 0.65% (if \( k \leq n-2 \)), we have

\[
|b_T| < \frac{1}{k} \sum_{i=0}^{2^k-1} \left[ \text{ones}(i) \times \frac{0.0065}{2^k} \right] = 0.00325.
\]

(41)

Eq. (41) reveals that, although \( D \) is not uniformly distributed, the bias of \( T \) is close to zero after truncation.

3) Bit Rate: The output bit rate after truncation is simply expressed as \( k_f \). According to Eq. (32) and Eq. (39), increasing \( f_s \) will decrease \( k \). \( f_s \) can be increased by \( 10 \times, 100 \times \), etc, while \( k \) is decreased by only a few bits. Consequently, to maximize the bit rate, \( f_s \) should be as high as possible and close to \( \frac{1}{\tau_{\text{min}}} \). If the von Neumann corrector is applied instead of bit truncation, the bit rate is lower than \( \frac{1}{k} n f_s \). Calculation based on Eq. (39) reveals that, even if \( \rho_U(t_s) = 0.95 \), \( k < \frac{n}{4} \) holds only when \( n < 4 \), leading to an impractical result \( k < 1 \). This reveals that the bit truncation scheme has a higher bit rate than the von Neumann corrector in practice.
C. Numerical Results

The representative TRNG scheme shown in Fig. 10 with a $5 \times 10$ transistor array is simulated. The 22nm PTM [32] is used. The widths of all the transistors are 50nm. $V_{dd}$ of all the pMOSFETs and the nMOSFET are 0V and 0.8V, respectively. We use the following RTN parameters for pMOSFETs: $(N_i) = 2$, $(\Delta V_{th}) = 20$V, $\tau_{0,max} = 10$ms, and $\tau_{0,min} = 1$µs. The RTN profile is randomly generated according to these parameters. The resolution of the ADC is $n = 8$. Since the purpose of this test is to verify the proposed randomness model, we assume that the amplifier can be adjusted such that $\mu_U \pm 3\sigma_U$ matches the full-scale range of the ADC. As a result, the ratio of $\bar{V}$ which can affect $\mu_U$ and $\sigma_U$ will have little impact on the final output. We use $\bar{V}_T = \bar{V}_s$ in this experiment. $V_d$ of the nMOSFET is utilized as the randomness source. Fig. 11 shows an example (different runs give different examples) of the simulated distribution and the fitted normal distribution of $V_d$ of the nMOSFET, in which the simulated distribution is converted from an 100-bin histogram. The fluctuation caused by multiple traps-induced RTN shows a nearly perfect normal distribution. Fig. 12 shows the simulated and theoretical PSDs of $V_d$ of the nMOSFET. The theoretical PSD is calculated by Eq. (28), where $A \approx 1.202 \times 10^{-4}$V$^2$ is calculated from simulation. The simulated PSD is generally consistent with the theoretical PSD with a small difference. The difference mainly comes from the approximation of Eq. (28).

Fig. 13 shows the theoretical and simulated autocorrelation coefficients, under different sampling frequencies and $k$. The theoretical autocorrelation is predicted by Eq. (32). The simulated autocorrelation before truncation is consistent with the predictions. The bit truncation scheme significantly eliminates the autocorrelation. The optimal $k$ values estimated by Eq. (39) are marked as bold and red on the labels of the X-axis. When $f_s \leq \frac{1}{70 \text{min.}}$, the optimal $k$ ensures that the autocorrelation after truncation is less than 5%. However, when $f_s = 5$MHz which is larger than $\frac{1}{70 \text{min.}}$, the autocorrelation is larger than 5% even if only 1 bit is kept.

Table I shows the simulated ApEn, MC $\pi$ values, bias and bit rate. The ApEn results further reveal that the optimal $k$ values calculated by Eq. (39) are correct. For the MC $\pi$ value, a significant error is observed when $f_s = 5$MHz due to the high autocorrelation. The biases of these cases are all at the magnitude of $10^{-4}$, revealing that the bit truncation scheme can ensure a near-zero bias. According to these results, it can be concluded that when $f_s \leq \frac{1}{70 \text{min.}}$, we can select an optimal $k$ by Eq. (39) such that the autocorrelation is small enough to generate high-quality random numbers.

For the bit rate, clearly, among all of these cases, the maximum bit rate such that the randomness is guaranteed is achieved when $f_s = 1$MHz and $k = 5$, which gives a bit rate of 5 Mbit/s. In this case, if the von Neumann corrector is applied instead of bit truncation, the bit rate is 1.92 Mbit/s, which is much lower than that of bit truncation.

### Table I: Simulation results of the TRNG based on multiple traps-induced RTN.

<table>
<thead>
<tr>
<th>$f_s$ (MHz)</th>
<th>$k$</th>
<th>ApEn</th>
<th>MC $\pi$</th>
<th>Bias</th>
<th>Bit rate (Mbit/s)</th>
</tr>
</thead>
<tbody>
<tr>
<td>$50$</td>
<td>5</td>
<td>0.6926</td>
<td>3.1393</td>
<td>7.44E-04</td>
<td>0.25</td>
</tr>
<tr>
<td>$100$</td>
<td>5</td>
<td>0.6927</td>
<td>3.1376</td>
<td>8.47E-04</td>
<td>0.3</td>
</tr>
<tr>
<td>$200$</td>
<td>5</td>
<td>0.6929</td>
<td>3.1358</td>
<td>4.37E-04</td>
<td>0.5</td>
</tr>
<tr>
<td>$500$</td>
<td>5</td>
<td>0.6930</td>
<td>3.1344</td>
<td>3.96E-04</td>
<td>0.6</td>
</tr>
<tr>
<td>$1K$</td>
<td>5</td>
<td>0.6929</td>
<td>3.1321</td>
<td>4.73E-04</td>
<td>0.6</td>
</tr>
<tr>
<td>$1.5K$</td>
<td>5</td>
<td>0.6927</td>
<td>3.1337</td>
<td>4.22E-04</td>
<td>0.5</td>
</tr>
<tr>
<td>$2K$</td>
<td>5</td>
<td>0.6896</td>
<td>3.1299</td>
<td>5.07E-04</td>
<td>0.25</td>
</tr>
<tr>
<td>$2.5K$</td>
<td>5</td>
<td>0.6864</td>
<td>3.0508</td>
<td>4.40E-04</td>
<td>0.25</td>
</tr>
</tbody>
</table>

D. Summary

In this section, we have derived a theoretical randomness model for multiple-traps induced RTN. We have demonstrated an interesting conclusion. When generating random numbers from multiple traps-induced RTN, the sampling frequency can be close to the switching frequency of the fastest trap. The high autocorrelation can be almost completely eliminated by truncating a few high-order bits from the converted digital words. The bias after truncation is also close to zero. We have provided a closed form to decide the optimal truncation.

VI. Case Study: An Oscillator-Based TRNG

In this section, we study an RO-based TRNG scheme and present how to determine key parameters for this TRNG, based on the proposed randomness models.
The TRNG scheme which contains a 25-stage inverter-based RO is shown in Fig. 14. Each stage has a load capacitance. Due to the RTN effect, the RO period will be random. A $K$-bit counter is used to count the rising edges of a high-frequency clock. The output of the RO clocks $k$ ($k \leq K$) D-type flip-flops (DFF). The inputs of the $k$ DFFs are connected to the $k$ low-order bits of the counter, and the $K - k$ high-order bits of the counter are discarded. Clearly, the digital output of the counter is sampled at the end of each period of the RO output, so the RO period is converted to digital numbers. Due to the randomness in the RO period, the output of the $k$ DFFs is also random. Actually, the theory of this scheme is the same as the representative TRNG based on multiple traps-induced RTN. The RO period is the randomness source which is affected by multiple traps-induced RTN, and the counter and the DFFs can be regarded as an ADC. We will explain how to determine $k$ for this scheme in the next subsection.

### B. Numerical Results

We test the RO-based TRNG at two technology nodes using the 22nm and 32nm PTM [32]. Parameters used in this test are listed in Table II. The widths of nMOSFETs and pMOSFETs are 80nm and 40nm, respectively. The load capacitance of each stage of the RO is 5pF. The counter runs at 2GHz. The width of the counter $K$ is 20. Please note that $K$ is not equivalent to the resolution of the ADC $n$ shown in Fig. 10. We will show how to calculate the equivalent $n$ for this scheme in the following.

Fig. 15 shows the simulated distributions of the RO period and the fitted normal distributions, under different $\frac{\tau_c}{\tau_e}$. The RO period shows approximate normal distributions. The ratio of $\frac{\tau_c}{\tau_e}$ can affect the mean value and variance of the distribution. Since $\frac{\tau_c}{\tau_e}$ depends on many low-level factors and the manufacture process, it is difficult to know the exact $\frac{\tau_c}{\tau_e}$ by theoretical analysis. We consider three different values of $\frac{\tau_c}{\tau_e}$ in this test.

The standard deviations of the three cases at the 22nm node are 28ns, 50ns and 35ns, respectively. For the 32nm node, they are 18ns, 27ns and 16ns, respectively. Obviously, $\tau_c = \tau_e$ yields the maximum variance of the RO period.

![Fig. 14: RO-based TRNG based on multiple traps-induced RTN.](image)

**TABLE II: Parameters used in the RO-based TRNG.**

<table>
<thead>
<tr>
<th>Parameter</th>
<th>Value</th>
</tr>
</thead>
<tbody>
<tr>
<td>$V_{dd}$ (nMOSFET/pMOSFET)</td>
<td>0.8V/0.9V</td>
</tr>
<tr>
<td>$V_{th}$ (nMOSFET/pMOSFET)</td>
<td>0.503V/0.461V/0.494V</td>
</tr>
<tr>
<td>$t_{rise}$</td>
<td>10ns/1µs</td>
</tr>
<tr>
<td>$t_{fall}$</td>
<td>10ns/1µs</td>
</tr>
<tr>
<td>$(N_I)$ (nMOSFET/pMOSFET)</td>
<td>1.0/2.0</td>
</tr>
<tr>
<td>$(\Delta V_{th})$ [2], [28]</td>
<td>20mV</td>
</tr>
<tr>
<td>$V_{c}$</td>
<td>0.7/1.6</td>
</tr>
</tbody>
</table>

**TABLE III: Simulation results of the RO-based TRNG.**

<table>
<thead>
<tr>
<th>Parameter</th>
<th>22nm</th>
<th>32nm</th>
</tr>
</thead>
<tbody>
<tr>
<td>$\tau_c/\tau_e$</td>
<td>1.0 1.0</td>
<td>0.1 0.1</td>
</tr>
<tr>
<td>Autocorrelation</td>
<td>0.062% 0.015% -0.25% -0.055% -0.087% -0.23%</td>
<td>0.6929</td>
</tr>
<tr>
<td>Bias</td>
<td>9.6E-05 6.1E-05 7.6E-05 -5.6E-05 -7.7E-05 -9.1E-04</td>
<td>0.6929</td>
</tr>
<tr>
<td>ApEn</td>
<td>0.693 0.693 0.693 0.693 0.693 0.693</td>
<td>0.6929</td>
</tr>
<tr>
<td>MC $\pi$</td>
<td>3.1311 3.1411 3.1376 3.1400 3.1543 3.1512</td>
<td>3.1512</td>
</tr>
<tr>
<td>$k$</td>
<td>7 6 6 6 5 5</td>
<td></td>
</tr>
<tr>
<td>Bit rate (Mbit/s)</td>
<td>1.80 1.45 1.61 1.49 1.21 1.28</td>
<td></td>
</tr>
</tbody>
</table>

Now we show how to calculate the equivalent $n$ and $k$ in the RO-based TRNG, based on the proposed randomness model for multiple traps-induced RTN. The equivalent $n$ is determined such that $[0, 2^n - 1]$ can cover most (i.e., 99.73%) of the digital outputs before truncation. Take the case of $\frac{\tau_c}{\tau_e} = 0.1$ at the 22nm node as an example. When the counter runs at 2GHz, $n \approx \log_2(28 \times 10^{-9}) \times (2 \times 10^9) \times 6) = 8.4$. The RO period is equivalent to the sampling period. According to Eq. (32), the autocorrelation of the RO period is about 80%. According to Eq. (39), we get $k = 6$, which means that we need 6 DFFs to sample the output of the counter. $k$ values calculated by Eq. (39) are shown in Table III.

Table III shows the simulation results at the two technology nodes under different $\frac{\tau_c}{\tau_e}$ ratios. These results reveal that the generated random numbers of these cases are all of high quality and good randomness. The bit rate at the 32nm node is lower than at the 22nm node, mainly due to the smaller variance in the RO period at the 32nm node.

**C. Randomness Test**

The NIST test suite [23] is adopted to evaluate the randomness of the generated random numbers. For each case, we generate 500 random bit sequences by 500 independent simulations (each simulates a time interval of 50ms), and then feed them into the NIST test suite. Table IV lists the pass rates of the reported p-values. A p-value larger than 0.01 indicates that the test is passed. High pass rates are observed from Table IV, indicating good randomness of the generated random
TABLE IV: Pass rates (%) obtained from 500 runs of NIST.

<table>
<thead>
<tr>
<th>22nm</th>
<th>32nm</th>
</tr>
</thead>
<tbody>
<tr>
<td>( \tau_e / \tau_c )</td>
<td>1.0</td>
</tr>
<tr>
<td>ApproximateEntropy</td>
<td>98.4</td>
</tr>
<tr>
<td>BlockFrequency</td>
<td>98.8</td>
</tr>
<tr>
<td>CumulativeSums( a )</td>
<td>98.9</td>
</tr>
<tr>
<td>FFT</td>
<td>98.4</td>
</tr>
<tr>
<td>Frequency</td>
<td>98.6</td>
</tr>
<tr>
<td>LinearComplexity</td>
<td>98.4</td>
</tr>
<tr>
<td>LongestRun</td>
<td>99.2</td>
</tr>
<tr>
<td>NonOverlappingTemplate( a )</td>
<td>98.8</td>
</tr>
<tr>
<td>OverlappingTemplate</td>
<td>99.0</td>
</tr>
<tr>
<td>RandomExcursions</td>
<td>98.0</td>
</tr>
<tr>
<td>RandomExcursionsVariant( a )</td>
<td>98.6</td>
</tr>
<tr>
<td>Rank</td>
<td>99.6</td>
</tr>
<tr>
<td>Runs</td>
<td>99.0</td>
</tr>
<tr>
<td>Serial( a )</td>
<td>98.8</td>
</tr>
<tr>
<td>Universal</td>
<td>Insufficient length of bit sequence</td>
</tr>
</tbody>
</table>

\( a \) These tests report multiple p-values, so each reported pass rate is the mean value of all the p-values from the corresponding test.

![Histograms of p-values obtained from 500 runs of the ApproximateEntropy test in NIST](image1)

Fig. 16: Histograms of p-values obtained from 500 runs of the ApproximateEntropy test in NIST (the X-axis is the p-value and the Y-axis is the count).

numbers. The distribution of p-values can also be utilized to evaluate the randomness of random numbers [23]. In theory, the distribution should be uniform. Fig. 16 shows histograms of p-values obtained from the ApproximateEntropy test (the block length is 5) in the NIST test suite. The six subfigures all show approximate uniform distributions, indicating high randomness of the generated random numbers, as well.

VII. COMPARISON

In conventional noise-based TRNGs [16], [17], device noises are periodically sampled, amplified and compared with a reference voltage to generate random bits. Since the thermal noise and \( \frac{1}{2} \) noise are both tiny (typical magnitudes are from \( nV \) to \( \mu V \) [38], strong amplifiers are required. On the contrary, RTN offers significant fluctuations in advanced technologies. Measured data have shown that \( \Delta V_{th} \) caused by a single trap can be larger than 70mV at the 22nm node [1]. Such a large fluctuation can be easily converted to digital bits without suffering from variations, e.g., signal coupling problems.

Conventional jitter-based TRNGs [19] typically use a slow and jittery clock to sample a fast clock. Due to the jitter of the slow clock, the fast clock is sampled at random positions so that random bits are generated. To achieve high randomness, the jitter must be larger than the period of the fast clock. However, measured data have shown that the jitter-to-mean period is at the magnitude of only \( 10^{-3} \) at the 0.18\mu m node [18], which is very small. Actually, the variation of the RO period caused by RTN can also be regarded as a “jitter”. As shown in Fig. 15, the jitter-to-mean period caused by RTN is at the magnitude of \( 10^{-2} \), such a big jitter allows that multiple bits can be generated from one sampling, resulting in a higher bit rate.

In summary, compared with conventional noise- and jitter-based TRNGs, the advantages of RTN-based TRNGs mainly come from the large fluctuations in advanced technologies. In addition, several studies have shown the increasing RTN effect due to the shrinking of the feature size [15], [27], [39], so RTN is becoming more significant. In practice, the RO jitter should be caused by all possible randomness sources, including RTN, \( \frac{1}{2} \) noise, white noise, and etc. However, measured data have shown that at the 22nm node, RTN is the major noise source and much more important than \( \frac{1}{2} \) noise [2]. This conclusion reveals that, the RO jitter at advanced technology node is mainly caused by RTN.

VIII. CONCLUSIONS

In this paper, we have derived fundamental randomness models for RTN-based TRNGs. We have given theoretical models for the autocorrelation coefficient, bias, and bit rate of TRNGs based on both single trap- and multiple-traps-induced RTN. We have given theoretical methodologies to determine key parameters for designing RTN-based TRNGs, such as the sampling frequency and the number of truncated bits. Table V briefly summarizes the most important points of the proposed randomness models. The proposed models have been verified by numerical simulations. An RO-based TRNG at two technology nodes is studied based on the model of multiple traps-induced RTN. The proposed randomness models will be verified by fabricated chips in our future work.

APPENDIX A

DERIVATION OF EQUATION (15)

Let \( \{X_n\} \) be the sampled binary sequence, and \( \{Y_n\} \) be an intermediate sequence

\[
Y_n = \begin{cases} 
X_{2n}, & \text{if } X_{2n} \neq X_{2n+1} \\
2, & \text{if } X_{2n} = X_{2n+1}.
\end{cases}
\]
If all the twos in \{Y_n\} are discarded, we get a new sequence \{X_n\}, which is the output of the von Neumann corrector for \{X_n\}. When \( f_s < \frac{1}{\pi^2} \), the second-order autocorrelation coefficient between \( X_{n+1} \) and \( X_{n+2} \) is less than 5%, so \{X_n\} can be approximately regarded as a Markov chain, and thus, we have the following probabilities for \{Y_n\}:

\[
P(Y_{n+1} = 1 \cap Y_n = 1) \approx P_{10}P_{01}P_{10} \Delta y_{11}
\]

\[
P(Y_{n+1} = 2 \cap Y_n = 1) \approx P_{10}(P_{00}P_{01} + P_{01}P_{11}) \Delta y_{21}
\]

\[
P(Y_{n+1} = 1 \cap Y_n = 2) \approx P_{10}(P_{00}P_{01} + P_{01}P_{11}) \Delta y_{12}
\]

\[
P(Y_{n+1} = 2 \cap Y_n = 2) \approx \frac{P_{00}P_{00} + P_{01}P_{11}}{P_{00}P_{01} + P_{11}} \Delta y_{22}. \tag{43}
\]

Clearly, “11” in \{Z_n\} is generated from “11”, “121”, “1221”, · · · in \{Y_n\}. However, the probabilities of observing “11” in \{Z_n\} and observing “11”, “121”, “1221”, · · · in \{Y_n\} are different, because the lengths of \{Y_n\} and \{Z_n\} are different. The difference of the probabilities equals the ratio of their lengths, which can be obtained from Eq. (18). Consequently, we have

\[
P(Z_{n+1} = 1 \cap Z_n = 1) \approx \frac{f_s}{2R_{eN}} \left[ y_{11} + y_{12}y_{21} + y_{12}y_{22} + y_{12}y_{22} + y_{12}y_{22} + \cdots \right]
\]

\[
= \frac{f_s}{2R_{eN}} \left[ y_{11} + y_{12}y_{21} \left( 1 - y_{12} \right) \right]. \tag{44}
\]

According to Eq. (17), the probabilities of ones and zeros in \{Z_n\} are balanced, and thus, the mean value and the variance of \{Z_n\} are 0.5 and 0.25, respectively. The autocorrelation coefficient of \{Z_n\} is expressed as

\[
\rho_{eN}(t_s) = \frac{P(Z_{n+1} = 1 \cap Z_n = 1) - 0.5 \times 0.5}{0.25} \approx 2 \left[ 1 - \frac{1}{2} - P_0P_1 + 3P_0P_1 e^{-\frac{t_s}{\sigma^2}} - P_0P_1 e^{-\frac{3t_s}{\sigma^2}} ight] - 1. \tag{45}
\]

The terms with respect to \( e^{-\frac{t_s}{\sigma^2}} \) and \( e^{-\frac{3t_s}{\sigma^2}} \) can be ignored since they are quite small, and then we can simply get Eq. (15).

**APPENDIX B**

**EXPLANATION OF EQUATION (35)**

We first consider Eq. (34). The physical meaning of Eq. (34) is illustrated in Fig. 17. The two-dimensional region constructed by the PDF of the normal distribution \( \phi_{v}(\mu_U, \sigma_U; x) \) and the interval \( [\mu_U - 3\sigma_U, \mu_U + 3\sigma_U] \) on the X-axis is divided into \( 2^{n-k} \) groups. All the groups have the same length \( \frac{6s}{\pi^2} \) on the X-axis. Each group is further divided into \( 2^k \) grids. All the grids have the same length \( \frac{6s}{\pi^2} \) on the X-axis. Each grid has a coordinate from 0 to \( 2^k - 1 \). Let \( s(i) \) be the area of grid \( i \). It is clear that \( s(i) = P(D = i) \), so we have

\[
P(T = i) = \sum_{z=0}^{2^{n-k}-1} s(i + z2^k), 0 \leq i \leq 2^k - 1. \tag{46}
\]

Due to the symmetry of the PDF, we have

\[
P(T = i) = \sum_{z=0}^{2^{n-k-1}-1} \left[ s(i + z2^k) + s((z+1)2^k - 1 - i) \right], \tag{47}
\]

It is easy to check that

\[
P(T = i) = P(T = 2^{k-1} - 1 - i), 0 \leq i \leq 2^{k-1} - 1. \tag{48}
\]

So we only need to consider the difference between half of all the \( P(T = i) \)'s. The difference between \( P(T = i) \) and \( P(T = j) \) is expressed as

\[
P(T = i) - P(T = j) = \sum_{z=0}^{2^{n-k}-1} \left[ s(i + z2^k) - s(j + z2^k) + s((z+1)2^k - 1 - i) - s((z+1)2^k - 1 - j) \right], \tag{49}
\]

The four area terms on the right side of Eq. (49) are in the same group for the same \( z \). If \( n - k \) is big enough, the length of each group will be small enough, such that the PDF curve within each group can be approximated by a straight line with little error, as shown in Fig. 17. If the approximation has no error, the algebraic sum of the four area terms is exactly 0, and thus, \( P(T = i) = P(T = j) \) (for \( 0 \leq i, j \leq 2^k - 1 \)). Of course the PDF curve within each group is not an exact straight line, so we have \( P(T = i) \approx P(T = j) \) in practice. Since the integral of a normal distribution PDF has no closed form, numerical experiments have verified that when \( n - k = 2 \), the maximum relative error of the approximation shown in Eq. (35) is about 0.65%. According to the above explanation, increasing \( n - k \) (i.e., decreasing \( k \)) makes the approximation more accurate so the error is smaller.

**APPENDIX C**

**HEURISTIC DERIVATION OF EQUATION (39)**

We first consider the physical meaning of Eq. (37). Like the two-dimensional case shown in Appendix B, the three-dimensional region constructed by the PDF of the bivariate normal distribution \( \phi_{v}(\mu_U, \sigma_U, \mu_V, \sigma_V; x, y) \) and the square region \( [\mu_U - 3\sigma_U, \mu_U + 3\sigma_U] \times [\mu_U - 3\sigma_U, \mu_U + 3\sigma_U] \) on the X-Y plane is divided into \( 2^{n-k} \times 2^{n-k} \) groups. The bottom of each group is a square and the length of one side is \( \frac{6s}{\pi^2} \). Fig. 18 shows the four nearest groups to the central point \( (\mu_U, \mu_V) \). Each group is further divided into \( 2^k \times 2^k \) grids. The bottom of each grid is also a square and the length of one side is \( \frac{6s}{\pi^2} \). Each grid has a local coordinate \((i, j)\) \((0 \leq i, j \leq 2^k - 1)\) in its group. If the contour of the PDF...
The projection of the bivariate normal distribution \( \phi_2(\mu_U, \sigma_U, \rho_U(t_s); x, y) \) is projected onto the X-Y plane, as shown in Fig. 18. The equation of the projected ellipse is given by

\[
(x - \mu_U)^2 + (y - \mu_U)^2 = 2\rho_U(t_s)(x - \mu_U)(y - \mu_U)
\]

\[
= -2\sigma_U^2(1 - \rho_U^2(t_s)) \ln \left( \frac{2\pi\epsilon}{1 - \rho_U^2(t_s)} \right),
\]

where \( \epsilon \) is the relative height of the contour, \( \epsilon \) should be near zero to get accurate results. A bigger \( \rho_U(t_s) \) leads to a higher eccentricity of the ellipse, i.e., the ellipse looks narrower.

Clearly, \( P(T_s = i \cap T_{s+t_s} = j) \) (Eq. (37)) equals the total volume of all the grids whose local coordinates are \( (i, j) \) (0 \( \leq i, j \leq 2^k - 1 \)). We can ignore all the grids out of the ellipse, since the cumulative probability out of the ellipse is quite small if \( \epsilon \approx 0 \).

Like the two-dimensional case explained in Appendix B, the PDF surface within each grid can be approximated by a plane, such that all \( P(T_s = i \cap T_{s+t_s} = j) \)'s are almost equal with an exception when the ellipse cannot cover sufficient grids. It can be analyzed from Fig. 18 that, if the ellipse cannot fully cover the two shaded groups, the volumes of the shaded grids in the ellipse cannot be balanced even if the approximation by planes is accurate enough, leading to a big difference between \( P(T_s = i \cap T_{s+t_s} = j) \)'s. If this happens, the autocorrelation coefficient of \( T \) will be high according to Eq. (38). Consequently, the optimal \( k \) to ensure a low autocorrelation coefficient should be selected such that the ellipse can just fully cover the two shaded grids, i.e., the two points \( (\mu_U + \frac{2\sigma_U}{\sqrt{2\pi\epsilon}}, \mu_U - \frac{6\sigma_U}{\sqrt{2\pi\epsilon}}) \) and \( (\mu_U - \frac{2\sigma_U}{\sqrt{2\pi\epsilon}}, \mu_U + \frac{6\sigma_U}{\sqrt{2\pi\epsilon}}) \) are both on the curve of the ellipse. Substituting either point into Eq. (50) will yield Eq. (39).

References


Boxun Li (S’13) received the B.S. degree from the Department of Electronic Engineering, Tsinghua University, Beijing, China, in 2009, where he is currently pursuing the M.S. degree.

His current research interests include energy-efficient hardware computing system design and parallel computing based on GPUs.

Yu Wang (S’05-M’07-SM’14) received the B.S. and Ph.D. (with honour) degrees from the Department of Electronic Engineering, Tsinghua University, Beijing, China, in 2002 and 2007, respectively.

He is currently an Associate Professor with the Department of Electronic Engineering, Tsinghua University. His current research interests include parallel circuit analysis, application specific hardware computing (especially on the brain related problems), and power/reliability aware system design methodology. He has authored and co-authored over 130 papers in refereed journals and conferences. He received the IBM X10 Faculty Award in 2010, the Best Paper Award in ISVLSI 2012, the Best Poster Award in ISLPED in 2012, and 6 best paper nominations in ASPDAC, CODES+ISSS, and ISLPED.

Xin Li (S’01-M’06-SM’10) received the Ph.D. degree in Electrical and Computer Engineering from Carnegie Mellon University, Pittsburgh, PA in 2005, and the M.S. and B.S. degrees in Electronics Engineering from Fudan University, Shanghai, China in 2001 and 1998, respectively.

He is currently an Associate Professor in the Department of Electrical and Computer Engineering, Carnegie Mellon University, Pittsburgh, PA. His current research interests include integrated circuit and signal processing. He received the NSF Faculty Early Career Development Award (CAREER) in 2012, the IEEE Donald O. Pederson Best Paper Award in 2013, a Best Paper Award from Design Automation Conference (DAC) in 2010, two IEEE/ACM William J. McCalla ICCAD Best Paper Awards in 2004 and 2011, and a Best Paper Award from International Symposium on Integrated Circuits (ISIC) in 2014.

Yongpan Liu (M’07) received the B.S., M.S., and Ph.D. degrees from the Department of Electronic Engineering, Tsinghua University, Beijing, China, in 1999, 2002, and 2007, respectively.

He is an Associate Professor in the Department of Electronic Engineering, Tsinghua University. His research interests include low power design, emerging circuits and systems, and electronic design automation. He published over 60 papers and designed several sensor chips, including the first nonvolatile processor (THU1010N). He served as a technical program committee member in DAC, ASPDAC, ASSCC. He received the ISLPED in 2012, ISLPED in 2013 Design Contest Award, and several best paper nominations.

Huazhong Yang (M’97-SM’00) received the B.S. degree in microelectronics in 1989, M.S. and Ph.D. degree in electronic engineering in 1993 and 1998, respectively, all from Tsinghua University, Beijing, China.

In 1993, he joined the Department of Electronic Engineering, Tsinghua University, where he is a specially-appointed professor of the Cheung Kong Scholars Program. His current interest includes wireless sensor networks, data converters, parallel circuit simulation algorithms, nonvolatile processors, and energy-harvesting circuits. He has authored and co-authored over 300 technical papers and 70 granted patents.

Xiaoming Chen (S’12-M’15) received the B.S. and Ph.D. degrees from the Department of Electronic Engineering, Tsinghua University, Beijing, China, in 2009 and 2014, respectively.

Since 2014, he has been working as a postdoctoral researcher in Electrical and Computer Engineering, Carnegie Mellon University, Pittsburgh, PA, USA.

He is also a part-time researcher of the Department of Electronic Engineering, Tsinghua University. His current research interests include efficient hardware computing system design and parallel computing based on GPUs.

Lin Wang received the B.S. degree from School of Mathematical Sciences, Nankai University, Tianjin, China, in 2009, and the Ph.D. degree from Academy of Mathematics and Systems Sciences, Chinese Academy of Sciences, Beijing, China, in 2014.

She is a lecturer in School of statistics, Capital University of Economics and Business, Beijing, China. Her current research interests include probability and mathematical statistics, computational biology, and bioinformatics.