Linear Complexity Pruning Rule Enables Efficient Buffer Insertion Considering Process Variations

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ABSTRACT

Advanced process technologies call for a proactive consideration of process variation to ensure effective timing yield and keep the manufacturing cost down. Buffer insertion critical for almost any high performance IC designs nowadays, however, has not gained enough attention in attacking the correlated process variations because of the lack of an efficient *pruning rule* that defines the dominance relationship between two randomly but dependent solutions. The major contribution of this work is an efficient twoparameter (2P) based pruning rule for buffer insertion considering both inter-die and intra-die spatially correlated process variations. Under the normality assumptions on distribution, we theoretically prove that the 2P-based pruning rule has linear complexity (both time and space). Experiment results confirm the linear scalability of such an algorithm, and the capacity of the algorithm is also increased by more than thousands of times compared to the state-ofthe-art technique. We further apply the algorithm for timing optimization considering correlated process variation. We show that buffer insertion without considering spatial correlated variation reduces the timing yield by more than 50%, thus alarm the importance of developing efficient algorithms for IC designs to attack the process variation effects actively.

1. INTRODUCTION

Advanced process technologies impose significant challenges for modern IC designers as we move into the ultra-deep submicron era where manufactured circuits exhibit substantial performance variations. The proactive consideration of process variation during the design stage is critical to ensure effective timing yield and keep the manufacturing cost down.

Current studies on process variations have been mainly focused on the variability modeling and statistical timing analysis (STA) [1, 2, 3]. There are very few papers that actively consider the process variation effects for design optimization. Critical for almost any high performance IC designs nowadays, the buffer insertion problem has been studied extensively with different objectives in literature [4, 5]. However, none of them has considered the effect of process variations yet. Moreover, almost all of these algorithms to some degree follow the same dynamic programming paradigm in solving the buffer insertion problem. This is mainly due to an important result owning to [4] that the dynamic programming based buffer insertion problem can be solved optimally in polynomial complexity by properly defining the *dominance* relationship (or pruning rule) between solutions.

However, when process variation is considered, solving the same buffer insertion problem "optimally" becomes tricky. One of the difficulties lies in the fact that solutions are no longer constant values, but random variables. Moreover, these solutions are not independent but dependent with correlations. Different from the correlation discussed in the STA community, such correlation is not due to path reconvergence, but mainly due to the way we compute the solutions. In dynamic programming based buffer insertion, solutions are computed recursively from downstream nodes, rendering solutions from the same subtrees are correlated in nature. When intra-die spatial variation is considered, solutions from different subtrees will also exhibit certain correlations, which further complicates the problem. Another difficulty is how to define the *pruning rule* between different solutions in the presence of process variations. Because without a properly defined pruning rule, a straight-forward implementation of dynamic programming would make the algorithm complexity to increase exponentially.

To the best of our knowledge, there are three pieces of recent work in literature [6, 7, 8] that have attempted to solve the buffer insertion problem with consideration of different process variations. However, none of them has addressed the above two difficulties with definite answers. [6] studied the buffer insertion problem considering only the effect of wire length variation, albeit the fact that wire length variation is not a typical process variation. It was assumed that there was no correlation between different solutions. Moreover, three heuristic pruning rules were proposed with none of them can bound the complexity of the algorithm. The largest benchmark has 1260 sinks and no runtime was reported. [7] proposed to capture the correlation between solutions by using the joint probability density function (JPDF), and compute the JPDF recursively. A four-parameter based pruning rule was proposed to reduce the runtime complexity. However, the complexity and the effectiveness of the four-parameter based pruning rule is not clear, as only small benchmarks with no more than 9 sinks were reported. [8] studied a similar buffer insertion problem with a one-parameter based pruning rule, which is a simplified version of the fourparameter based pruning rule. Moreover, both [7] and [8] only considered the random device variations, but not the spatial correlated variation.

The contribution of this work is as follows. We propose an efficient two-parameter (2P) based pruning rule in the context of buffer insertion. Under the normality assumptions,

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we theoretically prove that the 2P-based pruning rule has linear complexity (both time and space), which enables efficient implementation of buffer insertion algorithm with consideration of both inter-die and intra-die spatial correlated variations. When solutions are not normally distributed, we give a theoretical necessary condition and experimentally verify that the two-parameter based pruning rule still has linear complexity in practice. Equipped with the efficient pruning rules, we employ a first-order process variation model to solve the buffer insertion problem considering both random device variation, inter-die variation and intradie spatial correlated variations. Experiment results show that the algorithm based upon the 2P pruning rule can increase the buffer insertion capacity to more than sixty thousands of sinks, which is more than $60 \times$ improvement over [6] and $1000 \times$ improvement over [7]. Moreover, our experiment results also show that buffer insertion without considering spatial correlated variation reduces the timing yield by more than 50%, thus alarm the importance of developing efficient algorithms for IC designs to attack the process variation effects.

The rest of this paper is organized as follows. In Section 2, we propose a two-parameter based pruning rule with consideration of process variation. In Section 3, we present a first-order process variation model to consider different types of correlated variations. In Section 4, we discuss the buffer insertion algorithm with correlated process variations. We report experiment results in Section 5, and conclude in Section 6. Proofs of Theorems will be included in a technical report.

2. VARIATION AWARE PRUNING

2.1 Review of Deterministic Pruning

For a given buffered routing tree, two figure-of-merits are associated with every legal buffer position t in the tree: i.e., the *input loading capacitance* (or *downstream loading capacitance*) L_t and the *required arrival time* T_t . The basic buffer insertion problem formulation is to find the locations of buffers in the given routing tree such that the required arrival time (RAT) at the root is maximized [4]. In the context of dynamic programming, we first traverse the routing tree in the reverse topological order and propagate solutions bottom up while book-keeping all intermediate solutions. At the end (root), we pick the optimal solution with the largest RAT. By backtracking the chosen optimal solution, we can determine the optimal solutions for each sub-tree recursively.

The major complexity of dynamic programming based buffer insertion lies in the merging of two sets of solutions obtained from two different sub-trees. In general, the total number of possible combinations for merging is $n \cdot m$, where n and m are the number of solutions from two sub-trees, respectively. If all combinations are kept at each merging node, the number of solutions will grow exponentially towards the root. To avoid this problem, [4] proposed to define the *dominance relationship* (or *pruning rule*) between two solutions such that solution (L_1, T_1) dominates solution (L_2, T_2) if condition $L_1 < L_2$ and $T_1 > T_2$ are satisfied. In another word, solution (L_2, T_2) is redundant and can be removed. [4] proved that there would be no more than n+mnumber of solutions after pruning instead of $n \cdot m$. Moreover, pruning on N number of sorted solutions can be done in O(N) time, not quadratic time $O(N^2)$ for pair-wise com-

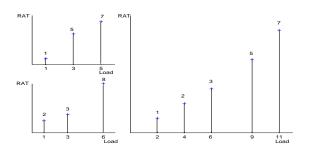


Figure 1: Linear merging O(n+m).

parison between solutions.

Even though pruning based upon the dominance relation can help reduce the total number of solution, in general we still have to pay the price of $O(n \cdot m)$ locally in order to obtain all possible combinations for merging. In deterministic buffer insertion, such a procedure can be further reduced to O(n + m) by using a merge sort like operation on the two sets of already sorted solutions. To see this, we give an example in Figure 1, where there are three solutions from each sub-tree, i.e., n = 3 and m = 3, and they are strictly sorted in terms of both L and T. By using a merge sort like operation on both solution sets while following the dominance rule, we can obtained all non-dominated solutions in linear time O(m+n). Moreover, the solutions after merging is also strictly sorted as shown in Figure 1.

Based upon the above two linear operations on pruning and merging, [4, 9] proved that by keeping only dominating solutions at every node, the dynamic programming based algorithm can solve the buffer insertion problem in $O(N^2)$ time without loosing optimality. When there are *B* types of buffers in the library, [10] proved that the deterministic buffer insertion problem can be solved optimally in $O(B \cdot N^2)$.

2.2 Review of Four-parameter Based Pruning

Even though the conventional deterministic buffer insertion problem can be solved optimally via dynamic programming [4], solving its variation-aware counterpart "optimally" is not that easy. The difficulty lies in the fact that L_t and T_t are no longer constant values, but two dependent random variables, as both are recursive functions of their downstream random variations. Moreover, different solutions, (L_1,T_1) and (L_2,T_2) , no matter whether or not they come from the same subtree, can be also dependent in the presence of spatial correlated variations.

By relating the dominance relation with designers' willingness of accepting uncertainty for a given design, [7] proposed the following *four-parameter* based (4P) dominance rule. Each parameter π_{α} gives a measure of designers preference for certainty in choosing the design parameter x in the presence of variations, such that the final design would have x less than π_{α} with $(100\alpha)\%$ certainty. Mathematically, this is given by

$$\alpha = \int_{-\infty}^{\pi_{\alpha}} f(x) dx, \qquad (1)$$

where f(x) is the probability density function (PDF) of x. Given four parameters, π_{α_l} and π_{α_u} for L_t , and π_{β_l} and π_{β_u} for T_t , such that $0 \le \alpha_l < \alpha_u \le 1$ and $0 \le \beta_l < \beta_u \le 1$, solution (L_1, T_1) is said to dominate solution (L_2, T_2) if the following conditions are satisfied:

$$\pi_{\alpha_u}^{(1)} < \pi_{\alpha_l}^{(2)} \tag{2}$$

$$\pi_{\beta_l}^{(1)} > \pi_{\beta_u}^{(2)}.$$
 (3)

Despite of its intuitive definition, such a dominance relation is hard to use in practice for large designs, because under such a dominance relation, we still have to pay the price of $O(n \cdot m)$ in computing all possible combinations for merging solutions, and $O(N^2)$ for pruning. Moreover, there is no guarantee that the number of solutions after pruning is linearly, which makes it less efficient for large designs as we will show in the experiment parts.

2.3 Two-parameter Based Pruning

We observe that for deterministic buffer insertion, the linear time operations for pruning and merging are made possible because the following two properties hold: (1) for any two given solutions, we can always order them by comparing their deterministic values, i.e., T_1 is either greater than T_2 or less than T_2 ; (2) there exists a *transitive ordering* between solutions, i.e., if $T_1 > T_2$ and $T_2 > T_3$, then $T_1 > T_3$ (similarly, if $L_1 < L_2$ and $L_2 < L_3$, then $L_1 < L_3$). If we ensure that the above two properties hold for solutions with variations, we can achieve similar linear time complexity for both pruning and merging¹. In the following, we propose a new variation aware pruning rule that would enable us to keep both merging and pruning operations in linear complexity even in the presence of process variations.

To eliminate uncertainty when comparing two solutions, we can extend the deterministic dominance relation between (L_1, T_1) and (L_2, T_2) by enforcing:

$$P(L_1 \le L_2) = 1,$$
 (4)

$$P(T_1 \ge T_2) = 1. \tag{5}$$

The physical interpretation of this extension is that solution (L_1, T_1) has 100% chances (almost always) to result in a larger required arrival time but with a less loading capacitance when compared to solution (L_2, T_2) . In another word, T_1 is said to be greater than T_2 if $P(T_1 \ge T_2) = 1$. We have the following Lemma².

LEMMA 1. Given T_1 , T_2 and T_3 as three **dependent** random variables with arbitrary distributions, if $P(T_1 > T_2) =$ 1, $P(T_2 > T_3) = 1$, then $P(T_1 > T_3) = 1$.

Proof: Let $X = T_1 - T_2$ and $Y = T_2 - T_3$ and the JPDF of X and Y be f(x, y). As $P(T_1 > T_2) = P(X > 0) = \int_0^{+\infty} x \int_{-\infty}^{+\infty} f(x, y) dy = 1$, we have $\int_{-\infty}^0 dx \int_{-\infty}^{+\infty} f(x, y) dy = 1$. $\int_0^{+\infty} dx \int_{-\infty}^{+\infty} f(x, y) dy = 1$ -1=0. Because $f(x, y) \ge 0$ for all x and y, we have f(x, y) = 0 for x < 0. Similarly, from $P(T_2 > T_3) = P(Y > 0) = 1$, we have f(x, y) = 0 for y < 0. Therefore, we have $\int_0^{+\infty} dy \int_0^{+\infty} f(x, y) dx = 1$.

 $\begin{array}{l} F(T_2 > T_3) = F(T > 0) = 1, \text{ we have } f(x,y) = 0 \text{ for } y < 0. \\ \text{Therefore, we have } \int_0^{+\infty} dy \int_0^{+\infty} f(x,y) dx = 1. \\ \text{Then } P(T_1 > T_3) = P(X + Y > 0) = \int_{-\infty}^{+\infty} dy \int_{-y}^{+\infty} f(x,y) dx = \\ \int_0^{+\infty} dy \int_{-y}^{+\infty} f(x,y) dx \ge \int_0^{+\infty} dy \int_0^{+\infty} f(x,y) dx = 1. \\ \text{As we know } P(T_1 > T_3) \le 1, \text{ we must have } P(T_1 > T_3) = 1. \\ \Box \end{array}$

Lemma 1 shows that comparison between T_1 and T_2 based upon $P(T_1 \ge T_2)=1$ enforces the transitive ordering property between solutions. However, for any given two solutions, it is not always possible to order them. Moreover, in practice, such a 100% chance requirement is too restrictive for pruning. Therefore, we relax such a requirement by adding two parameters such that solution (L_1, T_1) is said to dominate solution (L_2, T_2) if the following two conditions hold:

$$P(L_1 < L_2) > \overline{p_L},\tag{6}$$

$$P(T_1 > T_2) > \overline{p_T},\tag{7}$$

where $\overline{p_L}$ and $\overline{p_T}$ are two parameters between 0.5 and 1. In other words, the probability of L_1 being less than L_2 is greater than $\overline{p_L}$, while the probability of T_1 being greater than T_2 is less than $\overline{p_T}$. We call the pruning rule as defined by (6) and (7) as *two-parameter* (2P) based pruning rule in the following. We have the following lemma:

LEMMA 2. Given T_1 and T_2 as two **dependent** random variables with arbitrary distributions, we have either $P(T_1 > T_2) \ge 0.5$ or $P(T_1 < T_2) \ge 0.5$.

Proof: The proof follows directly from the fact that $P(T_1 > T_2) + P(T_1 < T_2) = 1$. \Box

Lemma 2 shows that comparison between T_1 and T_2 based upon $P(T_1 \ge T_2) > 0.5$ is always feasible for any two random solutions. Therefore, the remaining problem is whether or not such definition preserves the transitive ordering property between solutions.

Unfortunately, for arbitrary distributions, we can show that such definition in general do not preserve the transitive ordering property. However, when the distribution between random solutions is a joint normal distribution, we do have the transitive ordering property, which is stated by the following Lemma:

LEMMA 3. Given T_1 , T_2 and T_3 as three **dependent** random variables with joint normal distributions, if $P(T_1 > T_2) > 0.5$, $P(T_2 > T_3) > 0.5$, then $P(T_1 > T_3) > 0.5$

Proof: To see this, we assume both T_1 and T_2 (similarly L_1 and L_2) are normal, and we have the following closed form to evaluate the probability of $T_1 > T_2$: according to [12],

$$P(T_1 > T_2) = \Phi(\frac{\mu_{T_1} - \mu_{T_2}}{\sigma_{T_1, T_2}}), \tag{8}$$

where Φ is the cumulative density function (CDF) of standard normal distribution, μ_{T_1} , and μ_{T_2} are the mean values of T_1 and T_2 , respectively. σ_{T_1,T_2} can be computed by

$$\sigma_{T_1,T_2} = \left(\sigma_{T_1}^2 - 2 \cdot \rho_{T_1,T_2} \sigma_{T_1} \sigma_{T_2} + \sigma_{T_2}^2\right)^{1/2},\tag{9}$$

where $\sigma_{T_1}^2$ and $\sigma_{T_2}^2$ are variance of T_1 and T_2 , respectively; and ρ_{T_1,T_2} is the correlation coefficient of T_1 and T_2 .

Because any CDF function is a non-decreasing function, and for the standard normal distribution $\Phi(0) = 0.5$, then we have $\Phi(x) > 0.5$ for any x > 0. Therefore, to have $P(T_1 > T_2) > 0.5$ is equivalent to have $\frac{\mu_{T_1} - \mu_{T_2}}{\sigma_{T_1,T_2}} > 0$. Because σ_{T_1,T_2} is positive according to (9), hence we have $\mu_{T_1} > \mu_{T_2}$. Knowing $P(T_1 > T_2) > 0.5$ and $P(T_2 > T_3) >$ 0.5, we have $\mu_{T_1} > \mu_{T_2}$ and $\mu_{T_2} > \mu_{T_3}$. Therefore, we have $\mu_{T_1} > \mu_{T_3}$, which is equivalent to $P(T_1 > T_3) > 0.5$. \Box

According to the proof of the above lemma, it is easy to see that the following equivalent statements:

¹Note that the recently proposed predicative pruning [11] also exploits the same strictly ordering property of deterministic solutions.

²For simplicity, we only use T to illustrate the idea. It is understood that same results can be applied to L as well.

LEMMA 4. Given T_1 and T_2 as two **dependent** random variables with joint normal distribution, for $P(T_1 > T_2) > 0.5$ to hold, it is necessary and sufficient to have $\mu_{T_1} > \mu_{T_2}$ with μ_{T_1} and μ_{T_2} being the mean for T_1 and T_2 , respectively.

Based upon the above discussion, we have the following theorem regarding the complexity of dynamic programming based buffer insertion algorithm:

THEOREM 1. By utilizing the 2P-based dominance rule as defined in (6) and (7) with $\overline{p_L}=0.5$ and $\overline{p_T}=0.5$, the variation aware buffer insertion problem under the normality distribution assumption can be solved in $O(B * N^2)$, where B is the buffer types in the library, and N is the number of legal buffer positions.

Proof: According to Lemma (2) and (3), we can compare two random solutions and order them much the same way as in deterministic approach. Therefore, following similar arguments as in [4, 9] and [10], we conclude that our dynamic programming based variation aware buffer insertion problem can be solved in $O(B * N^2)$. \Box

Here are a few comments about the 2P-based prunning rule. The values of the two parameters, $\overline{p_L}$ and $\overline{p_T}$, in (6) and (7) determine how good the two desired properties (ordering property and the transitive ordering property between solutions) are approximately true. For arbitrary random distributions, on the one hand when $\overline{p_L} = 1$ and $\overline{p_T} = 1$, the transitive ordering property holds but not necessary the ordering property; on the other hand, when $\overline{p_L} = 0.5$ and $\overline{p_T} = 0.5$, the ordering property holds but not necessary the transitive ordering property according to Lemma 1. However, when the solutions follow certain distributions, the two proprerties can hold simultaneoully for some choice of $\overline{p_L}$ and $\overline{p_T}$. According to Lemma 2 and 3, we have proven that when the random solutions follow a joint normal distribution and for $\overline{p_L} = 0.5$ and $\overline{p_T} = 0.5$, both properties hold simultaneously, which enables us to prove Theorem 1.

Next we discuss the extension of 2P prunign rule for other choices of $\overline{p_L}$ and $\overline{p_T}$ and see how the two desired properties are affected³.

In fact, we have the following theorem which proves that for $\overline{p_T}$ (or $\overline{p_L}$) between 0.5 and 1, the transitive ordering property always hold under the normality assumption.

THEOREM 2. Given T_1 , T_2 and T_3 as three **dependent** random variables with joint normal distributions, if $P(T_1 > T_2) > \overline{P_T}$, $P(\underline{T_2} > T_3) > \overline{P_T}$, then $P(T_1 > T_3) > \overline{P_T}$ for any constant $\overline{P_T}$ between 0.5 and 1, i.e., $0.5 \leq \overline{P_T} \leq 1$.

Proof: Define $X = T_1 - T_2$ and $Y = T_2 - T_3$, then we have $X + Y = T_1 - T_3$. Therefore, $P(T_1 > T_2) = P(X > 0)$, $P(T_2 > T_3) = P(Y > 0)$, and $P(T_1 > T_3) = P(X + Y > 0)$.

Because T_1 , T_2 and T_3 are joint normal, then X and Y are also normal. Denote the PDF of X as $N(\mu_x, \sigma_x)$ and the PDF of Y as $N(\mu_y, \sigma_y)$, where μ_x and σ_x (similarly μ_y and σ_y) are the mean and standard deviation for X (similarly Y), respectively. Hence we can obtain the PDF of X + Y, which is also a normal distribution, as $N(\mu_x + \mu_y, \sqrt{\sigma_x^2 + \sigma_y^2 + 2\rho\sigma_x\sigma_y})$ with ρ being the correlation coefficient between X and Y.

We have

$$P(X>0) = P(\frac{X-\mu_x}{\sigma_x} > -\frac{\mu_x}{\sigma_x}) = 1 - \Phi(-\frac{\mu_x}{\sigma_x}), \quad (10)$$

where Φ is the CDF of the standard normal distribution (N(0, 1)). According to the property of standard normal distribution, we have $\Phi(-t) = 1 - \Phi(t)$, therefore we have

$$P(T_1 > T_2) = P(X > 0) = \Phi(\frac{\mu_x}{\sigma_x}).$$
 (11)

As we already know $P(T_1 > T_2) > \overline{P_T}$, we hence have $\Phi(\frac{\mu_x}{\sigma_x}) > \overline{P_T}$. Since any CDF function is also a non-decreasing function, we have

$$\frac{\mu_x}{\sigma_x} > \overline{t},\tag{12}$$

where $\Phi(\overline{t}) = \overline{P_T}$. Moreover, for $0.5 \leq \overline{P_T} \leq 1$, we have $\overline{t} > 0$.

Similary, we have

$$\frac{\mu_y}{\sigma_y} > \overline{t}.\tag{13}$$

From (12) and (13), we have

$$\mu_x + \mu_y > (\sigma_x + \sigma_y)\overline{t} \tag{14}$$

$$\frac{\mu_x + \mu_y}{\sqrt{\sigma_x^2 + \sigma_y^2 + 2\rho\sigma_x\sigma_y}} > \frac{(\sigma_x + \sigma_y)t}{\sqrt{\sigma_x^2 + \sigma_y^2 + 2\rho\sigma_x\sigma_y}}$$
(15)

Because $-1 \le \rho \le 1$, it is easy to show that

$$\frac{(\sigma_x + \sigma_y)}{\sqrt{\sigma_x^2 + \sigma_y 2 + 2\rho\sigma_x\sigma_y}} > 1.$$
(16)

As $\overline{t} > 0$, by multiplying both sides of (16) by $\overline{t} > 0$ and then combining it with (15), we have

$$\frac{\mu_x + \mu_y}{\sqrt{\sigma_x^2 + \sigma_y 2 + 2\rho\sigma_x\sigma_y}} > \frac{(\sigma_x + \sigma_y)\overline{t}}{\sqrt{\sigma_x^2 + \sigma_y 2 + 2\rho\sigma_x\sigma_y}} > \overline{t} \quad (17)$$

Therefore, by the fact that Φ is non-decreasing function as well as (17), we finally have

$$P(T_1 > T_3) = P(X + Y > 0)$$

= $\Phi(\frac{\mu_x + \mu_y}{\sqrt{\sigma_x^2 + \sigma_y^2 + 2\rho\sigma_x\sigma_y}})$
> $\Phi(\overline{t}) = \overline{P_T}.$ (18)

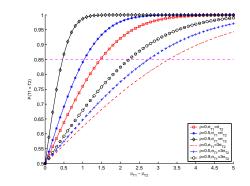


Figure 2: Probability of T_1 is greater than T_2 .

³Note, in the context of pruning, it is not interesting to take a value less than 0.5 for either $\overline{p_T}$ or $\overline{p_L}$.

In general, we could show theoreotically that the validity of the first property does not always hold for other choices of $\overline{p_L}$ and $\overline{p_T}$ except for 0.5 according to Lemma (2), however, we believe that in practice the validity of the first property still holds (or approximately holds) as we will show below. To see this, we plot the probability of T_1 greater than T_2 in Figure 2, according to (8), where the x-axis is the mean difference of T_1 and T_2 (μ_{T_1} - μ_{T_2}), and the y-axis is $P(T_1 > T_2)$. Three correlation coefficients, $\rho = 0, \rho = 0.5$ and $\rho = 0.9$, between T_1 and T_2 are plotted. The first three plots have $\sigma_{T_1} = \sigma_{T_2}$, and the rest three plots have $\sigma_{T_1} = 3\sigma_{T_2}$. According to Figure 2, we can see that when the difference between μ_{T_1} and μ_{T_2} becomes larger, the probability of T_1 being greater than T_2 is also increasingly larger. For a given required probability, say $\overline{p_L}=0.85$, it only requires μ_{T_1} being greater than μ_{T_2} by less than 4 time unit. In practice, for a general routing tree, such a small delay difference will likely exist among different solutions, either due to the difference in routing or due to the difference in buffering. Therefore, we conclude that pruning solution according to (6) and (7)will not loose much accuracy even when different $\overline{p_L}$ values are taken. Moreover, such an approximation becomes even better when two solutions have similar variance and have higher correlations, which are the cases for our buffer insertion problem because solutions from the same sub-tree or nearby sub-trees are highly correlated in nature.

Based upon Theorem 2 and the above discussion, we believe that in practice the 2P-based pruning rule could ensure the two desired properties, hence enabling us to carry out the variation aware buffer insertion algorithm much the same way as in the deterministic case. Therefore, similar runtime complexity conclusion can be drawn as shown in Theorem 1. Such a statement has also been verified experimentally in our experiment parts (Section 5).

3. PROCESS VARIATIONS MODELING

As has been shown in the STA community that blockbased STA analysis is suitable for increment computation of statistical timing. In the course of buffer insertion, such an increment computation feature is a must. Therefore, in the following, we discuss a first-order process variation model that incorporates both random device variation, inter-die variation and intra-die spatial correlated variations. Such a first-order modeling has been used for STA [3, 1], but none of them has explicitly considered all the above three type of variations. Moreover, their focus is mainly for timing analysis, while our focus in this paper is on design optimization.

3.1 Random Variation

We characterize a device (or buffer) in terms of its gate capacitance (C_b) , intrinsic delay (T_b) and output resistance (R_b) . Due to process variations, these values will no longer be fixed values. To simplify the model, we lump all variation effects into C_b and T_b while keeping R_b as a constant for a given device size. In general, devices characteristics are complicated (nonlinear) functions of the underlying physical parameters and sometimes are even hard to described in a closed form. Therefore, we resort to first-order approximation. The rational is that if the underlying parametric variations is small, any nonlinear relationship can be reasonably captured by a first-order approximation. Mathematically it can be described as:

$$C_b = C_{b0} + \sum \alpha_i \cdot X_i, \qquad (19)$$

$$T_b = T_{b0} + \sum \beta_i \cdot X_i, \qquad (20)$$

where C_{b0} and T_{b0} are nominal values of C_b and T_b , respectively; and X_i are the underlying parametric variations such as channel length, doping density, and gate oxide thickness. The coefficients α_i and β_i are sensitivity of C_b and T_b to the variation of X_i , respectively.

We run SPICE simulations to verify the accuracy of the above first-order modeling. For illustration purpose and also because of the lack of access to the real sources of foundry process variations, we only model the random L_{eff} variation using 65nm BSIM model in this section. We assume the variation of L_{eff} to be a symmetric normal distribution with the standard deviation as 10% of its mean value in our experiments. After extracting devices characteristics from SPICE simulation, we then use a least square curve-fitting technique to obtain (19) and (20).

Because of the nonlinear relationship between parametric variations (like channel length, doping density, gate oxide thickness, etc.) and the device characteristics, the latter's distributions are unlikely to be normal even if the underlying parametric variations are assumed to be normal. However, just as we have discussed above, if the underlying parametric variations is assumed to be small, the nonlinear relationship can be approximated by a first-order linear equation. Therefore, the device characteristics can also be approximated by a normal distribution. We validate this argument via Monte Carlo simulation. Figure 3 shows both the SPICE extracted PDFs of T_b and the normal distribution approximation from (20). It clearly shows that normal distribution is a reasonably good approximation of the real distribution because the two PDFs are very close to each other.

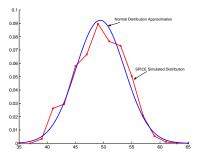


Figure 3: Normal distribution approximation of T_b .

3.2 Intra-die Variation

Current wafer manufacturing methods employ multiple identically oriented chips (usually four) within multi chip reticles during the stepping imaging process. Because of the optical lens distortion and wafer manufacturing processes, characteristics of devices on the same die are generally different and may take on the form of radial distortion [13], i.e., differences depending on distance from the center of the lens. Moreover, devices that are physically close to each other have higher correlation than devices that are far apart.

Therefore, to properly predict device characteristics, the following first-order model is proposed to capture the intradie spatial variations on device characteristics. We partition the die area into different regions, and associate each region with one independent random variables Y_i . When the variance of Y_i , σ_{Y_i} , are the same for all regions, we call such a model as *homogeneous* spatial variation model. When the variance of Y_i are different for different regions (and may even exhibit certain patterns), we call such a model as *heterogeneous* spatial variation model. Note that the spatial model proposed here is different from that of [14] in the sense that we use less number of independent variables but with more flexibility in capturing different spatial variations. We further note that associating each region with an independent random variable is also different from the principle component analysis (PCA) technique as employed in [1].

For a device located at a particular region R_t , we have

$$C_{b,t} = C_{b0} + \sum \alpha_i \cdot X_i + \sum_{i \in \mathcal{I}_t} \gamma_i \cdot Y_i, \qquad (21)$$

$$T_{b,t} = T_{b0} + \sum \beta_i \cdot X_i + \sum_{i \in \mathcal{I}_t} \theta_i \cdot Y_i.$$
 (22)

The index set \mathcal{I}_t defines the set of regions that spatial correlations matter for devices located at R_t , and the coefficient γ_i and θ_i further determine how strong the correlations are. In general, the larger the coefficients, the larger the correlation. Because the index set \mathcal{I}_t , and the coefficients γ_i and θ_i are region-dependent and different regions will have different \mathcal{I}_t , γ_i , and θ_i , by properly setting up these values, we can capture the spatial correlations between devices at different regions. For example, for two devices located at two nearby regions, they will share more number of common regions as decided by their common indexes in \mathcal{I}_{t1} and \mathcal{I}_{t2} , and thus would have larger corresponding correlations.

For example, Figure 4 shows a layout with regular defined regions. If we assume that each region is correlated with its closest neighboring regions (3×3 grids) with each region having an independent random variables Y_i , then the device's characteristics will be affected by Y_1 to Y_9 as given by (21) and (22). For two buffers that are physically closer to each other (e.g., buffer B1 and buffer B2), they share two number of correlated regions as defined by their respective correlation regions. On the other hand, for two buffers that are physically apart from each other (e.g., buffer B1 and buffer B5), they do not share any number of correlated regions, hence their spatial correlation is negligible.

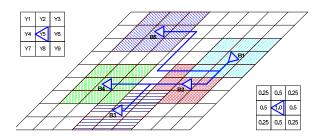


Figure 4: Modeling of the spatial correlations.

3.3 Inter-die Variation

As the inter-die variation affects all devices within the same die uniformly, we can model this variation by introducing another independent random variable, G and modify $\left(21\right)$ and $\left(22\right)$ as follows:

$$C_{b,t} = C_{b0} + \sum \alpha_i \cdot X_i + \sum_{i \in \mathcal{I}_t} \gamma_i \cdot Y_i + \xi \cdot G, \quad (23)$$

$$T_{b,t} = T_{b0} + \sum \beta_i \cdot X_i + \sum_{i \in \mathcal{I}_t} \theta_i \cdot Y_i + \eta \cdot G. \quad (24)$$

4. BUFFER INSERTION CONSIDERING PRO-CESS VARIATIONS

4.1 Key Operations for Buffer Insertion

In addition to the dominance rule between solutions, three other key operations are needed in order to solve the dynamic programming based buffer insertion problem. We review them briefly as follows. Denote c and r as interconnect's unit length capacitance and sheet resistance, respectively, we model each interconnect segment in the routing tree with length l as a π model. Under the Elmore delay model, the L_t and T_t can be computed recursively if we know the downstream node solutions.

If the solution at node t is obtained by adding a wire of length l at its direct downstream node n, then

$$L_t = L_n + c \cdot l \tag{25}$$

$$T_t = T_n - r \cdot l \cdot L_n - \frac{1}{2} \cdot r \cdot c \cdot l^2.$$
 (26)

If the solution at node t is obtained by adding a buffer at its direct downstream node n, then

$$L_t = C_b \tag{27}$$

$$T_t = T_n - T_b - R_b \cdot L_n. \tag{28}$$

If the solution at node t is obtained by merging two solutions from its two sub-trees rooted at nodes m and n, respectively, then

$$L_t = L_n + L_m \tag{29}$$

$$T_t = min(T_n, T_m). \tag{30}$$

4.2 Variation Aware Key Operations

In contrast to [7] where the process variation is modeled by the numerical JPDF, we employ the first order process variation model as discussed in Section 3 in this paper. We will report our comparison with [7] in Section 5. Below, we describe the changes that we need to make to the three key operations in solving the dynamic programming based buffer insertion problem.

When solutions at downstream nodes are modeled as random variables, the newly computed solutions (L_t, T_t) are also random variables. To make the recursive computation efficient, we keep the same first order form for L_t and T_t as their downstream nodes. In the following, we assume all downstream node solutions are known and can be represented by the following first order forms:

$$L_n = L_{n0} + \sum \alpha_{n,i} \cdot X_i \tag{31}$$

$$T_n = T_{n0} + \sum \beta_{n,i} \cdot X_i, \qquad (32)$$

where X_i can be any kind of defined variations. We then compute the solution at the current node t as follows. If the solution at node t is obtained by adding a wire of length l at its direct downstream node n, then

$$L_t = (L_{n0} + c \cdot l) + \sum \alpha_{n,i} \cdot X_i, \qquad (33)$$

$$T_t = (T_{n0} - r \cdot l \cdot L_{n0} - \frac{1}{2} \cdot r \cdot c \cdot l^2) + \sum (\beta_{n,i} - r \cdot l \cdot \alpha_{n,i}) \cdot X_i.$$
(34)

If the solution at node t is obtained by adding a buffer at its direct downstream node n, then

$$L_t = C_{b0} + \sum \alpha_{b,i} \cdot X_i, \qquad (35)$$
$$T_t = (T_{n0} - T_{b0} - R_b \cdot L_{n,0})$$

$$+\sum (\beta_{n,i} - \beta_{b,i} - R_b \cdot \alpha_{n,i}) \cdot X_i.$$
(36)

If the solution at node t is obtained by merging two solutions from its two sub-trees rooted at nodes m and n, respectively, then

$$L_t = (L_{n0} + L_{m0}) + \sum (\alpha_{n,i} + \alpha_{m,i}) \cdot X_i. \quad (37)$$

To express T_t after the *min* operation still to be a linear combination of underlying variations, we resort to the tightness probability idea from [3], and we have

$$T_{t} = t_{n,m} \cdot T_{n0} + (1 - t_{n,m})T_{m0} - \sigma_{n,m} \cdot \phi(\frac{\mu_{T_{m}} - \mu_{T_{n}}}{\sigma_{n,m}}) + \sum (t_{n,m} \cdot \beta_{n,i} + (1 - t_{n,m}) \cdot \beta_{m,i}) \cdot X_{i}, \quad (38)$$

where $t_{n,m}$ is the probability of T_n less than T_m , and can be computed by

$$t_{n,m} = \Phi(\frac{\mu_{T_m} - \mu_{T_n}}{\sigma_{n,m}}).$$
(39)

 ϕ and Φ are the PDF and CDF of the standard normal distribution, respectively; and $\sigma_{n,m}$ can be computed by

$$\sigma_{n,m} = (\sigma_n^2 - 2 \cdot \rho_{n,m} \cdot \sigma_n \cdot \sigma_m + \sigma_m^2)^{1/2}, \qquad (40)$$

where σ_n^2 and σ_m^2 are variance of T_n and T_m , respectively; and $\rho_{n,m}$ is the correlation coefficient of T_n and T_m . Knowing the first-order representation of T_n and T_m , we can compute σ_n^2 , σ_m^2 , and $\rho_{n,m}$ as follows.

$$\sigma_n^2 = \sum \beta_{n,i}^2 \sigma_{X_i}^2, \tag{41}$$

$$\sigma_m^2 = \sum \beta_{m,i}^2 \sigma_{X_i}^2, \tag{42}$$

$$\rho_{n,m} = \frac{\sum \beta_{n,i} \beta_{m,i} \sigma_{X_i}^2}{\sigma_n \sigma_m}, \qquad (43)$$

where $\sigma_{X_i}^2$ is the variance of X_i .

5. EXPERIMENT RESULTS

5.1 Experiment Setting

Two sets of benchmarks are obtained from the public domain for our experiments [11]. The characteristics of the benchmarks are shown in Table 1.

Because of the lack of access to the real wafer data, we derive the process variation data based upon the literature that addresses similar process variation issues but in the context of statistical timing analysis [14]. In our experiment, the 65nm BSIM technology is assumed. We budget the random device variation, inter-die variation, and intra-die variation

Bench	Sinks	Buffer Positions
p1	269	537
p2	603	1205
r1	267	533
r2	598	1195
r3	862	1723
r4	1903	3805
r5	3101	6201

Table 1: Characteristics of benchmarks.

all to be 5% of its nominal value, respectively. For the homogeneous spatial variation model, we uniformly distribute the budgeted 5% variation into different regions. For the heterogeneous spatial variation model, we distribute the budgeted 5% variation on the die from the South-West corner to the North-East corner in a linearly increasing fashion, i.e., devices located at the South-West corner has smaller spatial variation while devices located at the North-East corner has larger spatial variations. We divide the chip layout into different grids with the length of each grid as $500\mu m$. For devices located at a particular grid, their characteristics are affected by a set of nearby grids whose contribution weights form an isotropic stationary Gaussian process with the value tapers off at a distance about 2mm.

5.2 **Runtime Comparison**

We compare the runtime between our algorithm and [7] based upon RAT optimization (Section 5.3) as follows. Because there is no detailed runtime data from [7], we first compare the two algorithms runtime complexity in terms of the largest benchmark that each algorithm can handle. As reported in [7], the largest routing tree has only nine (9) sinks. In contrast, our algorithm can easily handle routing trees with more than three thousand sinks (> 3000), and it seems that nothing prevents our algorithm from handling even larger benchmarks⁴.

Bench	4P	2P	Speedup
p1	25.4	1.5	$17.3 \times$
p2	-	6.9	-
r1	-	25.0	-
r2	-	63.8	-
r3	-	131.5	-
r4	-	396.9	-
r5	-	922.8	-

Table 2: Runtime comparison in second.

Furthermore, we speculate that one of the main reasons in preventing [7] from trying larger benchmarks is that their 4P-based pruning rule is not very effective in both pruning and merging solutions. Another reason is due to the numerical complexity in computing JPDF, which requires the computation of high dimensional integral numerically. To verify this speculation, we have also reimplemented the algorithm of [7] based upon the same process variation models as used in this paper. In other words, we avoid [7]'s high complexity in computing the JPDF numerically, but rather to use the same first order process variation model to represent the JODI implicitly. Therefore, the only difference between our algorithm and that of [7] is the dominance rule. For the algorithm in [7], the 4P-based pruning rule as discussed in

 $^{^4\}mathrm{In}$ fact, the largest benchmark we have tested in house is an eight-level H-tree clock network with more than 64,000 sinks.

Section 2 is used, while our algorithm employs the newly proposed 2P-based pruning rule.

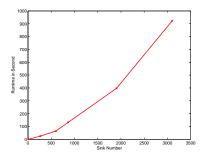


Figure 5: Runtime versus total number of sinks.

In Table 2, we report the runtime for both algorithms based upon the benchmarks we have tested. According to Table 2, we can see that the newly implemented algorithm of [7] now can handle much larger benchmarks than what was originally reported in [7], and the largest tested benchmark is p1 with 269 sinks. This improvement is mainly due to the avoidance of computing JPDF explicitly. However, we still fail to use the improved algorithm of [7] to run larger benchmarks. In fact, for the rest of tested benchmarks, it fails due to exceeding either memory capacity (2G) or tolerable time limit (4 hours in our setting). This observation is expected, because as we have explained in Section 2, the 4P-based pruning rule only imposes partially ordering between solutions, rendering the complexity of merging as $O(n \times m)$ and pruning as $O(N^2)$ for both memory and runtime. In contrast, by using the newly proposed 2P-based pruning rules, our algorithm can easily run through all benchmarks and for the largest benchmark r5, the runtime is less than 16 minutes. This significant runtime speedup is achieved because the 2P based pruning rule as discussed in Section 2 enforces a strict ordering between solutions, thus enables a linear complexity for both merging and pruning. This is further confirmed by Figure 5 which shows roughly the linear runtime scalability of our algorithms in terms of the number of sinks.

5.3 RAT Optimization

Enabled with the efficient implementation of buffer insertion considering both inter-die and intra-die variations, we can run our buffer insertion algorithm on the benchmarks for RAT optimization and study the effect of process variation on buffered interconnect design.

Before we report the experiment results, we first verify the accuracy of our model in predicting the RAT under process variations via the Monte Carlo simulation. Given a buffered routing tree with process variations, we run the algorithm to compute the PDF of RAT at the root. Such a procedure only requires the three key operations according to Section 4.2. After we obtain the RAT at the root represented in the form of (32), its PDF can be obtained by computing its mean and variance and it is approximated to be normally distributed. Figure 6 shows both the model predicted PDF and the Monte Carlo simulated PDF for the RAT at the root for one of the largest benchmarks (r5) in our experiments. We can clearly see that the first order process variation model is very accurate in predicating the PDF of

RAT, and therefore it can indeed be employed for RAT optimization.

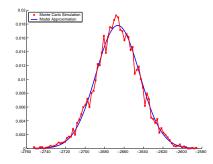


Figure 6: RAT at the root predicted by our model versus Monte Carlo simulation.

We compare three buffer insertion algorithms for RAT optimization in the following. The first one is the deterministic buffer insertion algorithm where all design variables are assumed to be nominal without variation, and we denote it as NOM. The second one is the variation-aware buffer insertion considering only random device variation and inter-die variation, but without considering the spatial variations. We denote such algorithm as D2D. The last one is the variation-aware buffer insertion algorithm considering all above variations, and it is denoted as WID. We run each algorithm on the given benchmarks and obtain the buffered routing tree. Two figure-of-merits are used to compare different algorithms. The first one is the 95% timing yield for RAT, which is defined as the 5%-tile RAT on the PDF such that the final RAT has the 95% chances of being larger than the 5%-tile RAT. The second one is the timing yield for a given required RAT at the root.

Table 3 compares the three algorithms using the heterogeneous spatial variation model as defined in Section 5.1. Column 2, 4 and 6 are the RAT at the 95% timing yield rate. The value in parenthesis are relative RAT degradation for NOM and D2D when compared to WID, respectively. According to Table 3, we can see that deterministic buffer insertion (NOM) without considering process variations always results in designs that require much higher RAT than WID and the increase can be as high as 23.1%. On average, NOM increases RAT by 9.7% compared to WID. From Table 3 we can also see that variation-aware design without considering spatial variation (D2D) also increases the RAT compared to WID, and the average increase is about 8.4%. It is interesting to note that compared to NOM, D2D achieves only about 2% improvement on average for RAT optimization. This confirms the common wisdom that global inter-die variation shifts the design space in the same direction while random device variation effect tends to diminish for a large design due to the variation canceling effects.

We also compare the timing yield for the three algorithms in Table 3. As we already observe that RAT from WID is about 10% better on average than both NOM and D2D, we set the mean RAT of WID with 10% reduction as the targeted RAT for all designs. We then compute the timing yield for all algorithms and report the results under Column 3, 5 and 7 in Table 3. According to Table 3, we see that NOMand D2D on average only achieve 42.2% and 46.7% timing yield and the yield loss is more than 50% when compared to WID.

Bench	NOM		D2D		WID	
	RAT $(\%)$	Yiled	RAT (%)	Yiled	RAT	Yiled
p1	-2673.5 (-2.4%)	99.6%	-2673.5 (-2.4%)	99.6%	-2611.7	100%
p2	-3791.3 (-7.7%)	99.9%	-3713.4 (-5.5%)	99.8%	-3519.3	100%
r1	-1240.7 (-15.9%)	0.5%	-1193.7 (-11.5%)	14.8%	-1070.3	100%
r2	-1808.1 (-23.1%)	0.1%	-1751.8 (-19.2%)	2.3%	-1469.2	100%
r3	-1658.8 (-9.3%)	48.5%	-1658.8 (-9.3%)	48.5%	-1518.0	100%
r4	-2475.7 (-10.7%)	27.6%	-2475.7 (-10.7%)	27.6%	-2236.0	100%
r5	-2934.9 (-8.6%)	83.5%	-2934.9 (-8.6%)	83.5%	-2703.3	100%
Avg	-9.7%	45.0%	-8.4%	47.0%	100%	100%

Table 3: RAT optimization under the heterogeneous spatial variation model.

Bench	NOM		D2D		WID	
	RAT (%)	Yiled	RAT (%)	Yiled	RAT	Yiled
p1	-2620.2 (-1.8%)	99.6%	-2620.2 (-1.8%)	99.6%	-2574.1	100%
p2	-3684.9 (-0.9%)	99.9%	-3685.9 (-0.9%)	99.8%	-3652.4	100%
r1	-1181.0 (-11.0%)	0.5%	-1141.0 (-7.3%)	14.8%	-1063.5	100%
r2	-1612.3 (-11.7%)	0.1%	-1575.9 (-9.2%)	2.3%	-1443.6	100%
r3	-1602.0 (-4.8%)	48.5%	-1602.0 (-4.8%)	48.5%	-1529.2	100%
r4	-2280.7 (-5.0%)	27.6%	-2280.7 (-5.0%)	27.6%	-2172.6	100%
r5	-2653.1 (-3.1%)	83.5%	-2653.1 (-3.1%)	83.5%	-2572.6	100%
Avg	-4.8%	45.0%	-4.0%	47.0%	100%	100%

Table 4: RAT optimization under the homogeneous spatial variation model.

Bench	NOM	D2D	WID
p1	$60 (1.09 \times)$	$60 (1.09 \times)$	55
p2	$156 (1.08 \times)$	$155 (1.08 \times)$	144
r1	$65 (1.44 \times)$	$61 (1.36 \times)$	45
r2	$135 (1.44 \times)$	$131 (1.39 \times)$	94
r3	$187 (1.42 \times)$	$187 (1.42 \times)$	132
r4	$375 (1.38 \times)$	$374~(1.38\times)$	272
r5	$608 (1.32 \times)$	$598 (1.30 \times)$	459
Avg	$1.15 \times$	$1.13 \times$	$1 \times$

Table 5: Number of buffers under different variation models.

We also run the same set of experiments but under the homogeneous spatial variation model as defined in Section 5.1 and report the experiment results in Table 4. Similar trends as in Table 3 can be observed. Moreover, by comparing Table 4 and Table 3, we observe that homogeneous spatial variation leads to in general a larger RAT for everv design. We further report the total number of buffers inserted for each design in Table 5. We find that for all designs, WID uses the least number of buffers compared to NOM and D2D. This results show that our WID algorithm can wisely insert buffers in the routing tree such that the final buffered solution achieves real RAT optimization in the presence of process variations.

Finally, we experimentally verify that different choices of $\overline{p_L}$ and $\overline{p_T}$ in (6) and (7) indeed have little impact on RAT optimization in practice. We have tried different combinations of $\overline{p_L}$ and $\overline{p_T}$ (from 0.5 to 0.95) for the same set of experiments as reported in Table 3 and 4. However, among all tested experiments, we see less than 0.1% difference in the final optimal RAT at the root. This observation is expected as we have discussed in Section 2.3.

CONCLUSION AND DISCUSSION 6.

An efficient two-parameter (2P) based pruning rule has been proposed for dynamic programming based buffer insertion. Under the normality assumptions on distribution, we have theoretically proved that the 2P-based pruning rule has linear complexity (both time and space), which enables

efficient implementation of the buffer insertion algorithm considering both inter-die and intra-die spatially correlated process variations. Experiment results have confirmed the linear scalability of our algorithm, whose capacity has been increased by more than thousands of times compared to [7]. We have applied the algorithm for timing optimization considering correlated process variations and concluded that process variation must be considered for real optimal designs, and buffer insertion without considering spatial correlated variation would reduce the timing yield by more than 50%, this calls for developing efficient algorithms for IC designs to attack the process variation effects actively.

In this work, we only considered timing optimization for signal nets. In the future, we intend to apply the same 2Pbased pruning rule and develop efficient algorithms for clock skew minimization.

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