

SOFT ERROR AWARE PHYSICAL SYNTHESIS

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To my wife, Reginara and my parents, Ailton and Telma for their infinite support

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CHAPTER 1

Introduction

Soft errors are events generated by the interaction of alpha particles, neutrons, protons, electrons, muons, or heavy ions with semiconductor regions in an Integrated Circuit (IC), potentially causing errors [1]. The growing concern of the semiconductor industry related to Soft Errors is attributed to the increased sensitivity of electronic devices occasioned by “Moore’s Law Scaling”. With the scaling of planar technologies from micrometer scale towards current nanometer scale transistors, parameters, such as transistors sizes, nodal capacitances, and interconnectivity resistances have decreased to the level where they directly impact the technology node sensitivity to radiation [2].

As charged particle transverses through a semiconductor region, it loses energy through coulombic interactions and generates electron-hole pairs. The effect of such charge generation inside a semiconductor region will be a function of both characteristics of the radiation event and physical details of the semiconductor region. The understanding of the detailed characteristics of the radiation event has an important role in the circuit soft error characterization. As it was shown by [3], ions with the same energy but different Linear Energy Transfer (LET) values have significantly different cross section. Moreover, ion-to-ion interaction, which can be predicted by Monte Carlo methods, can also impact the device error cross section.

In the context of transistors, the semiconductor technology process parameters and operational conditions have strong influence in how charge is collected by the individual transistors. As discussed by [1], lower operating bias, reduced switching energy, substrate doping and device cross-section have directly impacted on circuits soft error rate. The increase in transistor density on a die is also causing new

error modes and mechanisms. With the technology node scaling, more transistors can be fabricated in a given semiconductor region, allowing multiple transistors to collect charge from a single radiation strike, a phenomenon referred to as charge sharing [4]. This is an important issue that has been a focus of the research community for a long time, and is discussed in more details later.

In parallel to the increasing concern for Soft Errors in the reliability field, the Electronic Design Automation (EDA) research field has gained great attention from research community and industry. EDA has emerged as key science to allow the design of nanometer-size transistors and the integration of billions of electronic components in a single IC. Some of the reasons for the rapid growth and evolution of the EDA field were the increasing complexity of new technologies, extremely high density transistors in a single IC, and concerns over circuit designers productivity [5].

What is important to understand at this moment is that EDA has become a key part of the semiconductor industry to support the deployment of current and future semiconductor technologies. EDA allows the increasing complexity of new technologies to be hidden from semiconductor design engineers, with algorithms making key decisions that impact the electrical characteristics of a circuit. Without EDA tools, the latest generation of IC designs with billions of transistors [6] could not be possible. In current nanometer technology nodes it is impractical to consider a commercial design process without any kind of EDA tools to assist circuit designers.

EDA can be defined as a collection of methodologies, algorithms and metrics that automate the design, verification and test of electronic components [5]. This Ph.D. thesis is focused on design automation, more specifically the Physical Synthesis process. For the moment, it's enough to understand that Physical Synthesis is the process responsible for converting a list of logical gates and macro blocks into a physical representation of the IC that is placed, routed and electrically correct. The Physical Synthesis is one of the last steps of the Application Specific Integrated Circuit (ASIC) design flow process, before masks are built and sent to the IC foundries.

The case for the analysis of Soft Errors during the Physical Synthesis arises due to both cell placement and electrical correction techniques influencing the electrical characteristics of the circuit. Moreover,

because of the increase in the transistor density, charge sharing has become an issue and the relationship between adjacent electrical components impacts the overall circuit sensitivity to soft errors. The report [5] from the National Science Foundation (NSF), 2006, indicates soft error analysis as one of the key metrics to be integrated into the EDA flow.

Another important aspect is the commercial impact of the Physical Synthesis to the EDA industry. The Physical Synthesis by itself is responsible for most revenue by the EDA Industry, accounting for hundreds of millions of dollars [7]. On the other side, for companies, Non-Recurring Engineering (NRE) costs of Very Large Scale Integration (VLSI) designs are skyrocketing, and it has become critical for companies to avoid any kind of redesign effort. The report [5] from 2006, estimates an NRE of about \$30M/IC design for VLSI designs.

The integration of Soft Error awareness into the Physical Synthesis could potentially be used by companies to avoid redesign efforts in situations where Soft Error susceptibility needs to be reduced. Performing the soft error analysis automatically in the synthesis flow will be necessary due to the multitude of parameters affecting the physical synthesis and soft errors. It's the goal of this research to identify metrics for the soft error characterization and to investigate Physical Synthesis transformations that impact the soft error cross-section.

1.1 Related Work

The work published by [8, 9] implements a Placement aware of Soft Errors, using a Simulated Annealing algorithm, and later using Quadratic Modeling for the objective function optimization. These works use current waveforms to evaluate the circuit node soft-error sensitivity and propose the selective increase in the wire length, to increase the interconnection Resistance-Capacitance (RC), thus leading to the reduction of soft errors. Results from these papers indicate an average 27.12% [8] and 47.01% [9] soft error rate reduction according to their metric.

There are several differences between the work proposed by [8, 9] and the one proposed in this thesis. A key difference is that the selective wire length increase is not the objective constraint proposed in this work. Wire length increase is associated with both delay and power consumption increase, two objectives that are usually minimized by standard placement flows [7]. Another important difference is that the primary goal of this thesis is not to propose hardening techniques, but to present a detailed modeling approach to allow the integration of soft error analysis into the synthesis flow.

Other researchers have directed their effort to use commercial placement tools for Soft Error reduction [10, 11]. These projects usually refer to this methodology as a “Constrained Placement” where the jobs submitted to commercial placement tools receive additional constraints, usually used to group cells. Both the projects try to enhance the “Pulse Quenching” [12] effect by grouping cells into macros. Results from these papers indicate an average 35% [10] and 9-19% [11] soft error rate reduction. Table 1 is used to summarize the main aspects of the papers discussed above.

Table 1-1: Summary of main characteristics of related work

	Constraint	Objective	Soft Error Reduction
Placement Algorithm [8, 9]	Wire length	Increase interconnect RC	27.12%, 47.01%
Constrain Placement Tool [11, 10]	Cell grouping	Increase charge sharing	35%, 9-19%

According to the metrics used by the papers referenced before, the use of the RC product to filter small SETs is able to achieve a soft error reduction in the range of 27-47%, and for placements enhancing SET Pulse Quenching, the sensitivity reduction is in the range of 9-35%. All of these papers assume a very simple model for radiation events based on a simplistic double-exponential hit-current model. Though such a model is good enough for approximate analysis, pulse-quenching effects require a precise hit-current model in space and time to accurately model effects of an ion hit.

1.2 Contribution

The primary goal of this work is to build a methodology capable of properly modeling the Soft Error event, starting with the interaction of particles with the semiconductor region to the circuit-level response. Detailed characteristics of the radiation event are embedded in the methodology, to allow the proper characterization of the circuit cross-section and the evaluation of the influence of both electrical correction and circuit hardening techniques to the cross-section. Such accurate cross-section modeling efforts will result in better physical synthesis flow. Individual contributions are:

1. Develop a methodology for soft error analysis in a computationally intensive process.
 - a. Develop analytical models to estimate Collected Charge and Charge Sharing effects given a particle energy deposition profile and technology node characteristics.
 - b. Develop a methodology to estimate the SET Pulse Width.
 - c. Develop a methodology to estimate the soft-error cross-section.
2. Evaluate the impact of electrical correction techniques to the soft-error cross-section.
 - a. Gate Sizing, Gate Cloning and Buffering
3. Propose the optimization of Tap Cell placement to reduce the soft-error cross-section.

The results of this work will allow a better understanding by the EDA community of how the Physical Synthesis impacts the circuit-level soft-error cross-section and the limitations of various hardening techniques. By integrating this analysis into the Physical Synthesis flow, semiconductor companies will be able to reduce the soft -error cross-section of circuits without the need of a re-design.

1.3 Organization

This dissertation is organized as follow:

- Chapter 2: A background review of soft errors is discussed. Mechanisms involved in the charge collection by a diode and the parasitic bipolar amplification are discussed.
- Chapter 3: The model to estimate the collected charge and charge sharing by both p-type and n-type transistors is discussed, and a comparison with other models is presented.
- Chapter 4: The SET pulse width estimation methodology is discussed, presenting results from several different technology nodes. The methodology to estimate the circuit soft error cross-section is discussed.
- Chapter 5: The state of the art in of Physical Synthesis process is summarized, along with methods and transformations that lead to electrical changes that can impact circuit sensitivity to radiation. A discussion over the ideal synthesis steps to integrate the soft error analysis is presented.
- Chapter 6: The impact of the electrical correction techniques: Gate Sizing, Gate Cloning and Buffering are discussed and their impact to the circuit soft-error cross-section is evaluated. Circuit hardening based on Gate Sizing, Cloning and Tap Cell Placement techniques are evaluated.
- Chapter 7: Main thesis conclusions are discussed.

CHAPTER 2

Radiation Effects – Soft Errors and Mechanisms

Soft errors are transient events generated by the interaction of radiation with the semiconductor regions. These errors are considered transient because they will eventually disappear with circuit operation and do not directly cause permanent damage. There is an extensive range of error modes characterized as permanent, but these are not the focus of this work. For a broader perspective over permanent errors caused by radiation, please consult references [13, 14].

At this moment, it is important to identify the nomenclature used by the community to classify error modes associated with soft errors in digital circuits. The following are the error modes described by the research community that are important to the understanding of this work:

- Single-Event Transient (SET): The term SET is used to refer to a voltage transient that is generated in a logic gate, usually in a combinational logic block, and is able to propagate through the circuit [1].
- Single-Event Upset (SEU): An SEU refers to a soft error in a sequential element (memory). The SEU is able to flip the logic state stored in the feedback loop of the memory element [1].
- Multiple Bit Upset (MBU), Multiple Event Transients (MET): Both MBU and MET refers to multiple combinational and sequential elements been affected by a single event [1].

As this research focuses on the Physical Synthesis process, the primary error modes of interest are the SET/MET events. The SEU/MBU error modes are not ignored, but as memory elements are pre-designed and included in the IC as Standard Cells, the Physical Synthesis will have little impact on the soft error performance of memory cells. The next sections summarize the fundamentals of the soft error phenomena, followed by the procedure used to model the radiation event in this work, and the

methodologies used to characterize both the semiconductor device and circuit netlist response to Single Event Effects (SEE).

2.1 Soft Error Mechanism

The main physical mechanisms involved in the soft error event are shown in figure 2-1. In the figure an n-p diode representing a transistor drain, during a radiation event is shown. The particle can transfer energy by direct ionization (coulomb interaction) or indirect ionization (nuclear reaction). Direct ionization will transfer energy to the semiconductor material, exciting electrons from the valence band to the conduction band, thus generating electron hole pairs (e-h pairs). Indirect ionization will occur if a particle, such as a neutron or a proton, strikes an atom in the lattice structure, thus generating secondary ions. These ions create e-h pairs through direct ionization.

In the event shown in Figure 2.1, the generated e-h pairs create a distortion of the electric field located in the depletion region boundary (charged region). The distortion of the electric field will extend the collection area of the region, thus increasing the collected charge [15]. Carriers will move towards the depletion region by drift due to the presence of an electric field and diffusion due to the gradient of carriers in the region.

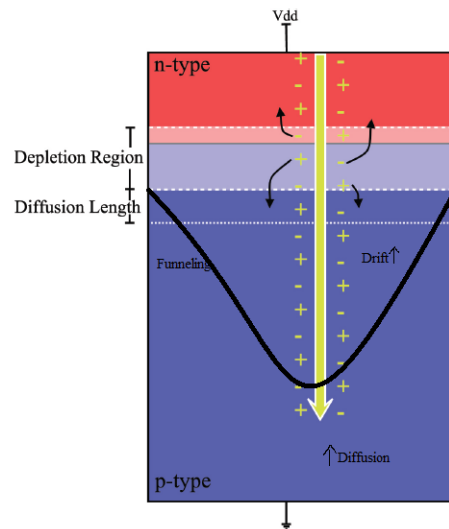


Figure 2-1 Electric field funneling in a junction during radiation event [16]

The charge collection profile can be represented with a double exponential current and is shown in figure 2-2. In the first few pico-seconds after the ion strike, most of the collected charge is due to the electric field. This is shown in the figure as the drift current component. After the first few pico-seconds, the charge collection due to diffusion dominates. Diffusion charge collection is a slow process and forms the long tail in the current waveform shown in the figure. This simple representation has been shown to be accurate for low LET particles, in the range of about 3-7 MeV*cm²/mg, but not accurate for high LET particles [16].

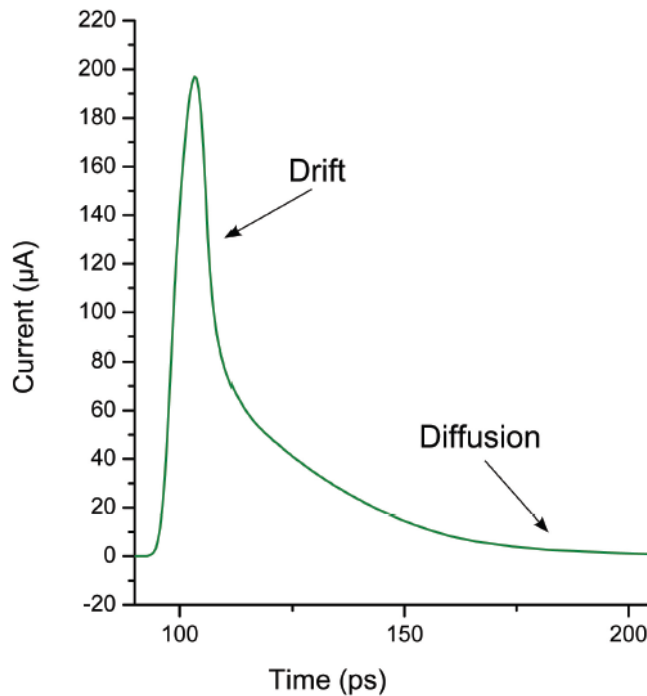


Figure 2-2 Current generated by the SEE [16]

The strength of electric field in the charge-cloud region, along with the depletion region width and device effective collection depth are important parameters for estimating the charge collection. Understanding both the radiation event and electronic device parameters are key to characterize the single event. The particle mass, energy and strike angle, along with the target electronic device characteristics like, device dimensions, doping, materials and bias are important to effectively characterize the device response.

The charge deposition process described in this section does trigger additional mechanisms that play an important role in the process. The scope of this problem from a simple event across a p-n junction

event has to be expanded to take into account effects of parasitic elements, multiple transistors, and standard cells. The well de-biasing and the parasitic bipolar amplification are going to be discussed next.

2.1.1 Well Debiasing and Parasitic Bipolar Amplification

The simple case of the n-p junction device is expanded to a model that includes additional devices with both p and n-type transistors forming a basic cell, as seen in figure 2-3. In this model, it is possible to identify additional mechanisms that affect the transistors and circuits during a radiation event. As discussed before, the ionization generates a cloud of e-h pairs that leads to a change in the electric field and a potential gradient in the well.

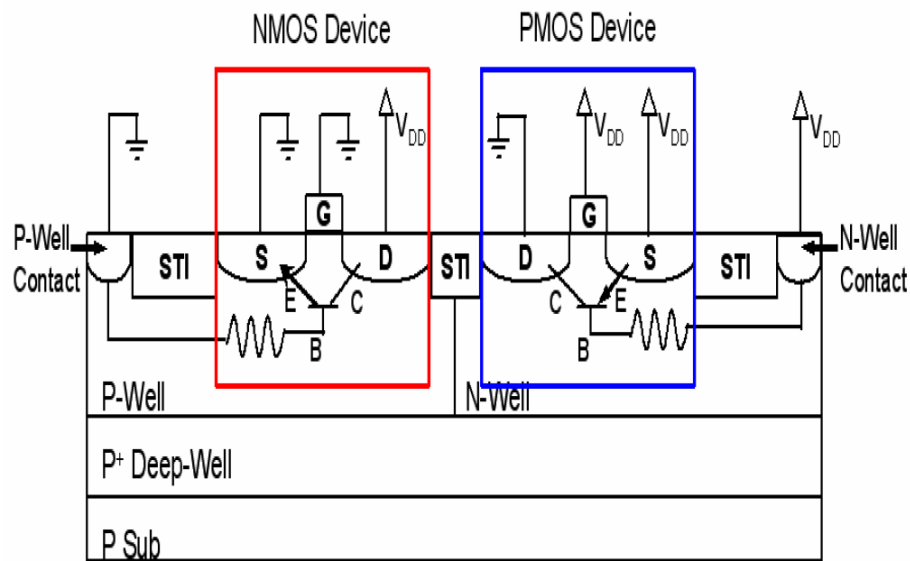


Figure 2-3 Parasitic bipolar amplification [17]

The localized potential drop in the substrate region of both transistors is caused by the deposited charge by an incident ion; this is referred to as well-debiasing or well-potential perturbation. Well-debiasing leads to the activation of the parasitic bipolar transistors, shown in Fig. 2-3, formed by the source/body/drain regions of transistors and shown in the figure by the Emitter(E)\Base(B)\Collector(C) terminals,. In the figure the two parasitic bipolar transistors, p+np+ and n+pn+, are identified. A potential drop in the channel (body) region, for instance for the PMOS transistor, can forward bias the source/body

junction leading to the activation of the parasitic bipolar transistor [17]. In this process carriers are going to be injected from the emitter to the body, and collected by the collector terminal, thus increasing the total amount of charge collected by the drain region during the radiation event. This charge, will then be represented at the circuit-level as a transient current.

It is important to identify the parameters affecting the perturbation in the well potential, as subsequently, they also control the bipolar transistor turn-on parameters. The P-Well and N-Well contacts are used to maintain a constant potential in the well. The vertical ($R_{vertical}$) and horizontal (R_{well}) well resistances are two layout dependent design parameters that can be used to control the size of the perturbation in the well potential, and the time required to revert back to the original well potential.

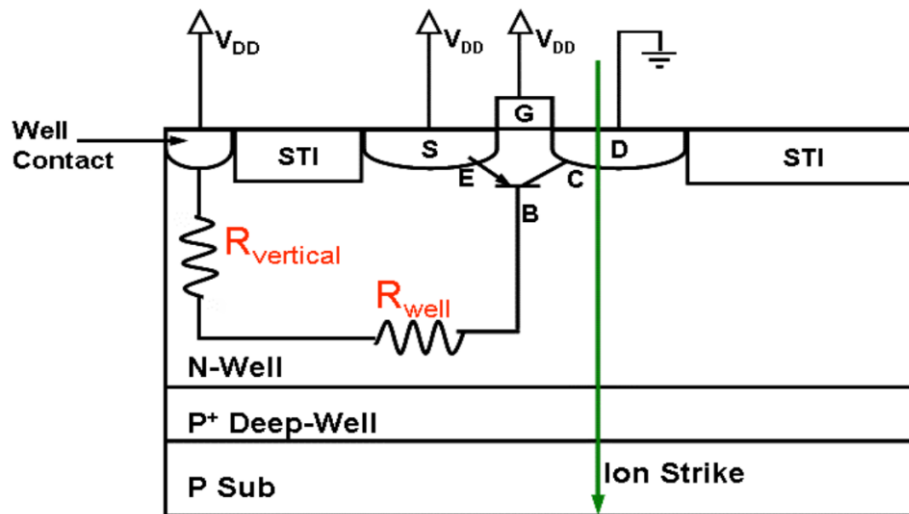


Figure 2-4 Resistances associated with the well potential [17].

The well resistance (R_{well}), can be controlled by the well contact distance. The closer the contact is to the base (channel region), the smaller will be the perturbation in the well potential, resulting in more charge required for the parasitic bipolar transistor activation. This will also result in smaller recovery time. The vertical resistance ($R_{vertical}$) is modulated by the contact area and well depth. Since well depth is not a parameter under designer's control, only the well-contact area is available to control $R_{vertical}$. The bigger the contact area, the less will be the resistance and the quicker will be the recovery of the well after the well de-bias event.

In dual well technology process with n-well, the p-channel transistor is considered to be less tolerant to the well-debias when compared with the n-channel transistor [18], due to both the fact that holes have lower mobility when compared to electrons and the confinement of charge in the well structure. For an additional insight related to well engineering techniques that can be used to reduce the well de-bias, please consult reference [19].

Now that a basic understanding of how the well de-bias can result in more charge collection by semiconductor regions, it's time to consider the Charge Sharing event. It turns out that in real circuits, several transistors are placed in close proximity of each other, sharing the same well. With scaling, the distances between these transistors are becoming increasingly smaller, and the transistor density is increasing. On the top of that, to increase the density of transistors, the number of well contacts in the IC is reduced. At advanced technology nodes, well and substrate contacts are designed as a separate cell and placed at a distance dictated by Design rules. Standard cells that do not contain well/substrate contacts are referred as tap less cells, and the cell containing the well/substrate contact is referred as Well Tap cell or Tap cell. These design practices leads to increased sensitivity of transistors to the well de-bias effects.

Due to the proximity of the transistors, charge sharing has become a reality. In nanometer technology nodes, a single radiation event can cause multiple logical nodes to collect charge, resulting in possible Multiple Event Transients (MET) propagating through the circuit. The charge-sharing event was investigated by [20, 21] in 130-nm and 90-nm technologies to show how it affects single-event sensitivity and hardening techniques for these and future technologies. Charge sharing may also be enhanced when a particle incident at a sharp angle traverses through multiple transistors.

With the understanding of important physical mechanisms required to model the transistor/circuit response during a radiation event, the next step will be to develop computationally efficient models for soft error analysis in this thesis. To achieve this goal, first the collected charge by a node needs to be estimated, followed by transformation of the collected charge into an electrical pulse at each of the node being affected. Simply put, we need to estimate the charge collection at all affected nodes and then convert the collected charge into multiple SET pulses in the circuit. The next section will discuss how the

radiation event is modeled, followed by the charge collection and supporting models to estimate standard cell response.

2.2 Radiation Event Modeling

Modeling the radiation event is a challenging process, due to its complexity. One of the main problems associated with the characterization of the radiation event is the fact that the radiation environment plays an important role in the characteristics of particles involved in the process. As it was discussed before, the radiation event can be classified between direct and indirect ionization, this in turns leads to the fact that different radiation environments can have different prevalent mechanisms to generate SEE. Discussing the different radiation environments and details over direct and indirect ionization are not the objective of this thesis, consult reference [3] for a better understanding of radiation environment characterization.

What is important to understand at this moment is that direct ionization can be directly described by the Linear Energy Transfer (LET), but not indirect ionization. To make the situation more complicated, in the terrestrial environment indirect ionization by neutrons play a major role in the characterization of the single-event effects. Modeling the radiation event in a way that allows both direct and indirect ionization to be integrated into the design/simulation flow is key for the soft error analysis.

A solution to bypass this problem is to use external tools and frameworks to characterize the event and model both direct and indirect ionization as a profile of energy deposition. This profile can be extracted using Monte Carlo simulation tools, like SRIM 2008 [22] and the Geant4 framework [23]. These simulators allow the description of the target material along with characteristics of the radiation particle in detail, and have been used before for this purpose. In figure 2-5, images from a Geant4 application developed to estimate the range and the energy deposition profile of positive and negative muons in a muon detector are shown.

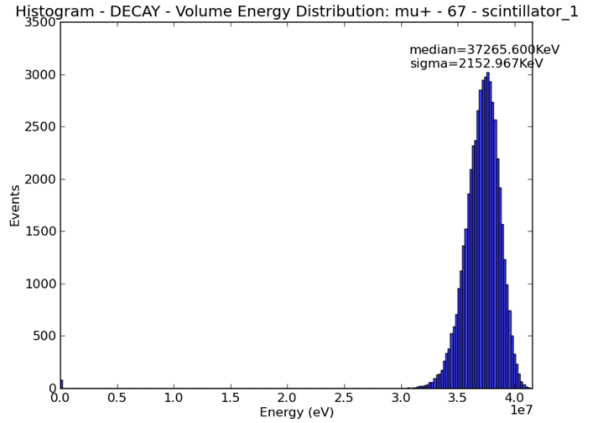
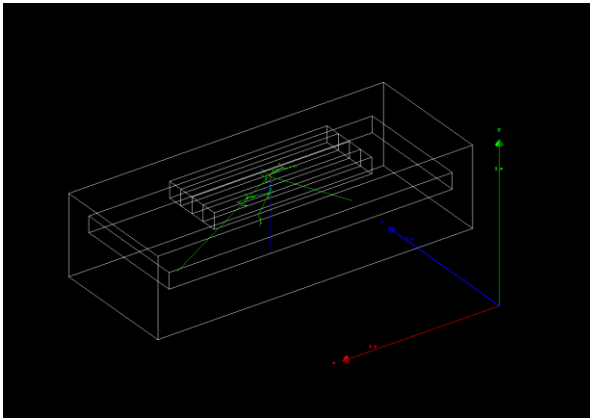


Figure 2-5 (a) Detector hit by a 67 MeV positive muon (μ^+). Blue and green lines indicate the path of positive and neutral charge respectively. (b) Energy distribution in the detector.

The radiation event is characterized in this work by a distribution (Uniform, Gaussian or Constant) of energy deposition profiles (dE/dx), and can be taken from simulations like the one shown in the last figure. Both Geant4 and SRIM2008 (which doesn't model subatomic particles) are able to estimate these energy deposition profiles given the particle type and energy, along with target structure dimension and material information. Be aware that there are significant differences between Geant4 and SRIM2008 tool, with the first being a more advanced and specialized software package. Due to its simplicity, the SRIM 2008 tool was used during most of this work, when necessary to estimate both Linear Energy Transfer (LET) using Eqn. 2.1 and Stopping Power (STP) with Eqn. 2.2 of heavy ions. These relationships are useful to estimate the amount of e-h pairs generated in a device by the radiation event.

$$LET = \frac{dE}{dx} < \text{MeVcm}^{-1} > \quad (2.1)$$

$$STP = \frac{dE/dx}{\rho} < \text{MeVcm}^2/\text{mg} > \quad (2.2)$$

With the stopping power and the target material known, one can also estimate the total number of ion pairs generated along the particle track. By knowing the amount of energy required to generate an e-h pair in Silicon material, one can estimate the amount of e-h generated in the track. For Si, the amount of energy required to generate an e-h pair is about 3.6 eV/e-h pair [24]. Thus the number of Ion Pairs (IP) generated by the ion can be estimated by using Eqn. 2.3.

$$I.P = \frac{dE/dx}{\omega} , \omega = eV/i.p , (2.3)$$

This relationship has been used by others [25, 26] to estimate the collected charge using the IRPP model. For this thesis, the LET of particles using SRIM2008 was calculated and used as input to the collected charge estimation method presented in the next chapter. Metal lines and the packaging material were not considered in these simulations, but could be easily included in the analysis.

Note that the data calculated from these tools are converted into lookup tables, with pre-computed values. These tables can be easily updated when necessary. For instance, the methodology presented in this work can receive data from the MRED tool [27] to improve accuracy as MRED is known for having great accuracy when predicting energy deposition profiles. With the assumption that the energy deposition profiles are known, the next step is to estimate the collected charge by the circuit transistors.

CHAPTER 3

Collected Charge Modeling

For advanced technologies, SEE analysis requires accurate estimation of collected charge at any given circuit node. Once collected charge is known, circuit simulators can use current sources or other models, such as the Bias-Dependent Model [28], to easily estimate voltage perturbations in order to predict circuit-level response to an incident ion. As the charge collected by a circuit node is a complex function of technology parameters (junction depth, doping densities, etc.) and ion characteristics (ion species, LET values, angle of incidence, etc.), this is one of the most challenging tasks for SEE analysis and predictions.

To enable the applicability of collected charge models to circuit analysis, the computational complexity of these models must be manageable without imposing significant loss in desired accuracy. This means low run time and memory space requirements. These requirements are especially important when models are to be integrated into an Electronic Design Automation (EDA) flow, a key achievement that could enable the use of soft error analysis in the standard ASIC design flow. This chapter discusses an accurate and computationally efficient model for estimating collected charge at a circuit node that could be used by the EDA community.

3.1 Background

Many analytical models, such as the RPP [29], IRPP [30], Messenger Double Exponential [31], electric field funnel model [32], and Ambipolar-Diffusion-with-Cutoff (ADC) model [33] have been proposed to estimate the collected charge, each with its own limitations and requirements. Some models, such as [30] and [32], are based on electric field funneling and are known to be limited to low LETs particles and short particle tracks [34]. Another major drawback for most of these models is their applicability to the latest generation of technology nodes.

For advanced technology nodes, the close proximity of transistors and small geometries result in increased charge-sharing between different semiconductor regions and parasitic-bipolar amplification within a transistor. Most of these models can not effectively model these two prominent mechanisms at advanced technology nodes. Unfortunately, the alternative to these shortcomings is to perform TCAD simulations for every transistor in the circuit, a task that requires significantly more computational and labor resources and time commitment to extract needed results. Since accurate estimation of collected charge is a key requirement in determining the response of circuits to an incident ion, faster techniques to estimate collected charge are of high interest to the radiation effects community.

In this chapter, the framework required to apply the ADC model to advanced technology nodes is discussed and analytical expressions based on the ADC model are developed for modeling charge-sharing. Comparison with published data for different technology nodes shows the efficacy of this approach in modeling collected charge and charge sharing, while also reducing the computational complexity of the problem. A comparison with both RPP and IRPP models shows a significant improvement in accuracy when using ADC.

3.2 ADC Model and Proposed Extension

The ADC was first proposed by Edmonds [33], for estimating charge collected by bulk diodes, and later extended to analyze multiple diodes [35]. The experimental validation of the ADC model and a methodology to estimate the required omega function used in the ADC model have also been reported [34]. Unlike other models for estimating collected charge, the ADC model does not attempt to model the single-event (SE) radiation event as a “funneling” of the junction electric field, but as a composition of regions within the semiconductor device with well-defined physical characteristics during the radiation event. Fig. 3-1 shows the three regions originally proposed in [33]: the Depletion Region (DR), the Ambipolar Region (AR), and the High-Resistance Region (HRR), characterized during the radiation event. Of these, the depletion region is the conventional depletion region associated with the p-n junction. The AR region is formed due to the high carrier density and weak electric field in the substrate region

satisfying the ambipolar diffusion equation [33]. The HRR region is formed at the bottom of the Quasi Neutral Region (QNR), where the concentration of carriers is much lower than the AR [33].

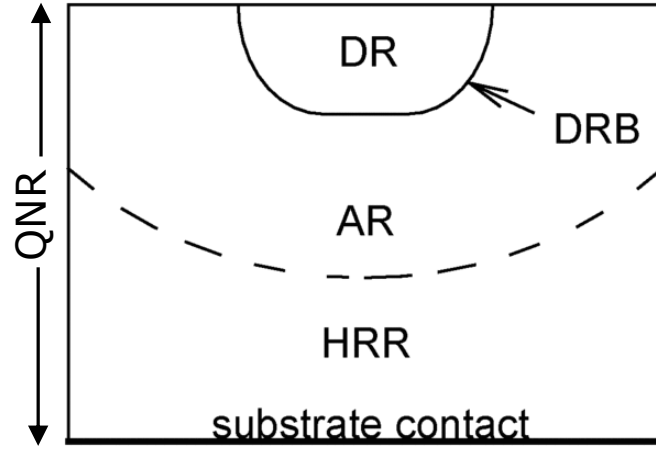


Figure 3-1 Different regions formed after the transit of an ion through a p-n junction [33].

These regions are modeled using drift and diffusion equations along with Poisson's equation to model the collected charge as shown below in Eqns. (3.1-3.3). Eqns. (3.1) and (3.2) yield the collected charge under high-level and low-level carrier injection conditions, respectively. For these equations, the limits of integration (Depletion Region (DR) and Quasi Neutral Region (QNR)) are shown in Fig. 3-1. Constants D_m and D_M are diffusion coefficient of minority and majority carriers, respectively, P_I is the density of e-h pairs along the track, and q is the elementary electronic charge and is equal to 1.602×10^{-19} C. Eqns. (2.1–2.3), are used to calculate the e-h pair density in the charge track, and further used to estimate total charge deposited.

$$Q_H = \left(1 + \frac{D_m}{D_M}\right) \int_{QNR} qP_I(\vec{x})\Omega(\vec{x})d^3x + \int_{DR} qP_I(\vec{x})d^3x \quad (3.1)$$

$$Q_L = \int_{QNR} qP_I(\vec{x})\Omega(\vec{x})d^3x + \int_{DR} qP_I(\vec{x})d^3x \quad (3.2)$$

$$\Omega(\vec{x}) = \left[\frac{Q_{TCAD}(\vec{x})}{Q_t}\right] \quad (3.3)$$

Software tools, such as MRED [27][9], SRIM2008 [22] and Geant4 [23] can be used to estimate P_I for a given particle and material as discussed in chapter 2. The omega function, $\Omega(\vec{x})$, is a position-dependent weighting factor, which determines how much charge would be collected in the terminals when carriers

are generated locally inside the device. Note that $\Omega(\vec{x})$ shown in (3.3) is derived using a point source under low level injection ($P_I(\vec{x})$ is a delta function) [34]. It satisfies the Laplace equation inside the QNR with appropriate boundary conditions [33] as:

$$\begin{aligned}\nabla^2\Omega(\vec{x}) &= 0 \text{ in QNR}, \quad \Omega(\vec{x}) = 1 \text{ on DRB}, \\ \Omega(\vec{x}) &= 0 \text{ on Substrate contact} \quad (3.4)\end{aligned}$$

All the device geometry information is contained in $\Omega(\vec{x})$ implicitly. It could only be solved analytically in certain simple cases, such as 1D simple diode and 3D isolated disk [33] or it could be probed through two-photon absorption laser experiment [34]. To evaluate the $\Omega(\vec{x})$ in device with complicated structure, TCAD simulations could also be used to construct the functional form by using the relationship shown in (3). After constructing $\Omega(\vec{x})$, it could predict the charge collection in low-level and high-level injection conditions based on Eqns. (3.1) and (3.2) assuming that $\Omega(\vec{x})$ is the same in both conditions. [34]. About 6 TCAD simulations were used to estimate the collected charge (QTCAD) and the total amount of charge generated by the event along the device depth (Q_t).

The above model was shown to work for a ($\sim 200 \mu\text{m} \times \sim 800 \mu\text{m}$) bulk silicon diode [34], but it is not directly applicable to transistors during a single-event. To apply the ADC model to the transistor case, an extension of the device model as used in the above discussion is necessary to take into account the physical differences as well as mechanisms occurring in a transistor, but not in a diode. These differences include the presence of a substrate contact, well/substrate boundaries, parasitic-bipolar amplification, charge sharing, and criteria to determine the low- and high-level injection conditions.

With the objective of characterizing the transistor device regions similar to those shown in Fig. 3-1 for the p-n diode, it was assumed that the bottom cutoff boundary for charge collection for p-type transistors is at the well boundary, where the well/substrate depletion region forms a secondary junction with the electric field limiting the collection of holes generated in the substrate. For the n-type transistors, this bottom cut-off boundary is given by the effective charge collection depth, assuming a dual-well process. Similar cutoff boundaries have been used before for TCAD SEE simulations [18].

For CMOS transistors, the presence of a parasitic bipolar transistor results in additional charge collection [20] and therefore must be taken into account in the model. As a result, the total collected charge for a PMOS transistor is the sum of the drift and diffusion currents inside the n-well and the parasitic-bipolar-amplification current. The parasitic bipolar amplification for NMOS transistors is assumed to be not significant for a dual-well process [20]. The parasitic-bipolar-amplification current is incorporated within the omega function for the proposed model as described in section V.

Lastly, to apply this model, it is important to identify device conditions and particle LET values required to create low and high-level injection. It was shown previously that charge deposition lower than 0.81 pC in the diode (n+p, p doped = $1 \times 10^{15} \text{ cm}^{-3}$) used by [34] creates a low-level injection condition. Although the specific value delineating high- and low-level injection conditions depends on the particular structure, this value provides a useful first-order estimate to be used in the proposed model for transistors. For particle LETs less than $30 \text{ MeV} \cdot \text{cm}^2/\text{mg}$, the total deposited charge in the n-well is less than 0.3 pC, which is much less than 0.81 pC. In addition, as the technology scales, the n-well doping increases, requiring higher LET particles to generate high-level injection conditions. Comparing the devices from [34] and the ones used in this work, the well/substrate doping is between 1 and 3 orders of magnitude higher, requiring even higher LETs to generate a high-level condition. Also, since the cross section saturates for most advanced technology nodes for particles with LET equal to or greater than $30 \text{ MeV} \cdot \text{cm}^2/\text{mg}$, the data showed in this work, and the proposed ADC extension, considers low-level injection conditions.

With the basic framework established to apply the ADC model to the transistor case, the model was extended to include charge-sharing effects. The objective is to estimate the charge collected by multiple transistors, given their distance to the ion hit location. Eqns. (3.5-3.7) show the equations used to accomplish this task.

$$Q_{qnr} = \int_{QNR} qP_1(\vec{x})(1 - \Omega(\vec{x}))d^3x \quad (3.5)$$

$$Q_{Ci} = Q_{qnr} * \Omega_d(\vec{d}) \quad (3.6)$$

$$\Omega_d(\vec{d}) = \left[\frac{Q_{TCAD}(\vec{d})}{Q_t} \right] \quad (3.7)$$

For the above equations, Q_{qnr} is the charge due to the diffusion current and the parasitic bipolar amplification and Q_{Ci} is the charge collected at node i . Q_{qnr} is modelled considering that the charge collected by a secondary transistor will be mainly influenced by the physical conditions of the AR and HRR regions, since in the DR region, $P_1(x)=0$. This assumption allows Q_{qnr} to be estimated using Eqn. (3.5). The charge collected at each node (3.6), is then a function of Q_{qnr} and the function $\Omega_d(d)$. The function $\Omega_d(d)$ is used to fit the parasitic-bipolar amplification and diffusion current, and thus is both technology and device dependent. To estimate Ω_d , a few TCAD simulations are required to estimate the collected charge vs. distance in an array of transistors. With this information, the ratio of collected charge and charge generated is calculated using Eqn. (3.7), and fitted to a curve to define the distance omega function, $\Omega_d(d)$. For the work presented in this paper, the function Ω_d was calculated using 8 TCAD simulations using the Sentaurus [36] simulator from the Synopsys tool set. The number of TCAD simulations required is small and the simulations need to be carried out only once for a given technology. A minimum of one simulation for each of the device regions shown in Fig. 3-1 is required for estimating the Omega function; more simulations will increase the accuracy of the model. More details related to the omega function estimation are provided in section V.

3.3. Devices, Technology Nodes and Radiation Event

To validate the accuracy of the model, 4 different technology nodes (130, 90, 65 and 40 nm) were analyzed, based on the availability of TCAD data for the comparison. Table I lists the technology parameters needed to characterize the devices to be evaluated by the proposed model. Except for the parameters listed in the table, no additional information regarding the technologies is required for the proposed model. These parameters are for dual-well, bulk CMOS planar technologies. Data for the 0.13 μm technology were acquired from [20], 90 nm from [21] and 40 nm from [18]. The data for 65 nm were estimated based on data from the 90 nm and 45 nm technologies.

The radiation event was modeled in this work as a single event, with charge deposition through coulombic interactions. Secondary interactions were not considered. To simplify the calculations, the density of electron-hole pairs was assumed to be constant along the particle track. For all calculations presented below, it was assumed that the particle was striking the center of the transistor drain (90° impact angle), crossing the device through the bottom cutoff boundary shown in Fig. 3-1.

The choice for a normal strike was due to the availability of TCAD data for the technologies and devices evaluated in this work. The capability of modeling angled strikes and changes in the strike location are two advantages that the ADC has when comparing with models that are based on the p-n junction electric field funneling mentioned in section I. These advantages are also shared by both RPP and IRPP models, since the injection depends mainly on the characteristics of the device region, particle track direction, length and energy. The work from [34] does show a comparison between the ADC applied to a diode with experimental Two-Photon Absorption (TPA) laser SEE measurement. For these experiments, particle tracks parallel to the junction were accurately modeled by the model.

Table 3-1- Technology Parameters

Tech\Parameters	130nm [20]	90nm [37]	65nm	40 nm [18]
Doping: Dif. N+	$2 \times 10^{20} \text{ cm}^{-3}$	$1 \times 10^{20} \text{ cm}^{-3}$	$1 \times 10^{20} \text{ cm}^{-3}$	$2 \times 10^{20} \text{ cm}^{-3}$
Doping: Dif. P+	$2 \times 10^{20} \text{ cm}^{-3}$	$1 \times 10^{20} \text{ cm}^{-3}$	$1 \times 10^{20} \text{ cm}^{-3}$	$2 \times 10^{20} \text{ cm}^{-3}$
Doping: Well N	$1 \times 10^{17} \text{ cm}^{-3}$	$1 \times 10^{17} \text{ cm}^{-3}$	$1 \times 10^{18} \text{ cm}^{-3}$	$1 \times 10^{18} \text{ cm}^{-3}$
Doping: Substrate P	$1 \times 10^{16} \text{ cm}^{-3}$	$1 \times 10^{16} \text{ cm}^{-3}$	$1 \times 10^{16} \text{ cm}^{-3}$	$1 \times 10^{16} \text{ cm}^{-3}$
Depletion Height	0.1 μm	60 nm	50 nm	30 nm
Well height	0.9 μm	0.75 μm	0.7 μm	0.6 μm
Effective Col. Depth	2.1 μm	1.56 μm	1.05 μm	1.03 μm

In the case of advanced technologies, shrinking dimensions require additional care for angled strikes. The single node collected charge can be estimated using the methodology presented in section III, but the pre-characterization step should be performed for different angled strikes for improved accuracy. For charge sharing estimation, the assumption of no charge generated in the DR of adjacent transistor may not hold for narrow angles, with $P_1(\vec{x}) \neq 0$ in the adjacent transistor depletion region. The charges in the DR of

nearby transistors have to be calculated with $\int_{DR} qP_I(\vec{x})d^3x$ for each DR in the particle track and added to the charge estimation (Q_{Ci}) calculated using Eqn. (3.6).

3.4. Pre-Characterization – Omega

As discussed in section 3.2 the solution of the Laplace's equation in the QNR can be approximated by Eqn. (3.3). With that in place there are two possible methods to pre-characterize the omega function for the ADC model: using experimental data or through TCAD simulations. Since experimental data for collected charge are very scarce and limited, published TCAD simulation results are used to estimate omega for the proposed model. The external data used in this work were carefully verified, taking into account details of the devices used in the simulations, along with the soft-error event characterization.

By using data from table I and 130 nm TCAD data from [20], the omega function was estimated for these technologies as a function of incident particle LET. It is important to notice that in this work only one pre-characterization was used for the 4 technology nodes described in 3.3, which will still provide accurate estimations. The reasoning for a single pre-characterization arises from the similar doping profiles used by these technology nodes. Fig. 3-2 shows the omega function generated for the 130-nm node for particle LET values of 1, 5 and 10 MeV*cm²/mg.

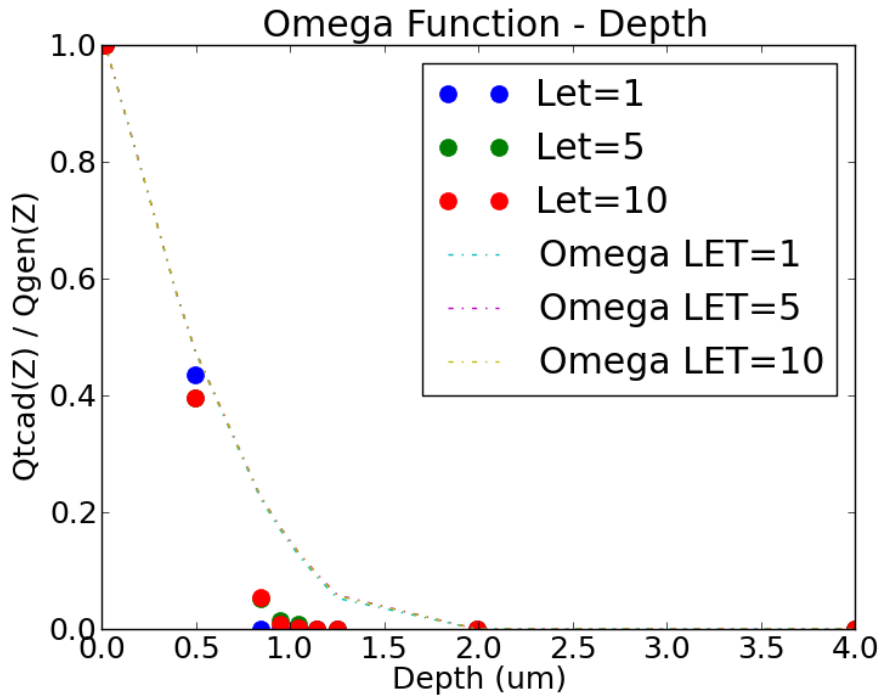


Figure 3-2 Omega Function as a function of depth. Filled circles represent TCAD results, dotted line is for the Omega function.

Using values of the omega function from Fig. 3-2 along with the density of e-h pairs corresponding to an LET of 10 MeV*cm²/mg, the collected charge as a function of depth is shown in Fig. 3-3. In the figure, within the depletion region the omega function is considered to be 1, corresponding to all charges generated in that volume being collected. From the DR boundary to the cut-off, which is considered to be the well boundary for PMOS transistors, the collected charge is calculated using $\Omega(x)$. The total collected charge is thus the integral of the curve shown in fig. 3-3.

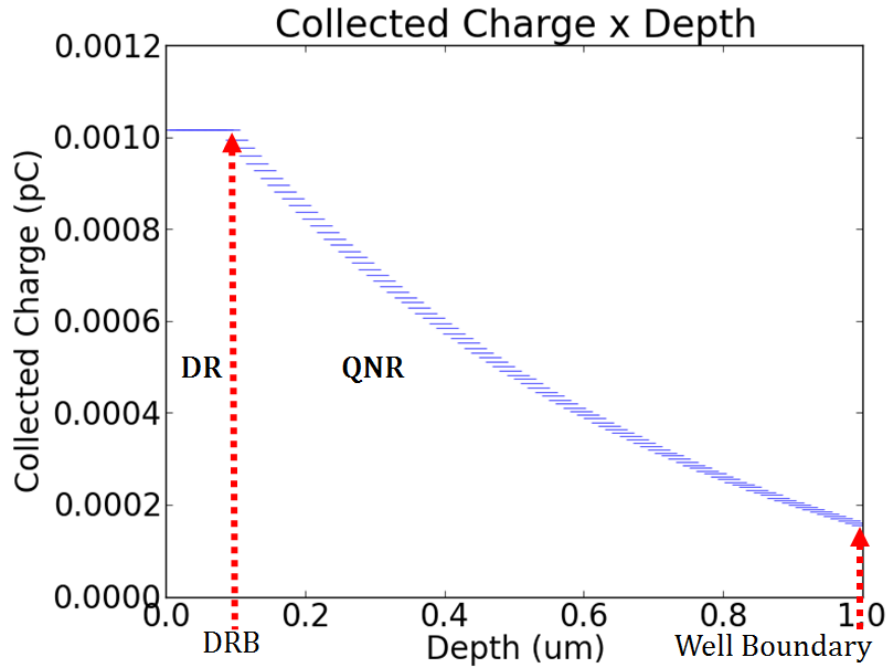


Figure 3-3- Collected Charge vs Depth for the 130 nm technology node parameters listed in Table I.

A similar procedure is used to estimate the omega distance function $\Omega_d(d)$ for the charge sharing Eqn. (9). Using TCAD data from the 130-nm node, the omega distance function was estimated by measuring the collected charge on transistors up to 2 μm from the hit location for different particle LET values. Eqn. (9) is then used to generate the data points for calibration. The secondary devices to collect charge are modeled in TCAD with the biasing setting the transistors in the “off-state”.

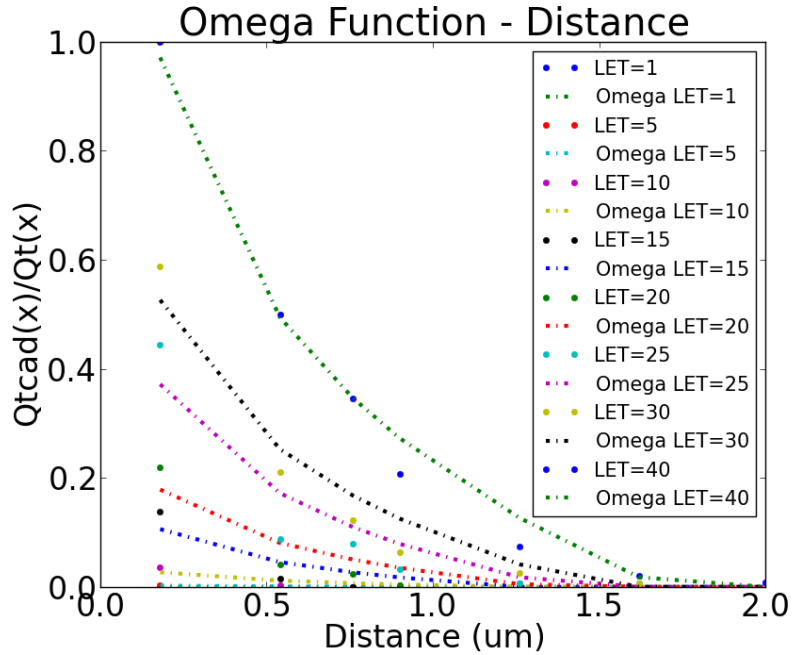


Figure 3-4: Omega function as a function of distance. Filled circles represent TCAD data, dotted line is for the Omega

By measuring the $\Omega_d(d)$ for multiple LETs the accuracy of the model can be improved by selecting the closest data set according to the LET used as input in the model.

3.5 Single Node Collected Charge

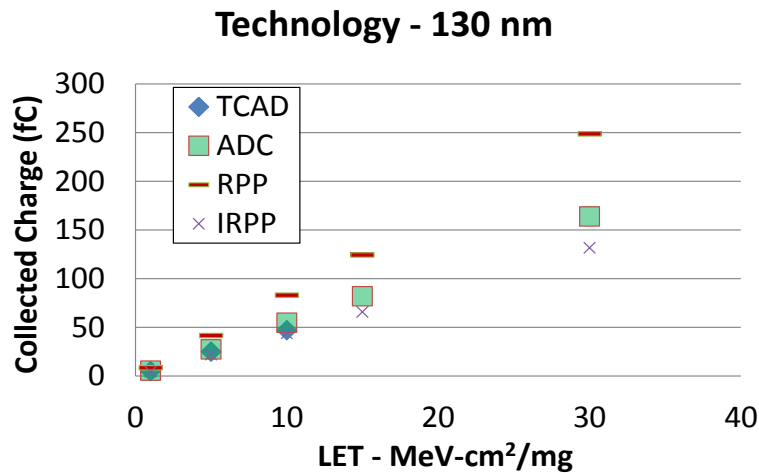
To validate the applicability of the proposed additions to the ADC model to estimate single node collected charge, the model is compared with TCAD simulation data and both RPP and IRPP estimations. The RPP and IRPP models can both be analytically represented with the nested sensitive volume Eqn. (3.8) [38], where α is a weighting coefficient and E the deposited energy. The first term of the equation is the energy to charge conversion factor assuming that 3.6 eV are required to produce an electron-hole pair [24].

Notice in this thesis the IRPP model is calculated considering multiple sensitive volumes. The IRPP model therefore requires the definition of sensitive volumes (SV) and collection coefficients (α_i). For this work sensitive volumes were defined based on the guidelines from [39], and the collection coefficients were estimated using 130nm technology data. For these devices a set of four SV, named as SV1, SV2,

SV3 and SV4 with collection coefficients 1, 0.62, 0.3 and 0.06, respectively, were used. The height of each SV has to be calculated for each technology, with the SV1 height being the DR height and SV2, SV3 and SV4 heights being approximately 30, 30, and 40%, respectively, of the distance from the DR to the well boundary for the p-type transistors (and effective collection depth for the n-type transistors). The RPP model can be seen as a single SV, with a constant collection coefficient.

$$Q_c \approx \frac{1pC}{22.5 MeV} \sum_{i=1}^N \alpha_i E_i \quad (3.8)$$

For a better comparison the RPP collection coefficient was fitted using 130nm technology data, and for these computations was set to be 0.8. The collection coefficients used by both RPP and IRPP were set using the same 130nm data, used by the ADC. In Fig. 3-5, the total collected charge for the devices from different technology nodes is shown, along with collected charge TCAD data from the literature and estimations using both RPP and IRPP models.



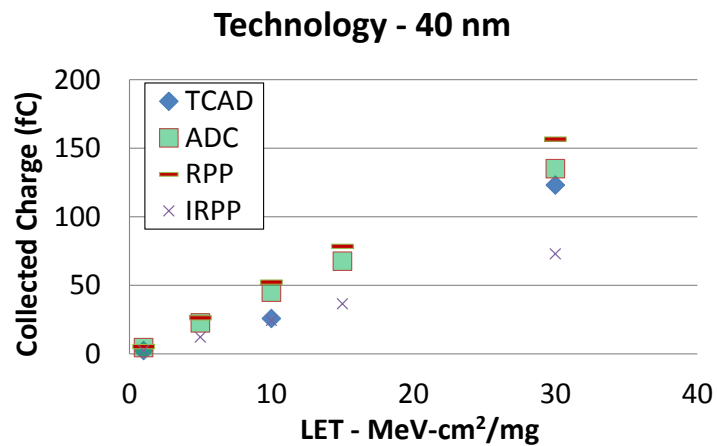
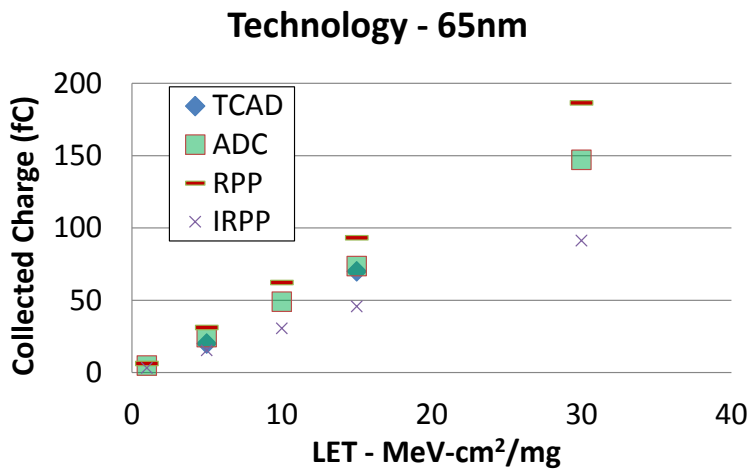
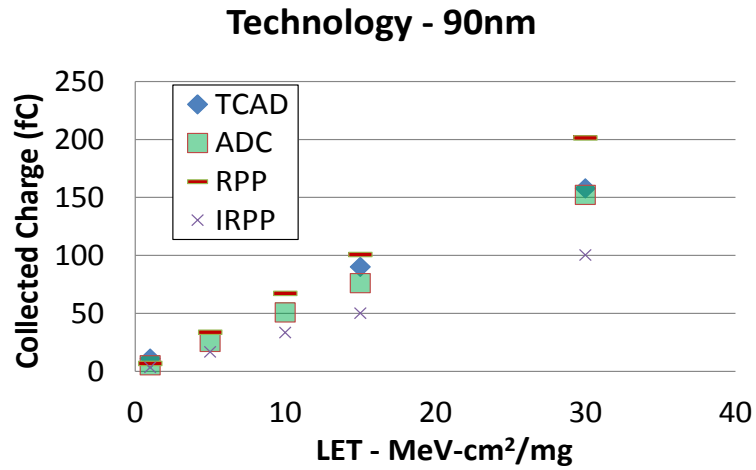


Figure 3-5. Simulation results for Collected Charge for proposed ADC model and TCAD data in the literature show excellent agreement for all technology nodes considered in this work.

The PMOS transistor data from the proposed model are compared with data from [26, 40, 41] and [20] for the 40, 90, 65, 130 nm technologies, respectively. The TCAD data mentioned before were validated, by verifying both the devices used in simulations and the physical models set in the simulations. These simulations were carried out using standard drift-diffusion models, carrier-carrier scattering models and both Shockley-Read-Hall and Auger recombination models. The heavy ion simulations were set with a track long enough to cross the devices, and the charge was distributed along the track with a Gaussian profile, a common practice when simulating soft error events.

For the technologies compared in this work, the results indicate that the values of collected charge estimated by the proposed ADC model are very close to the TCAD data reported in the literature. For very low particle LET values, $\sim 1 \text{ MeV}\cdot\text{cm}^2/\text{mg}$ all models show excellent agreement with TCAD data. However, as the particle LET values increase, the collected charge estimations of both RPP and IRPP models start to deviate from TCAD data.

3.6. Error Comparison between ADC, RPP and IRPP

For a better comparison between the three methodologies an error estimator (ϵ) is defined here as the absolute value of the difference between the charge estimation Q_c and the TCAD data Q_{tcad} and is given by $\epsilon = |Q_c - Q_{tcad}|$. The average error (μ) and sigma (σ) of the data shown in Fig. 3-5 are shown in Fig. 3-6.

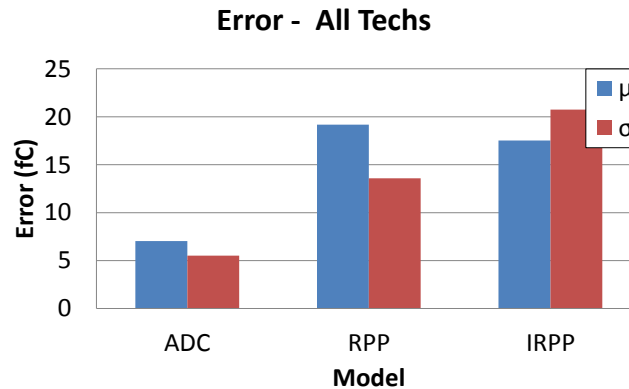


Figure 3-6: Mean and standard deviation for error in estimating the collected charge for the proposed ADC model, RPP model, and IRPP model when compared to TCAD results.

The expected error from the proposed ADC model is smaller compared with that of both the RPP and IRPP models. The proposed model shows a difference of approximately 7.1 fC, while the RPP and IRPP models show the same to be 19.1 fC and 17.5 fC, respectively. Fig. 3-7 shows the mean error values normalized to the proposed ADC model. The expected errors from the RPP and IRPP models were measured to be 2.72 and 2.49 times the error of the proposed ADC model. When analyzing these errors, it is clear that the RPP model consistently overestimates the collected charge, whereas the IRPP model consistently underestimates the collected charge for the not fitted technologies.

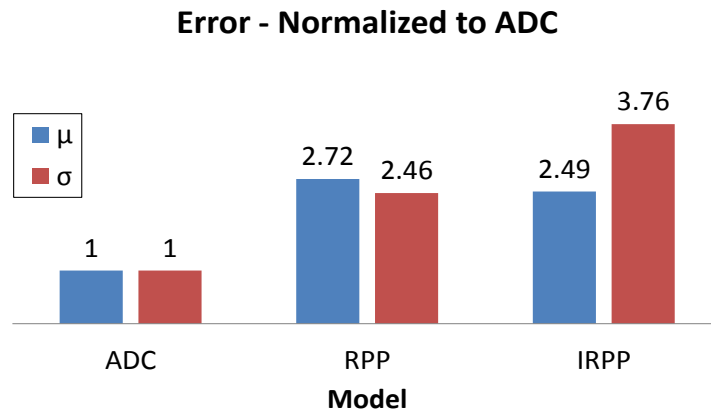


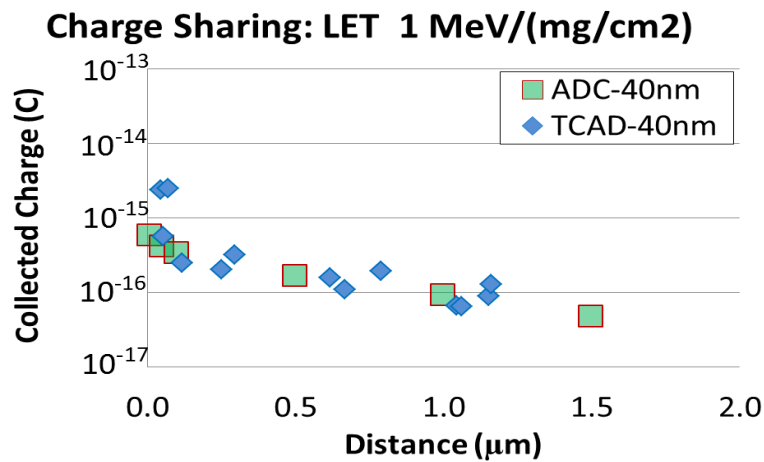
Figure 3-7: Normalized mean error of RPP and IRPP related to the ADC

Both RPP and IRPP were calibrated for improved accuracy with 130nm technology data. It is expected to have an overestimation of the RPP due to the use of a single collection coefficient for devices that have clearly more than one well defined physical region. This is exactly the improvement of the IRPP model, by allowing the user to define different sensitive volumes and respective collection coefficients. One of the problems of the IRPP is that there is no formalized methodology to define both the geometry of the sensitive volume and its collection coefficient, a problem that is solved by the ADC. Another key difference between these models is that regions defined by the ADC have clear and well defined physical characteristics, as discussed in section 3. The ADC model was able to consistently outperform both the RPP and IRPP methodologies with the same pre-characterization data step and without requiring any special assumptions or sensitive volume evaluations.

3.7 Charge Sharing

The proposed model is extended to include the charge sharing effects. In this case, the analysis is limited to comparison with 40-nm TCAD data due to the unavailability of data for comparison for other technology nodes. The goal is to estimate the amount of charge collected by transistors that are not directly hit by the primary radiation event

The charge-sharing mechanism for multiple transistors is evaluated using Eqns. (3.5-3.7). These calculations are performed for 1, 10 and 30 MeV*cm²/mg particles and results are shown in Fig. 3-8-a,b and c, respectively, along with published data [26]. Notice that the TCAD data used in Fig. 3-8 average the collected charge of both the PMOS and NMOS transistors. As a result, published data always show the average values for collected charge, whereas results for the proposed ADC model shows separate values for n-hit and p-hit. For low LET the model predicts with accuracy the collected charge, but its accuracy is decreased with the LET increase. The charge sharing case imposes additional complexity to the model, due to the more complex structure of the TCAD device. The data shown in Fig. 3-8, was extracted from 2D TCAD simulations, with an array of transistors containing both p+n and n+p devices.



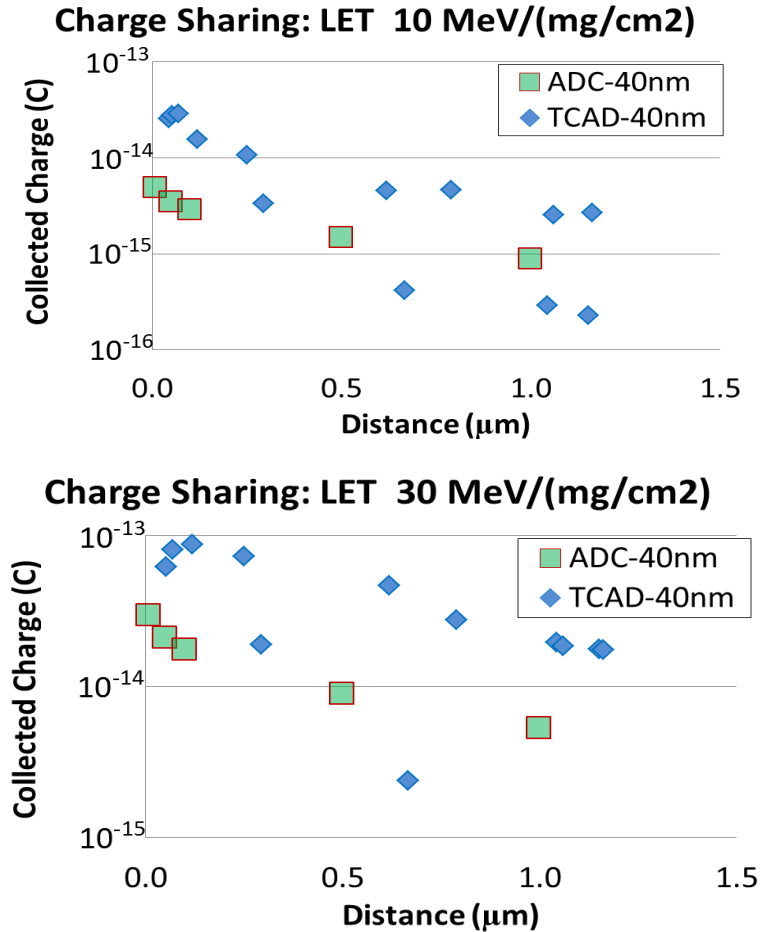


Figure 3-8 Collected Charge due to charge-sharing as a function of distance from the hit location for a 40 nm node. TCAD data from the literature for a particle with (a) LET equal to 1 MeV/mg/cm², and, (b) LET equal to 10 MeV/mg/cm², (c) LET equal to 30 MeV/mg/cm²

This in turns make the modeling more complicated since now the TCAD device has multiple junctions collecting charge and multiple well/substrate contacts influencing the parasitic bipolar amplification. The influence of the injected charge in the quasi-neutral region on the parasitic bipolar amplification will now be influenced by the position of multiple well/substrate contacts, in addition to other nearby junctions in the same well collecting charge. It was also noticed the collection of charge in devices in different wells. The ADC model is not able to accurately model all these conditions that will have a stronger influence with higher LETs. However considering the simplicity of the ADC, the accuracy does serve as a first order estimation for circuit level fault injection.

3.8 Conclusions

It is shown that the ADC model can be used to estimate the collected charge for multiple transistors in bulk nanometer scale technologies. Through technology characterization by TCAD simulations, a single omega function is shown to be adequate to model the collected charge for the hit node and other nearby nodes. When comparing the ADC model with both RPP and IRPP models, the error of these methodologies was measured to be about 2.72x and 2.49x higher respectively when compared with the ADC results. It was also shown that the proposed ADC extension was able to accurately estimate charge sharing, showing good agreement between estimates and 40-nm TCAD data. The simple formulation of the ADC model and its accuracy for collected charge estimation make this model a great candidate to be integrated into the EDA design flow and for circuit level fault injection simulations.

CHAPTER 4

Single Event Transient and Soft Error Cross-Section Modeling

With the collected charge at a circuit node calculated, the focus is shifted to the analysis of the standard cell response to the radiation event. For standard cell soft error analyses, the single-event transient pulse characteristics shown in Figure 4-1, are absolutely necessary as they are the dominant factors determining the circuit soft error performance. The model developed for this thesis calculates both the SET pulse width (measured at half VDD) and the voltage peak of the SET. For a given flip-flop design, if the SET pulse width exceeds the feedback loop delay, it will cause an upset [2]. For combinational logic errors, SET pulse width and clock frequency together determine the probability for an error. Also, SET pulse width strongly influences the electrical masking effects for logic errors. As a result, determination of SET pulse characteristics is of paramount importance for failure analysis, predictive capability, and mitigation of soft errors.

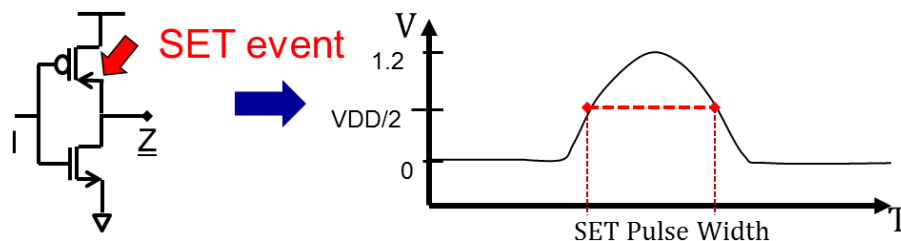


Figure 4-1 The SET pulse width generated by a Soft Error, measured at half VDD.

The SET pulse characteristics depend on the fabrication process parameters, incident ion characteristics, layout, and circuit topology. Simulation of all these parameters requires mixed-mode simulation capability with the ion-hit region and surrounding regions of interest modeled in 3D TCAD and the rest of the circuit using compact models. A single simulation run may take days to complete with hundreds of runs necessary to get statistically significant results. The computation complexity of such a task has turned researchers to experimentally measure SET pulse characteristics. Unfortunately, these

measurements only yield a final statistical distribution of the SET pulse characteristics. While extremely useful, these results do not yield any information regarding hit location, collected charge, charge sharing and circuit topology effects. In this chapter, analytical models have been developed to allow the estimation of SET pulse characteristics taking into account the parameters of interest and yielding the relevant information needed by the engineers. The proposed model is fast, accurate, and easy to integrate into existing design flows.

4.1 SET Pulse Width Model

The characterization of SET pulse requires accurate estimation of the collected charge at each circuit node, charge sharing, and the logic cell response to the injected current. The model proposed in this chapter requires the use of the ADC model presented in chapter 3 and fault injections to generate 3 lookup-tables. The first lookup-table holds the reference SET pulse width of minimum sized inverters, the second and third lookup table stores multiplying factors for the restoring current and load capacitance. The reference SET pulse width from an inverter (lookup table #1) along with the current and capacitance values (lookup tables #2 and 3) are used to estimate the SET pulse width for any given circuit node. The use of lookup-tables to estimate SET pulse widths had been used before [42, 43], but they were used to characterize each single standard cell, a task that requires a large number of lookup-tables as opposed to 3 tables as proposed in this thesis. These empirical parameters are evaluated only once for a given technology. Based on these parameters, SET pulse characteristics for all digital standard cells for a given technology can be estimated for different LET values. The SET pulse width is decomposed in a reference SET pulse width, restoring current factor and load capacitance factor, these are shown as SET influence factors in Figure 4-2.

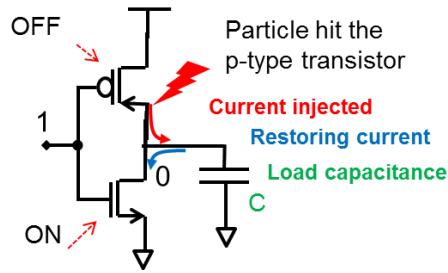


Figure 4-2 Circuit effects influencing cell response during an SEE

The lookup-table with SET pulse width from minimum sized inverters (Pw_{min}) is built by performing fault injections with different LET values (translated into collected charge by the ADC model) for each transistor in the cell. Then Eqn 4.1 is used to estimate the SET pulse width for any given node.

$$SET_{Pw} = Pw_{min}(Q_c) * Rest_{current}(R_T) * Cap_{load}(Cl) \quad (4.1)$$

The $Rest_{current}$ factor in Eqn. 4.1 maps the effect of the total resistance that connects the output node to a power rail to the SET pulse width. $Rest_{current}$ is estimated using an equivalent transistor for an array of transistors in series and parallel as shown Figure 4-3.

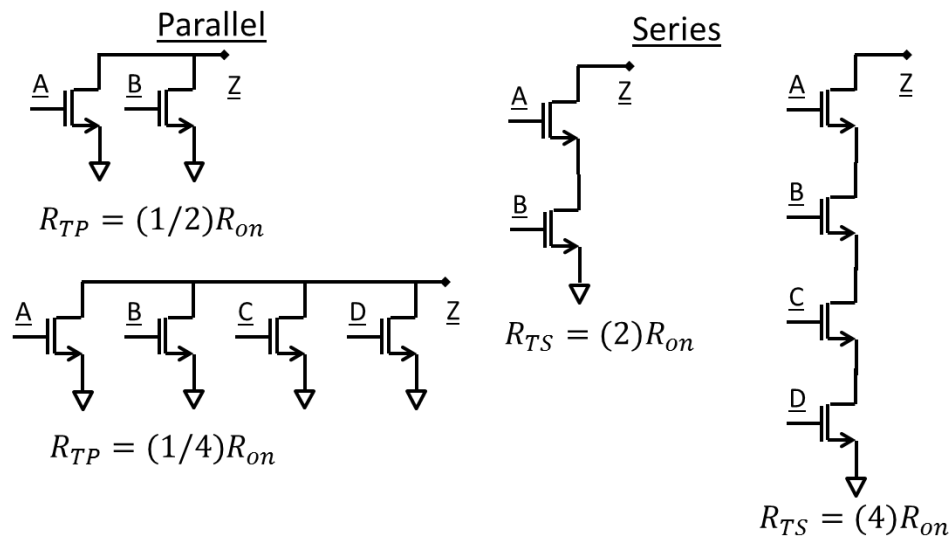


Figure 4-3 Circuits used for the technology characterization

In figure 4-4 the simulation results for the transistors in series used to generate the fit functions for the $Rest_{current}$ are shown for the evaluated technologies. Notice that for all technologies, and for both low and high collected charge values, the decrease in the restoring current (with increasing number of transistors in the path) will increase the SET pulse width.

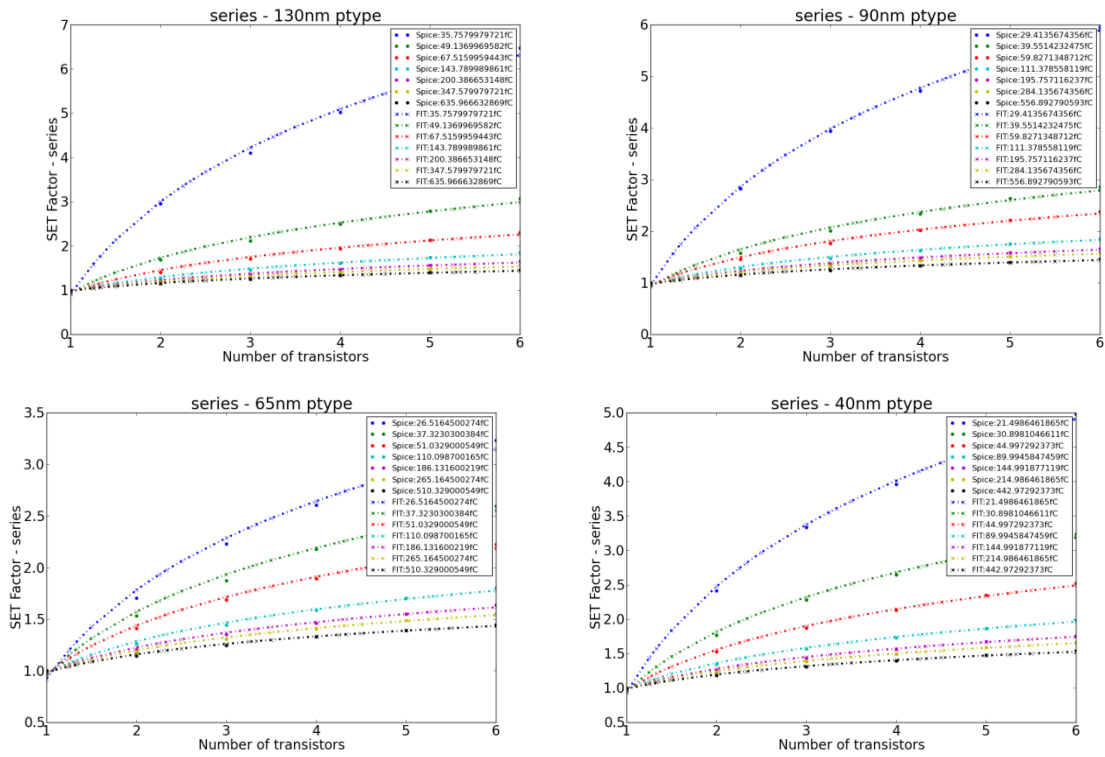


Figure 4-4 Restoring current factor for transistors in series

In figure 4-5 the simulation results for the parallel transistors to generate fit functions for the $Rest_{current}$ are shown for the evaluated technologies. The increase of restoring current (with increasing number of transistors in parallel) will produce a factor that decreases the SET pulse width. Note that this information is stored in the lookup table using the equivalent on-state resistance of these transistors.

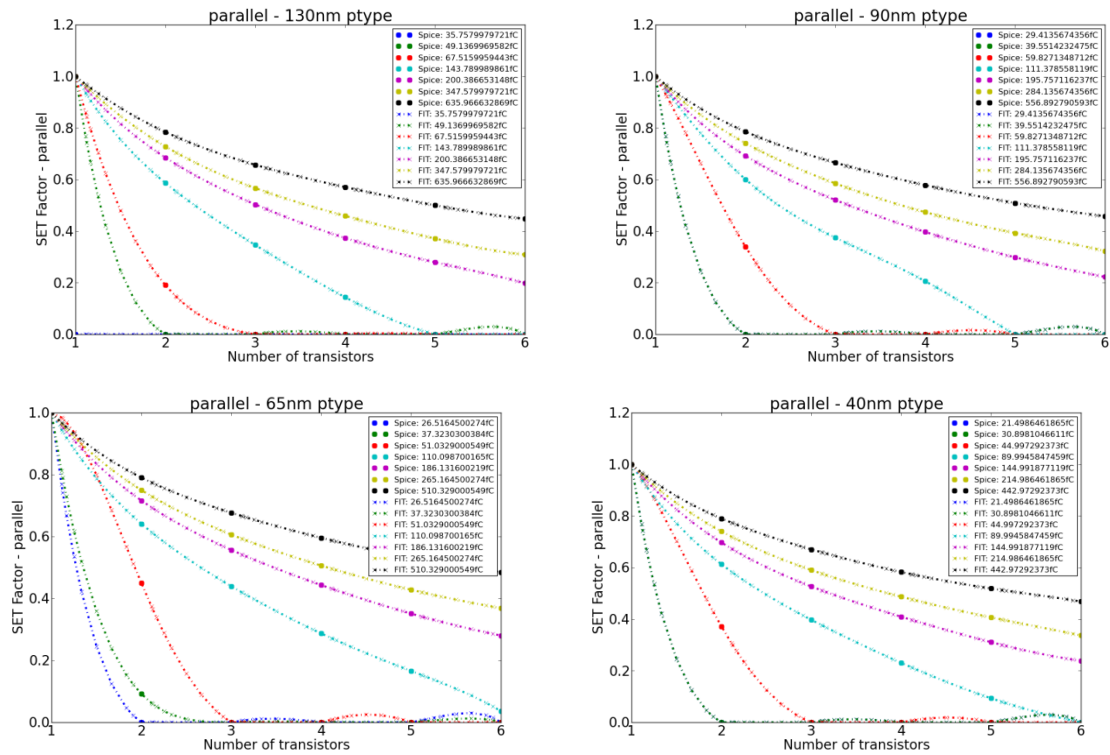


Figure 4-5 Restoring current factor for transistors in parallel

The Cap_{load} factor is the effect of the load capacitance at an output node on the SET Pulse width. Lookup Table #3 is used to map different capacitance values at an output node and the reference SET pulse width from an inverter to estimate the SET pulse width at the node in question. In figure 4-6 the multiplication factor as a function of Cap_{load} is shown for 130, 90, 65 and 40nm technologies.

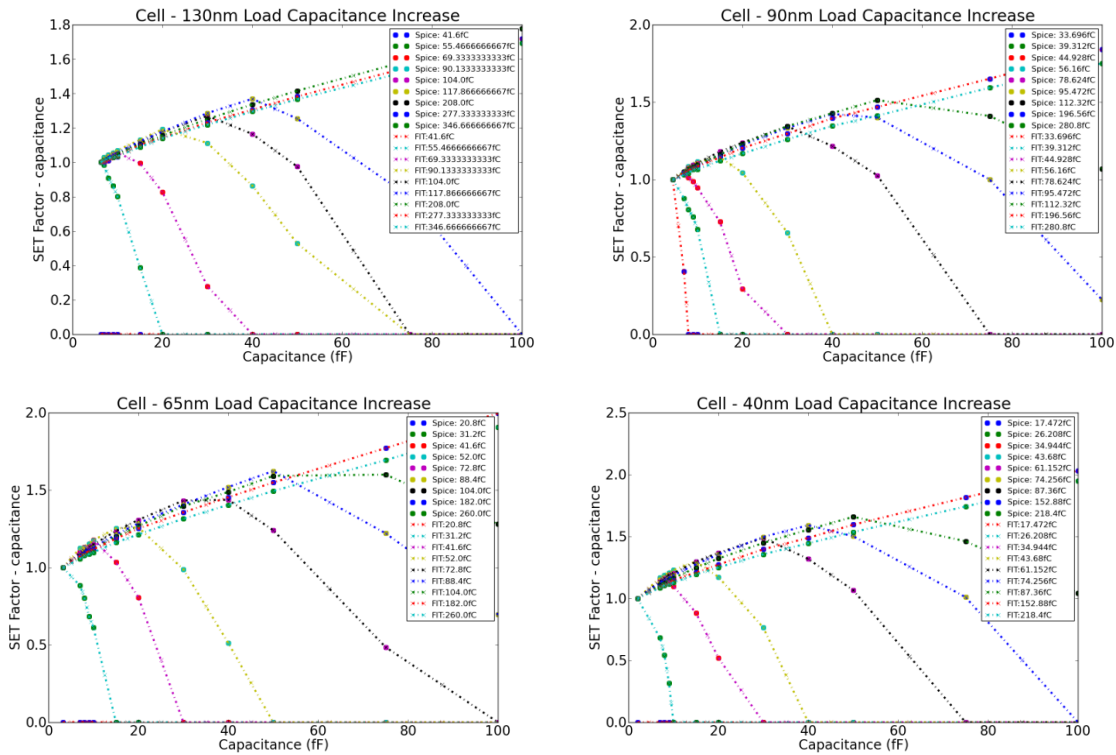


Figure 4-6 Load Capacitance Factor

Fig. 4-6 shows an important trend that should be carefully understood. By analyzing the results from these technologies it is possible to notice that the SET pulse width will be reduced if the load capacitance is increased or if the collected charge is decreased (near 17 fC). However, for energy depositions much higher than the 17 fC value, a threshold is seen beyond which increasing the load capacitance will decrease the SET pulse width. This in turns indicates that this is not a reliable parameter to be used for SET mitigation, since variations in the energy deposition will produce different trends. Also notice that for most of the technologies, a high load capacitance is required to reduce the SET pulse width. Also, increasing the capacitance to mitigate the SET pulses will increase the netlist delay, and may not guarantee the SET pulse mitigation.

An example of the SET pulse width analysis of an OAI21_X1 standard cell is shown in Figure 4-4. For the input conditions given, only two transistors (those with gates tied to A and C inputs) will be turned ON, resulting in the resistive network shown in the figure.

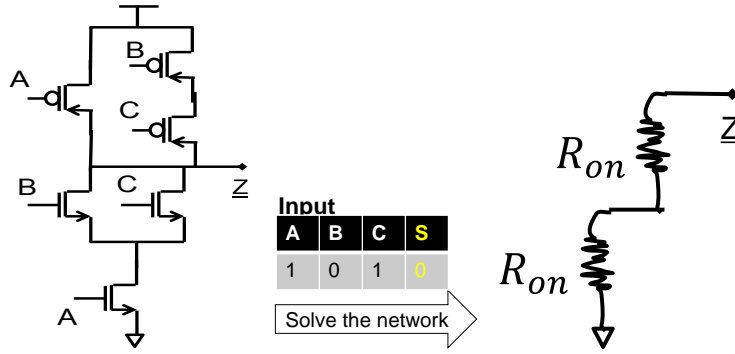


Figure 4-7 OAI21_X1 Cell on-state resistance modeling

Since these two transistors are the only ones connecting the output node to a power rail and they are in series, the equivalent resistance will be given by $(R_{on_c} + R_{on_a})$. Using such resistive network representation and the nodal capacitances, the SET is estimated using lookup-tables discussed previously. If multiple cells collect charge, it will result in multiple SETs. In the presence of multiple SET pulses within a cell (for example cells that contain more than just an inverter, NAND, or NOR) may have multiple SET's that either combine to yield a longer SET pulse, or may remain independent, or quench each other yielding a shorter SET pulse (this has been referred to as SET pulse quenching [12]) depending on circuit topology. All these factors are taken into consideration to yield a final SET pulse for a given ion incident at an angle at a given location for a cell operating at a given voltage to yield accurate single-event pulse characteristics.

Since generating additional SET pulses in the analysis is a straight forward process, the focus in this section is on pulse-quenching effects. To calculate the SET pulse quenching, the netlist is evaluated searching for an inversion of the signal within the cell logic stages. This condition was first evaluated by [12]. If from the logic stage i to $i+1$, the signal is inverted, the SET Pulse Quench might occur if enough charge is collected by both the stages. As discussed before, one can obtain the charge sharing at an adjacent node with Q_{ci} , and the quenching condition can be easily evaluated in the netlist graph. With this basic framework set in place, the SET pulse width of a cell subjected to SET pulse quench is given by (4.2).

$$SET_{Pw_quench} = \begin{cases} SET_{Pw}(Q_c) - SET_{Pw}(Q_{Ci+1}) & SET_{Pw}(Q_{Ci+1}) > 0 \\ SET_{Pw}(Q_c - Q_{Ci+1}) & else \end{cases} \quad (4.2)$$

In the equation the first case considers that enough charge caused a signal inversion in the node i+1 and for the second case an inversion didn't occur in the adjacent node.

4.1.1 Models and Operation Condition

For the fault injection simulations, the LET values of the particles were in the range from 1 to about 30 MeV*cm²/mg, as discussed in chapter 3. In table 4-1 few important parameters used to design the test benches used for the verification process and the model fitting are shown. The 130-nm, 90-nm, 65-nm and 40-nm technologies were evaluated in Virtuoso tool [44], using spice Predictive Technology Models (PTM) from ASU [45].

Table 4-1 Circuit and operation condition parameters

	130nm	90nm	65nm	40nm
Stand. p-transistor width	620nm	525nm	430nm	350nm
Stand. n-transistor width	250nm	210nm	170nm	140nm
Load capacitance	6.3fF	4.5fF	3.2fF	2fF
Power Supply	1.5 V	1.2 V	1.2 V	1.1 V

The fault injection was simulated using the Bias-Dependent model [28], with its input conditions provided by the ADC model, as discussed in chapter 3.

4.1.2 SET Pulse Width Results

To validate the accuracy of the SET pulse width model, the 4 technologies mentioned in the last section were used in simulations. The following results are from fault injections in standard cells, using the Bias-Dependent Model, and a SPICE-like simulator. For the next results, the p-type transistor was considered the target cell of the SEE event. Results in figure 4-8 show excellent agreement for predicting SET pulse characteristics using the proposed model, for several different standard cells.

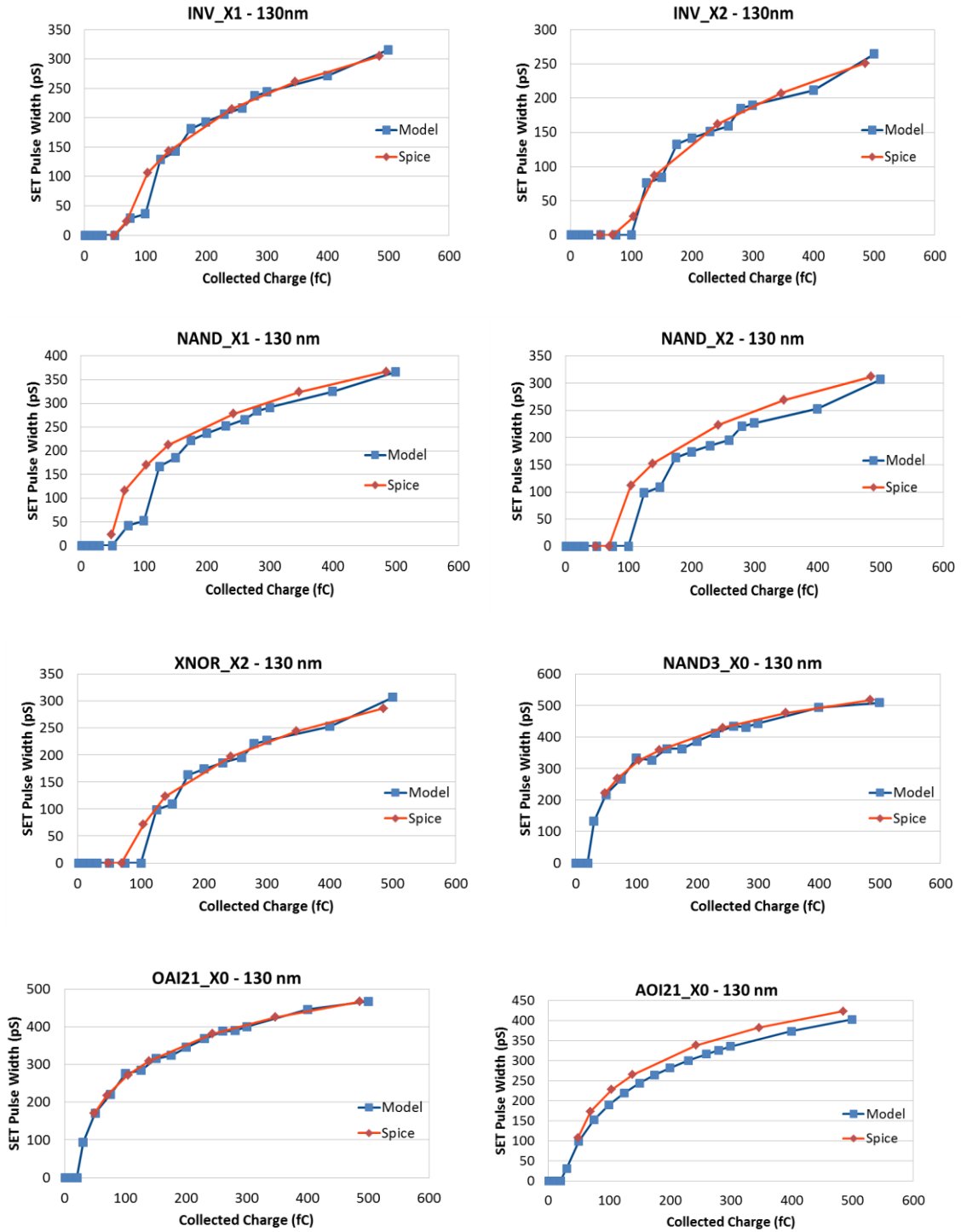
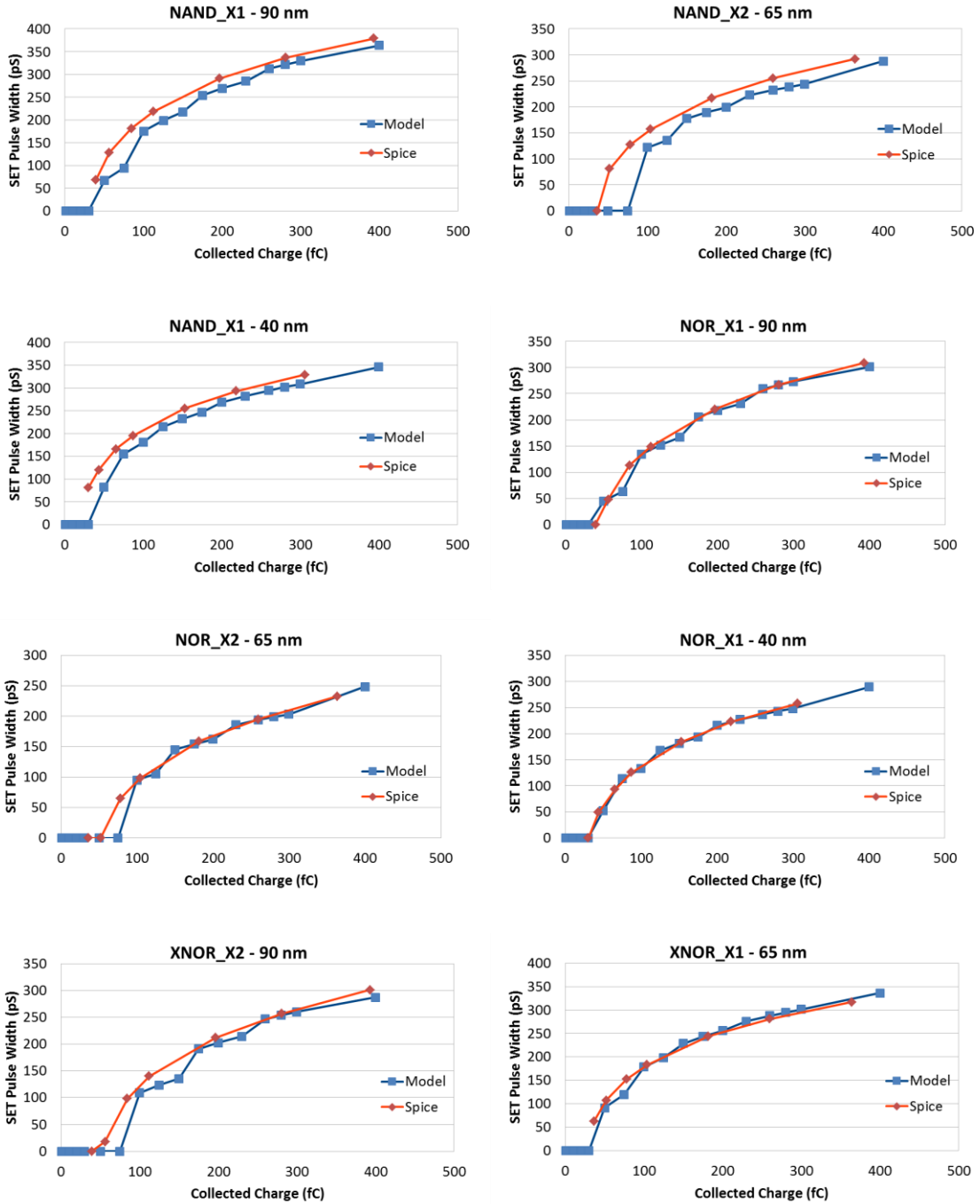
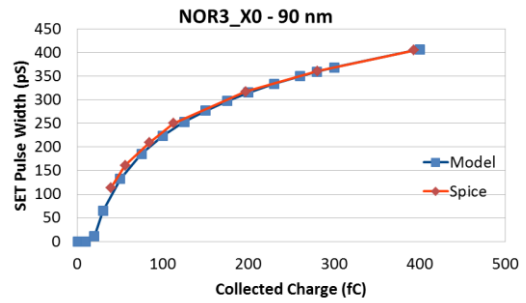
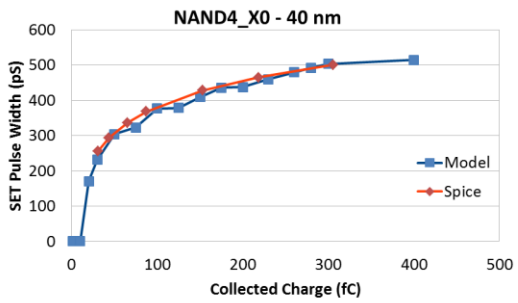
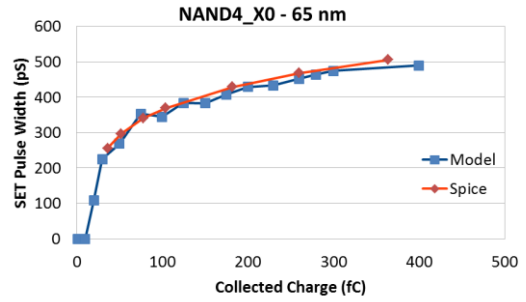
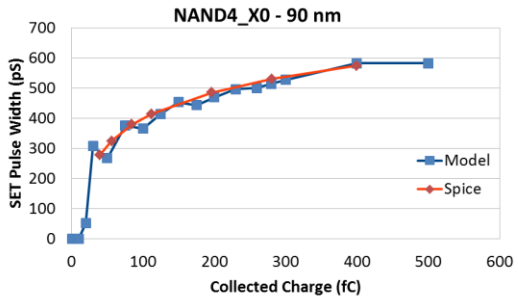
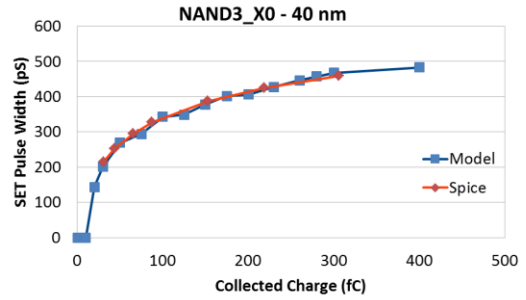
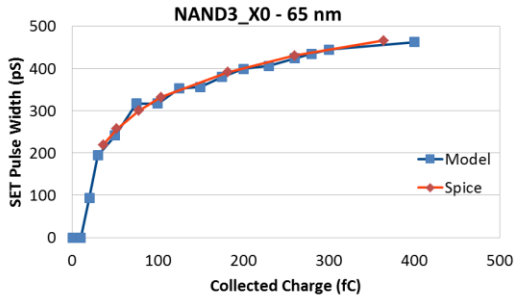
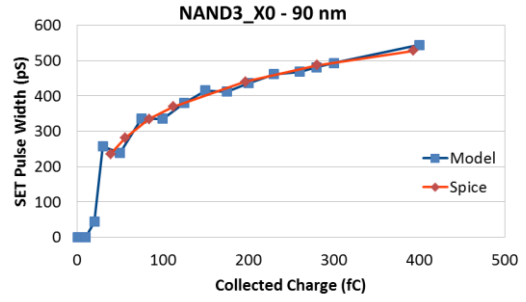
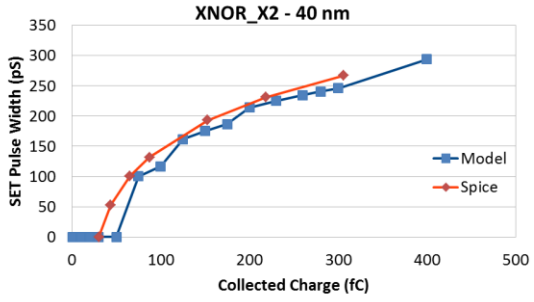


Figure 4-8 SET Pulse width from 130nm standard cells comparison between model and spice

Notice in figure 4-8 for 130-nm technology node that cells with different sizing and number of inputs were properly characterized by the SET model. In Figure 4-9, other standard cells that were evaluated for the 90-, 65- and 40-nm technology nodes are shown with the SET model displaying very good accuracy.





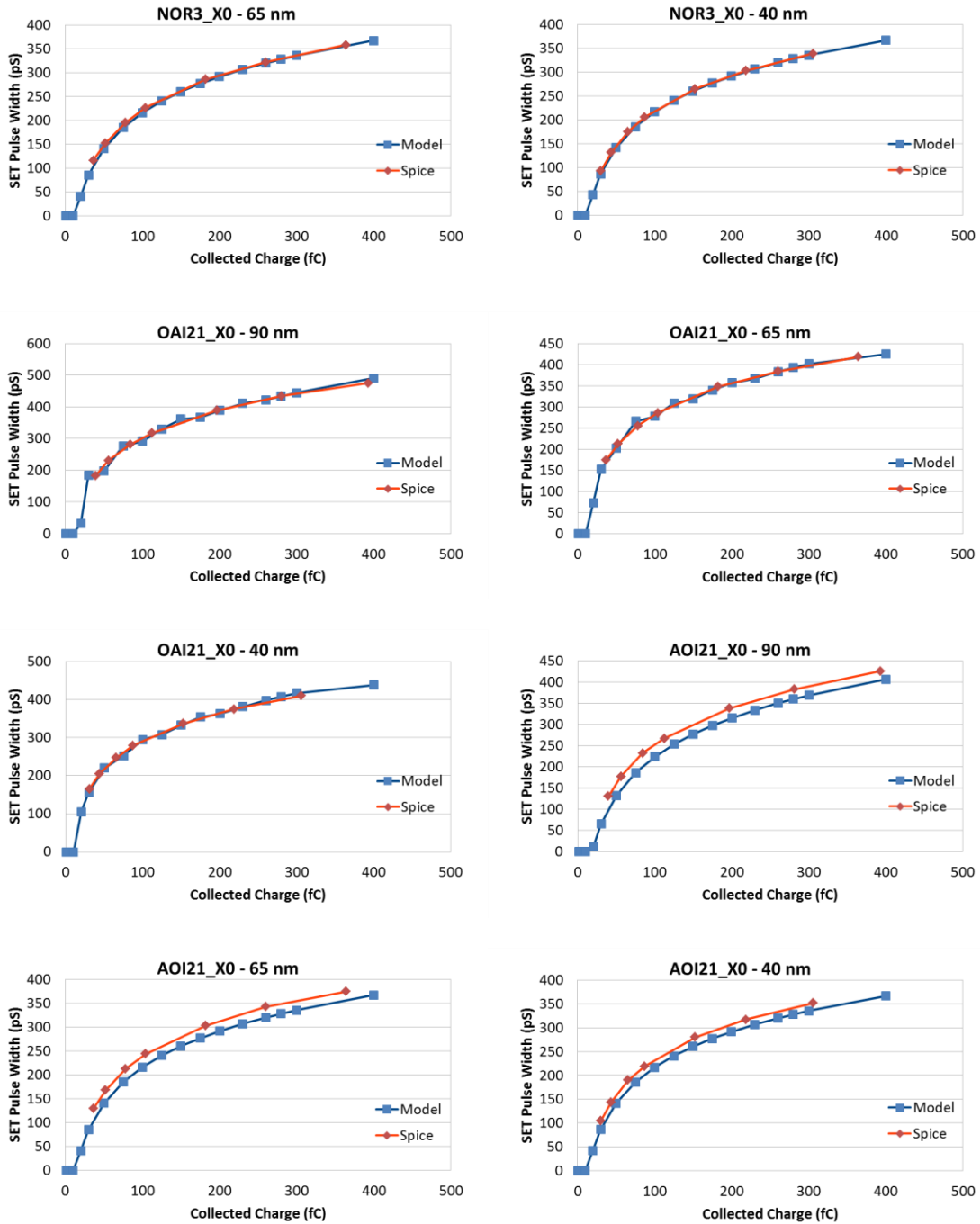


Figure 4-9 SET pulse width comparison for 90, 65 and 40nm technology nodes.

The SET pulse-quench effect model was validated by comparing fault injection in a sequence of two inverters and the estimation of the proposed model, and is showed in Figure 4-10. In this simulation the bias-dependent model injected a current in two stages of the logic simultaneously with the SET pulse width measured in the last chain stage.

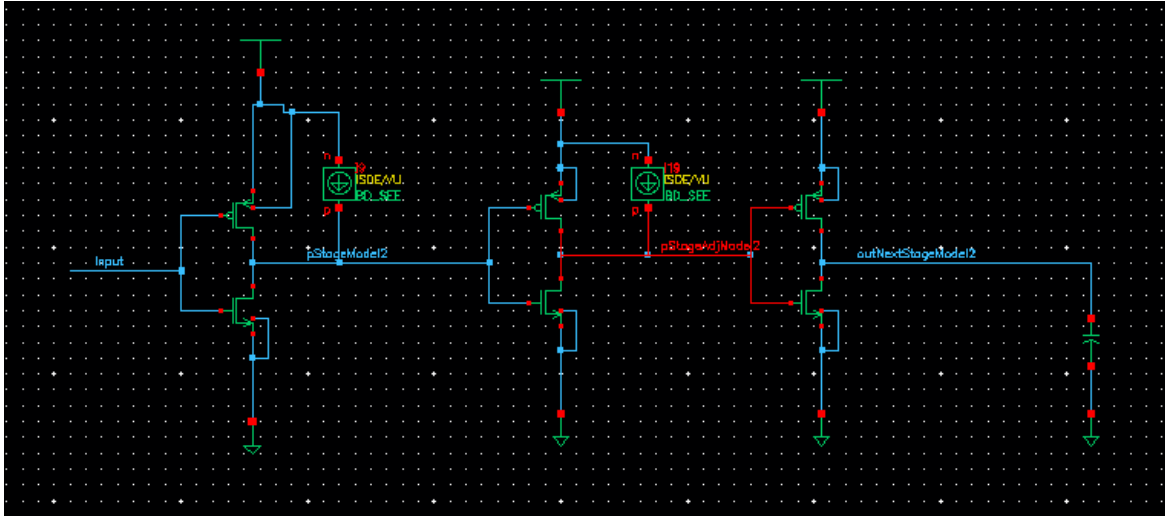


Figure 4-10 Spice benchmark used to evaluate the SET pulse quenching.

The charge collected by the first stage was calculated with Equation (3.2) and the same for the second stage was given by Equation (3.6). The results in figure 4-11 show a good agreement between the SET pulse-quench model and simulation results from the SPECTRE circuit simulator.

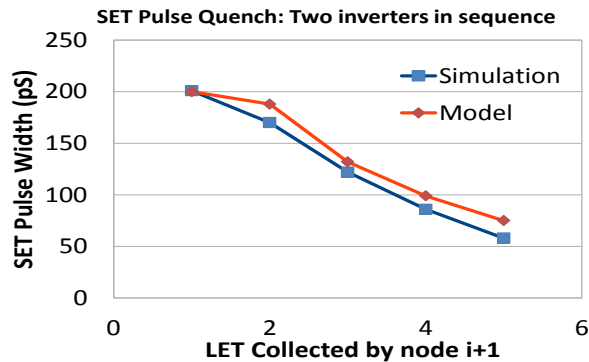


Figure 4-11 SET Pulse Quenching

The methodology was compared with experimental SET data from a 65-nm Test Chip, from [46]. In table 4-2 the data from the model and 65-nm experiments are shown, with an average error of about 29 ps, measured by averaging the absolute difference between model and experimental data. The experimental data showed in table 4-2 are from the average SET pulse widths.

Table 4-2 Model vs experimental data for 65nm inverters

LET (MeV*cm ² /mg)	Test Chip (ps)	Model (ps)
21.2	79.6	86.8
30.2	100.9	123.8
58.8	135.5	193.8

Analyzing table 4-2 data, it's possible to notice that for the LET range below 30 MeV*cm²/mg the model has very good accuracy, considering that the model was fitted with a predictive technology, and not with the commercial models used in the Test Chip. Above the LET range indicated, the error does increase but this energy range is not considered in this work. This work is geared towards terrestrial effects and most ions encountered in terrestrial environment will have an LET value of less than 30 MeV*cm²/mg.

4.1.3 SET Pulse Width Model Conclusions

A methodology to estimate the SET pulse width taking into account the charge sharing has been developed and demonstrated. The model was compared with both simulation and experimental results showing an excellent agreement between the model and the data. The analysis requires a technology characterization by performing a fault injection campaign on 3 custom circuits for parametric analysis to characterize the SET pulse width. The proposed model reduces the computational time by orders of magnitude while yielding accurate results for the SET pulse characteristics analysis. With this framework developed, the impact of SEE on circuit nodes can now be evaluated.

4.2 Circuit De-rating Factors

Before going over the details of the metrics to evaluate the circuit soft error cross-section, it is important to introduce circuit derating mechanisms that are going to impact the circuit response to SEE. The radiation effects community identifies the Soft Error Generation Probability, Electrical De-rating,

Logical De-rating and Time De-rating as the circuit probabilities to model SE effects in logic circuits [47, 48]. These de-rating probabilities are shown in figure 4-12, with its circuit representation.

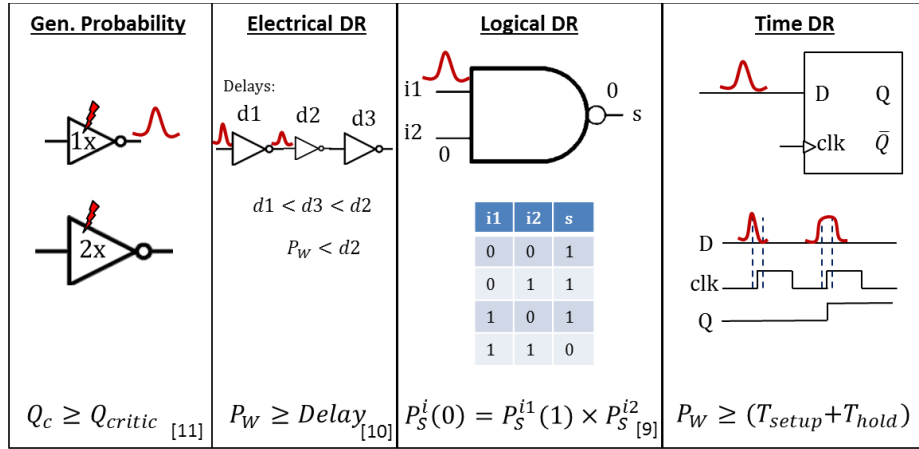


Figure 4-12 - De rating probabilities. Electrical, Logical and Timing DR respectively.

The Generation Probability gives the probability that a charged particle will generate an SET in the standard cell. The standard cell characteristics discussed in the last 2 chapters will impact this parameter, along with the load capacitance associated with the evaluated standard cell. For Flip-Flops the term Critical Charge is used to indicate the smallest amount of charge required to upset a sequential element, and for this work is used in the context of Standard Cells. The Generation Probability (Gen_p) is given by Equation 4.3, where SET_{Pw} is the resulting SET pulse width given the collected charge, SET_{Vpeak} is the SET voltage peak, and V_{Node} is the node voltage prior to the SEE. The generation probability is 1 if sufficient charge was injected to generate an SET and if the difference of voltage between the SET voltage peak and the node voltage prior to the SEE is greater than 70% of the supply voltage. The 0.7 factor is verified empirically by noting that the SET will not propagate if the SET voltage peak is not beyond a certain voltage, in this case 0.7, and it was verified by SPICE-like simulations.

$$Gen_p = \begin{cases} 1 & SET_{Pw} > 0 \text{ and } |SET_{Vpeak} - V_{Node}| \geq VDD * 0.7 \\ 0 & \text{else} \end{cases} \quad (4.3)$$

The Electrical De-rating probability, also commonly referred as Electrical Masking, calculates the probability that the circuit will electrically filter the SET pulse. The RC characteristics of interconnections and the transistors in the cell determine this probability. For instance, it is known that if the SET pulse

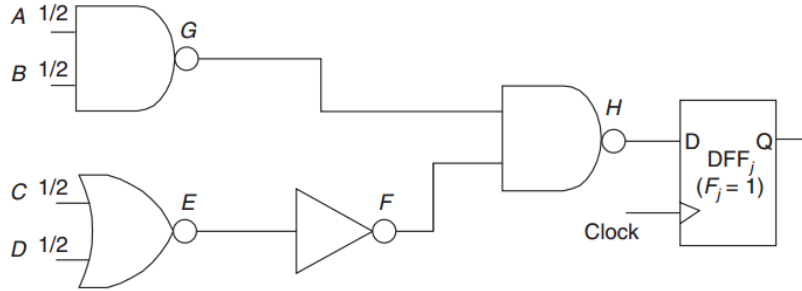
width (SET_{Pw}) is smaller than the gate delay, the SET will not be able to propagate to the next stage, and thus is masked. To calculate, the probability that an SET generated in a standard cell will propagate to the next stage requires the evaluation of the whole network/circuit. The computation of the electrical de-rating probability El_{DR} does require a delay analysis over the netlist, by analyzing the circuit graph from the primary inputs to primary outputs calculating the delay for each standard cell and interconnection. This analysis can be performed using a modified Breadth-First-Search (BFS) algorithm to traverse the circuit graph. From the input to the outputs, the interconnection delay and gate delay are calculated for each node in the graph with the maximum delay associated with each node output stored. The minimum delay path identifies the minimum pulse able to propagate through that path, yielding the El_{DR} equation (4.3).

$$\begin{aligned}
 k_i &= \{Delay_0, \dots, Delay_n\} \text{ defines a path} \\
 k_i \in K &= \{k_0, \dots, k_l\}, \text{ set of paths for the given cell} \\
 El_{DR} &= \begin{cases} 1, & SET_{Pw} \geq \text{Min}\{Max_{delay}(K)\} \\ 0, & \text{else} \end{cases} \quad (4.3)
 \end{aligned}$$

The Logical De-rating, or Logical Masking, refers to the probability that the SET will be masked by the circuit logic. One simple example is the two input NAND cell. If any one of the inputs has logic value 0, the output of the gate will be 1, no matter the state of the other input. If an SET pulse comes along the second input, it will not propagate to the NAND output, thus being masked. The Logical Derating factor is calculated using the methodology proposed by [48] and shown in figure 4-13. The algorithm works by first measuring the probability of each gate output for being in the logical state 0 or 1, starting from the primary inputs (PI) to the primary outputs (PO). Once the PO is reached, the algorithm traverses the netlist backward, starting from the PO calculating the propagation probability of each input. The propagation probability of an input is given by the probability of all the other inputs associated with an output being in a non-controllable state. For the PO, the propagating probability of connecting cells is 1, since they are directly connected to the output.

Step 1: Forward breadth-first search

- $Pr^E(1) = Pr^C(0) \times Pr^D(0) = 1/4$
- $Pr^G(1) = 1 - Pr^G(0) = 1 - [Pr^A(1) \times Pr^B(1)] = 3/4$
- $Pr^F(1) = Pr^E(0) = 1 - Pr^E(1) = 3/4$
- $Pr^H(1) = 1 - Pr^H(0) = 1 - [Pr^F(1) \times Pr^G(1)] = 7/16$



Step 2: Backward breadth-first search

- $P_{prop}^F = Pr^G(1) \times P_{prop}^H = 3/4$
- $P_{prop}^G = Pr^F(1) \times P_{prop}^H = 3/4$
- $P_{prop}^E = P_{prop}^F = 3/4$
- $P_{prop}^C = P_{prop}^E \times Pr^D(0) = (3/4)(1/2) = 3/8, P_{prop}^D = \dots = 3/8$
- $P_{prop}^A = P_{prop}^G \times Pr^B(1) = (3/4)(1/2) = 3/8, P_{prop}^B = \dots = 3/8$

Figure 4-13 Logical Derating Calculation [48]

For a 2 input NAND gate the probability of the output being in logic state zero is $P_r(0) = P_a(1) * P_b(1)$ and for being in the logic state one is $P_r(1) = 1 - P_r(0)$ as shown in figure 4-4. To calculate the cell state probability requires both the logic truth table of the respective standard cell and the logic states of the primary inputs.

Pr	A	B	S	
NAND	0	0	1	$P(0) = P_a(1)*P_b(1)$
	0	1	1	$P(1) = 1 - P(0)$
	1	0	1	
	1	1	0	

Figure 4-14 NAND gate logic state probability

And last, the Time De-rating, known also as Timing Masking, refers to the probability that the SET pulse is latched by a memory element. For a given rising edge Flip Flop (FF), the SET pulse width has to be able to reach the FF input with the clock edge, and have width equal or greater than the Flip Flop setup and hold time. The Timing de-rating is given by the Clock De-rating (C_{DR}) probability and the Flip Flop De-Rating probability (FF_{DR}). The C_{DR} is controlled by the clock period (T) and the FF_{DR} is controlled by the FF Setup time (T_{Setup}) and FF Hold time (T_{Hold}). These timing parameters define the time requirements for the SET_{Pw} to be latched by a Flip-Flop.

$$C_{DR} = \begin{cases} SET_{PW}/T & SET_{PW} \leq T \\ 1 & else \end{cases}$$

$$FF_{DR} = \begin{cases} 1 & SET_{PW} \geq T_{Setup} + T_{Hold} \\ 0 & else \end{cases}$$

$$T_{DR} = C_{DR} \cdot FF_{DR}. \quad (4.4)$$

With the last model discussed in this section, it is possible to develop metrics capable of analyzing the soft error impact on the whole circuit netlist, and not just for a single node/cell. This is the subject of the next section.

4.3 Soft Error Sensitivity and Cross Section Estimation

The models discussed in the last section are used to establish the basic framework to analyze the circuit soft-error cross-section and the circuit soft-error sensitivity index. The soft-error cross-section is the metric used to analyze the effective number of errors/chip area for a given circuit. The circuit soft-error sensitivity index is a simpler metric used to estimate the increase/decrease of the severity of the soft error event for the netlist. Notice that the soft error sensitivity index does not directly translate into a countable error.

The formulation of the circuit soft-error sensitivity index is shown in equations (4.5 and 4.6). These equations are a commonly used approach to evaluate the soft-error sensitivity of netlists in high abstraction models and it was used before by [48, 47]. This work expands the formulation by considering charge sharing, as shown in the second term of Eqn. 4.6. The circuit soft-error sensitivity index is given by the summation of individual nodes soft-error sensitivity as given by Eqn. 4.5. The nodal soft-error sensitivity index is given by the multiplication of the generation probability, electrical, logical and time de-rating factors.

$$N_{SP}(i) = Gen_p(i) \cdot El_{DR}(i) \cdot Lo_{DR}(i) \cdot T_{DR}(i) \quad (4.5)$$

$$C_{SP}(G) = \sum_{i=0}^N (N_{SP}(i) + \sum_{j=0}^K N_{SP}(j)) \quad (4.6)$$

In equation 4.6, for each standard cell evaluated, the neighboring standard cells placed there by the placement tool are also evaluated. Since their distance from the SEE hit location is known, models presented in this thesis can be used to estimate charge collected by them due to charge sharing. This is performed for both left and right cells of the hit node. A standard cell located above and below the hit cell is not considered in this computation since these cells are physically located in different regions and separated by routing channels.

This analysis is useful to quickly analyze the netlist and extract a reference soft-error measurement to make decisions during the design flow. This is possible because of the linear computational complexity $O(n)$ of the models developed in this work. However as discussed before, this equation can't be directly translated into an error cross-section. For this purpose another metric based on the de-rating factors is required. Since the main objective here is to mitigate soft error threat, estimation of cross-section can be performed at the end of the Physical Synthesis.

A novel methodology proposed in [49] and [47] provides guidance to estimate the circuit soft-error cross-section using probability matrices and fault-injection simulations. In this work, a Monte-Carlo-based methodology was adapted to remove the required high number of fault injections and it was expanded to include the charge sharing effect.

The work from [49] differ from other methodologies by analyzing the soft-error event in multiple clock cycles, and performing fault injection for the propagation analysis. It defines individual node error probabilities ($P_{C,N}^{Error}$) with a similar formulation as Eqn. 4.5 but fault injections are performed to analyze the netlist response to each clock cycle. This in turn results in an error probability matrix as shown in Figure 4-15, where each row represents a node sensitivity and each column represents the clock cycle of the respective analysis. An additional vector (A_N) is used to hold the nodal sensitivity area information to estimate the soft-error cross-section of individual nodes.

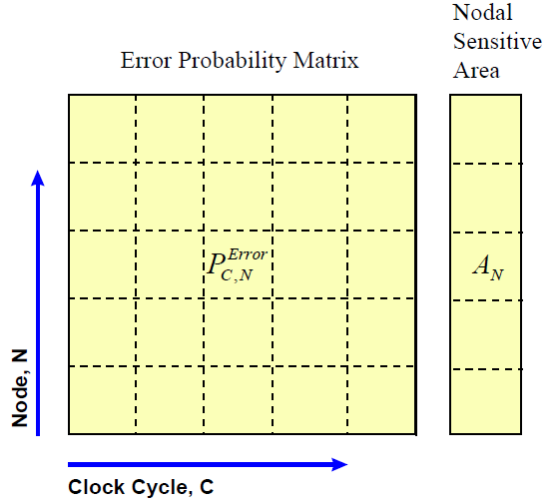


Figure 4-15 Error probability matrix of the SEUTOOL [49].

Once the error probability is calculated using Eqn. 4.7, the number of measured errors is averaged over the number of clock cycles (4.8), and the soft error cross-section is given by Eqn. 4.9.

$$P_C^{Error} = \sum_N P_{C,N}^{Error} \quad (4.7)$$

$$P^{Error} = \frac{\sum_C P_C^{Error}}{\sum_C 1} \quad (4.8)$$

$$\begin{aligned} \text{Cross - section} &= \frac{\# \text{ Observed Errors}}{\text{fluence}} \\ &= \frac{(P^{Error}) * (\text{fluence}) * (\text{Circuit Area})}{\text{fluence}} \\ &= (P^{Error}) \cdot (\text{Circuit Area}) \quad (4.9) \end{aligned}$$

This methodology was adapted in this thesis by first using an error probability matrix as shown in figure 4-15, but using equation 4.5 to estimate the individual node sensitivity. This in turns removes the required simulations used by [49]. As the analysis is required to be performed for multiple clock cycles, the state of the primary inputs used by $LO_{DR}(i)$ function will change, therefore all probabilities will have to be recomputed for each clock cycle. Another restriction is that the primary input probabilities can only have values of 1 (to indicate 100% probability of being in logic state 1) or 0 (to indicate 100% probability of being in logic state 0). This restriction allows the state probability and propagation probabilities

calculated by the LO_{DR} to be only 1 or 0, yielding to a deterministic metric that informs if the signal is able to propagate or not to the output.

Another requirement is to add an acceptance threshold in T_{DR} to convert the probabilistic information into a deterministic one, when the analysis is performed for low clock frequencies. Note that for the de-rating factors, with the LO_{DR} being guaranteed to be 1 or 0, the T_{DR} is the only non deterministic function in the equation. For high frequency and high LET particles, the condition $SET_{pw} \leq T$ is less likely, thus yielding T_{DR} to be 1. But this may not be true for all other conditions. A threshold is required to solve these cases. It may be simplified by solving the calculation of this threshold using a Monte Carlo analysis. During each algorithm iteration, a random number with a uniform distribution between 0 and 1 is sampled to estimate the time de-rating threshold. If the calculated T_{DR} is greater than the threshold the de-rating is considered to be 1 and 0 otherwise.

With these changes, the resulting calculation using Eqn. 4.5 will return 1 for an error predicted as captured by the output and 0 as masked. The charge-sharing is calculated by using Eqn. 4.5 for the respective adjacent node as discussed before. Note that this analysis considers only one SEE per clock cycle, thus the computation for matrix column will result in a matrix mostly filled by zeros, since the SEE is a local effect.

To calculate the error probability matrix, a Monte Carlo analysis is performed by defining first a vector of inputs to define the input states probabilities for each clock cycle (restricted to probabilities 1 or 0) and a reference energy deposition distribution function to be sampled by the analysis (Uniform, Gaussian or Constant). The analysis is performed by iterating over the vector of inputs and randomly selecting standard cells from the netlist in each clock cycle to be hit by the SEE.

For each cycle, the energy deposition distribution function is sampled and the ADC is used to estimate both collected charge and charge sharing. The charge is then injected in the selected standard cell and the adjacent nodes identified by the physical placement. With this information, the SET pulse width can be

estimated and equation 4.5 used, with the conditions discussed above. The resulting output is then checked and used to estimate the soft-error cross-section by using Eqns. 4.8 and 4.9.

To validate the soft error cross-section estimation the model was compared with experimental data from a Test Chip. The Test Chip was designed using a bulk, 45nm technology node and contains two combinational logic blocks, used to estimate the logic cross section. In figure 4-16 the 4 Bit Comparator cross-section is compared with results extracted by the model presented in this thesis. The 4-Bit Comparator was designed with 46 logical gates, and the soft error experiments were conducted using a flux of 5.5 MeV alpha particles from an Americium-241 source with an activity of 10 μCi [50]. The cross-section calculation considered a Gaussian energy deposition profile with average LET of 3.5 $\text{MeV}\cdot\text{cm}^2/\text{mg}$ and sigma of 0.7 $\text{MeV}\cdot\text{cm}^2/\text{mg}$. In this comparison the model underestimates the soft error cross-section by about 37.4%, a great result considering that the model was completely characterized using predictive technology nodes.

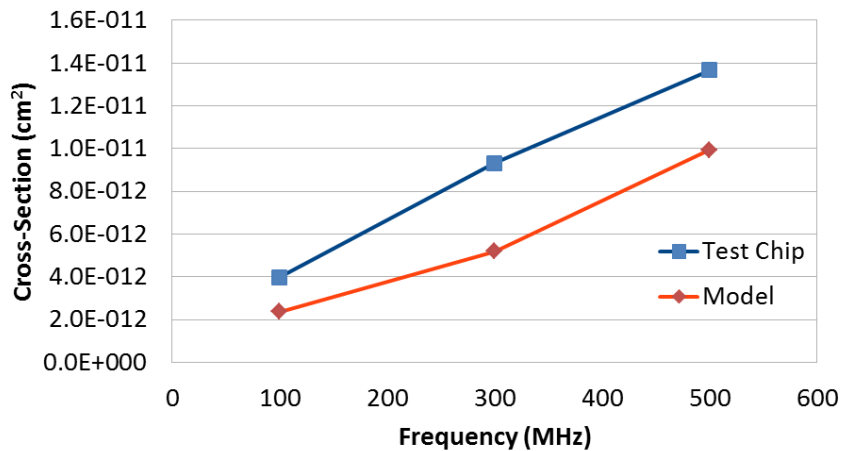


Figure 4-16 Test Chip vs Model Cross Section - 4 Bit comparator

In figure 4-17 the model is compared with the 72 Inverter chain circuit. In this circuit, chains of inverters with 12 stages are connected with OR gates to generate a single output. As observed in the figure the model also yields good results, with an average 40% overestimation of the soft-error cross-section.

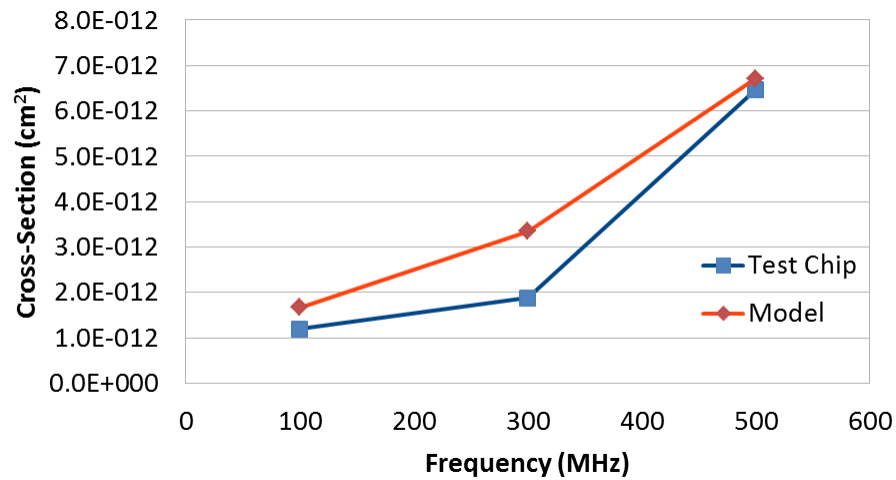


Figure 4-17 Test Chip vs Model Cross Section - 72 Inverter chain

These results show that the model is able to estimate the soft error cross-section with considerable accuracy. It is also important to indicate that the methodology was properly able to model the impact of the frequency variation, by correctly predicting the increase in soft error cross-section with the frequency increase.

4.4 Circuit Analysis Conclusions

Two metrics to evaluate the circuit response to SEE were presented and discussed, and will be used to report the soft error data in this work. With the framework to analyze soft errors for both standard cells and circuits developed, the focus now is on the Physical Synthesis process.

CHAPTER 5

Physical Synthesis

As discussed before, EDA constitutes an important research field to enable the scaling and the adoption of new technologies by semiconductor companies. The Physical Synthesis process is part of the automation flow used to convert a high level description of a circuit into an accurate map of layers to be manufactured by semiconductor foundries [5]. A typical EDA flow for an Application Specific Integrated Circuit (ASIC) circuit is shown in figure 5-1. The first two steps, Modeling and Design Verification consist of functional modeling\simulation of the design, and can be done using behavioral circuit descriptions.

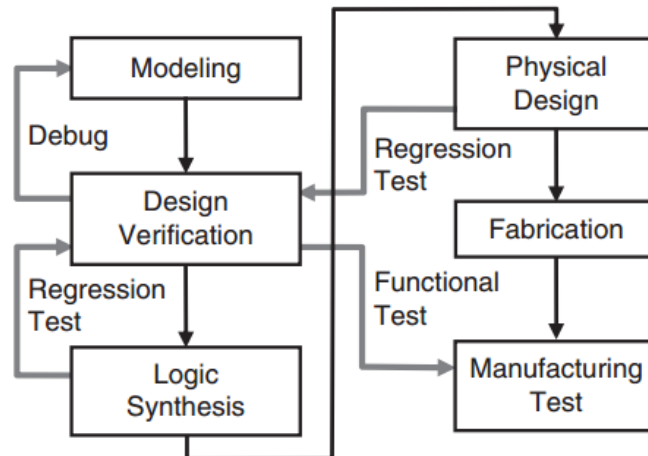


Figure 5-1 - Typical EDA flow [51]

Once the design is verified, the Logical Synthesis is performed with the objective of mapping the behavioral description into a logic description of the circuit using cell libraries. The output of the Logical Synthesis is a netlist that contains a list of standard cell instances and nodes representing interconnectivities, as seen in figure 5-2. Note that this is an important step in the synthesis process since it is responsible for identifying how the logic will be mapped into actual cell libraries. It's also worth noticing that this step will influence the circuit-level soft-error cross-section, since it defines the actual

cells used from a library. The evaluation of soft errors in the Logical Synthesis was the objective of another thesis [52], where important constraints and hardening techniques applicable to the Logical Synthesis were discussed. As this thesis focuses on physical synthesis and effects of physical synthesis on soft-error cross-sections, the Logical Synthesis step is considered, beyond the scope of this work. The input of the Physical Synthesis is the output of logic synthesis step in the form of a netlist description, as shown in figure 5-2, along with the standard cell library geometry information.

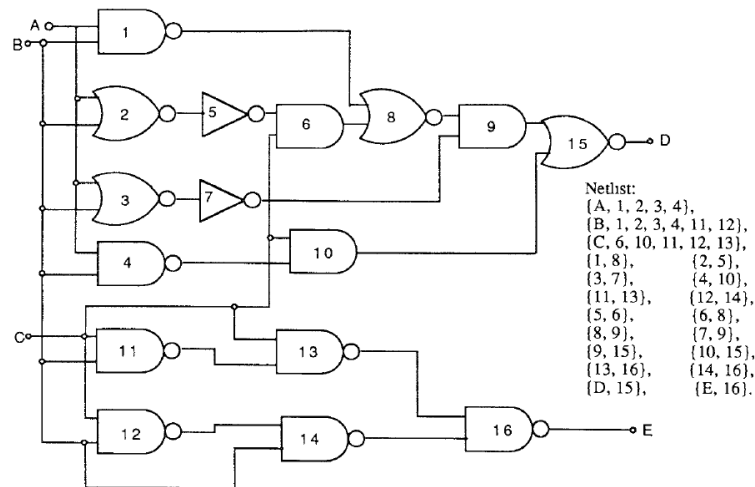


Figure 5-2 - Netlist and its circuit representation [53]

For the netlist shown in figure 5-2, standard cell instances are indicated with numbers and letters for Primary Inputs (PI) and Primary Outputs (PO). Interconnectivities are generated for each instance set by analyzing the resulting graph. Once the netlist is ready, the Physical Design\Synthesis process uses this information to generate a physical representation of the design that is placed, routed and electrically checked for correctness. The output of the Physical Synthesis can be seen in the diagram of Figure 5-3.

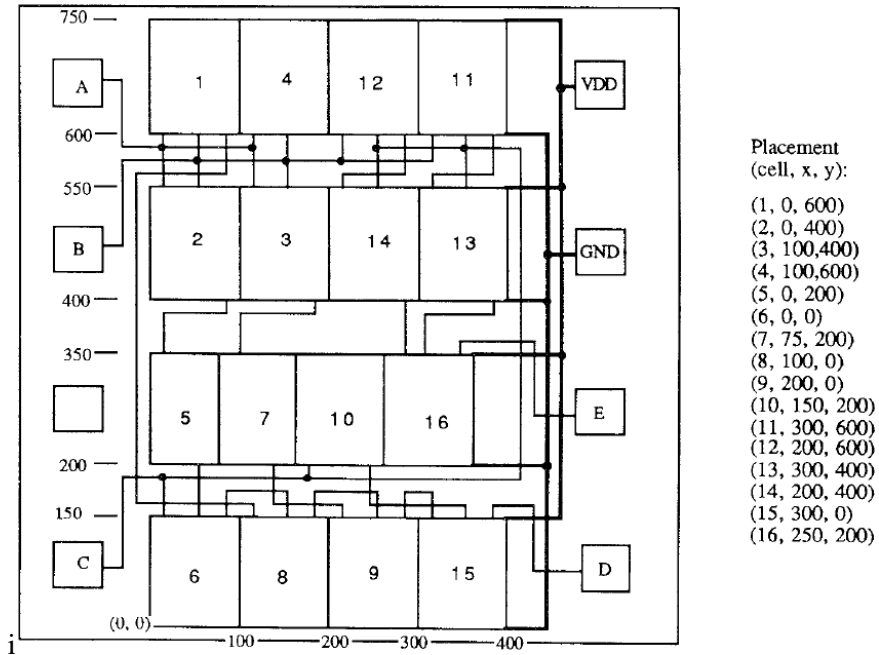


Figure 5-3 - Physical Synthesis output [53]

Once the Physical Synthesis is completed, the design is verified and masks are sent for production. With the physical IC manufactured, EDA tools are again used for manufacturing test and verification, before the foundry will send the ICs to their clients. As each of these steps have specific goals and address different problems, this work does not further discuss other synthesis steps and focus only on Physical Synthesis step.

5.1 Physical Synthesis Flow

The main objective of the Physical Synthesis process is to guarantee that the manufactured IC is electrically correct according to constraints provided by the ASIC designer. The Timing Closure is a term used by the community to refer to the timing constraints of the design that have to be matched for the correct circuit operation. The physical synthesis process can be sub-divided in steps shown in figure 5-4.

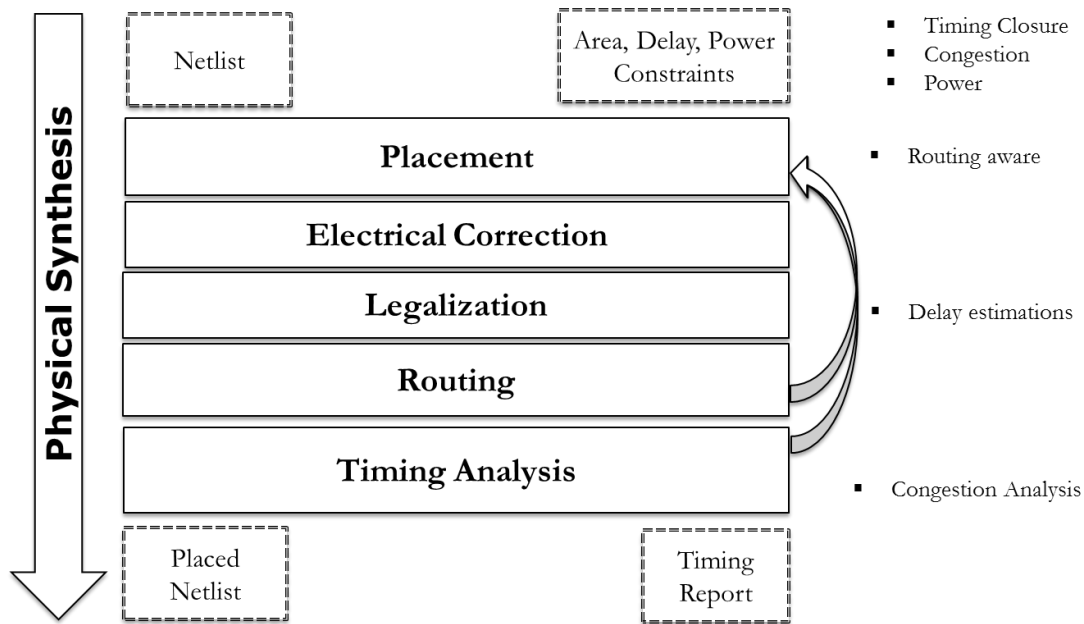


Figure 5-4 - Physical Synthesis based on a commercial design flow [54]

The placement is responsible for defining the position of each standard cell and input/output PADS in the provided IC area. The electrical correction and legalization steps are responsible for adjusting the netlist and the position of standard cells, with the objective of reaching the timing closure with the constraints provided by the designer [7]. The Electrical Correction step use techniques like Gate Sizing, Gate Cloning, Gate Relocation and Buffering in attempt to reach the timing closure [55]. The legalization step guarantees that there are no physical overlaps between the standard cells, thus not occupying the same region on an IC.

Once the placement is considered legalized, the Routing is performed to create the interconnectivities of the design. This step has less control over the design and is strongly influenced by the placement process. Due to this reason, routing has little influence on soft-error cross-section. The only effect routing will have on soft-error cross-section will be from parasitic interconnect capacitances. Usually, the only important information related to soft errors (routing capacitance) can be easily estimated using wire lengths. Wire length estimations are also needed for timing analysis. For this project, interconnect capacitances and delay are estimated using wire length models. Once a design is routed, a timing analysis is performed to verify if the timing closure is reached with the constraints provided by the user. If the

design does not reach the timing closure, the design has to be re-placed, re-routed, and passed through the electrical correction step again.

The electrical correction techniques do influence the electrical characteristics of the circuit, and thus will influence the soft error response of the IC. It's important to understand the techniques used by EDA to solve the circuit delay problems and how they will impact the circuit soft error response. A review of the main electrical correction techniques applied in the Physical Synthesis is discussed next.

5.2 Electrical Correction Techniques and Soft Errors

The main electrical correction techniques used in the Physical Synthesis are the Gate Sizing, Gate Cloning, Gate Relocation, and Buffering as discussed by [55]. These techniques are commonly used by the synthesis to achieve the timing closure after the initial netlist placement is available.

The Gate Sizing technique [56] consists of replacing standard cells that are not able to drive the expected load (higher load capacitances than expected), with a cell with identical function but higher current driving capability. This is shown in figure 5-5. In a standard cell library, higher driving capability is achieved through adding transistors in parallel to the ones already present in the cell. This operation, increases the width of the standard cell, but not the cell height. The addition of transistors in parallel reduces the total resistance from the output to the power rail, thus increasing the cell drive current ($I=V/R$). The usage of Gate Sizing as a technique to mitigate Soft Errors was proposed by [57] with an algorithm to compute the necessary gate size to mitigate SETs modeled by the double exponential equation [31]. The increased driving-current capability at the output node reduces the SET pulse width at that node, thereby reducing the circuit soft-error vulnerability. The effectiveness of transistor sizing to mitigate soft errors was further evaluated by [58] using a TCAD tool for a 90-nm technology.

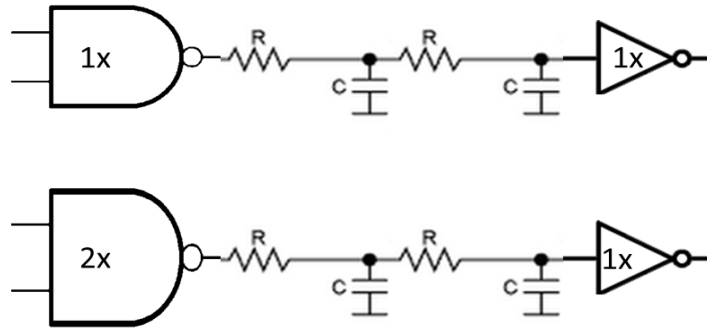


Figure 5-5 - Gate Sizing technique

The Gate Cloning [59] technique consists of duplicating the standard cell and splits the connections between the cell and its duplication. This in turns reduces the load capacitances associated with the cloned cell as shown in figure 5-6. The operation is usually performed in cells with high fan-out and long interconnections associated with its output. In figure 5-6 on the left shows a NAND2 cell driving 4 inverters and its interconnect parasitic RC network. The gate cloning technique duplicates the cell, keeping the inputs shorted and splits the load capacitances, so now each cell drives only 2 inverters, as shown on the right side of the figure.

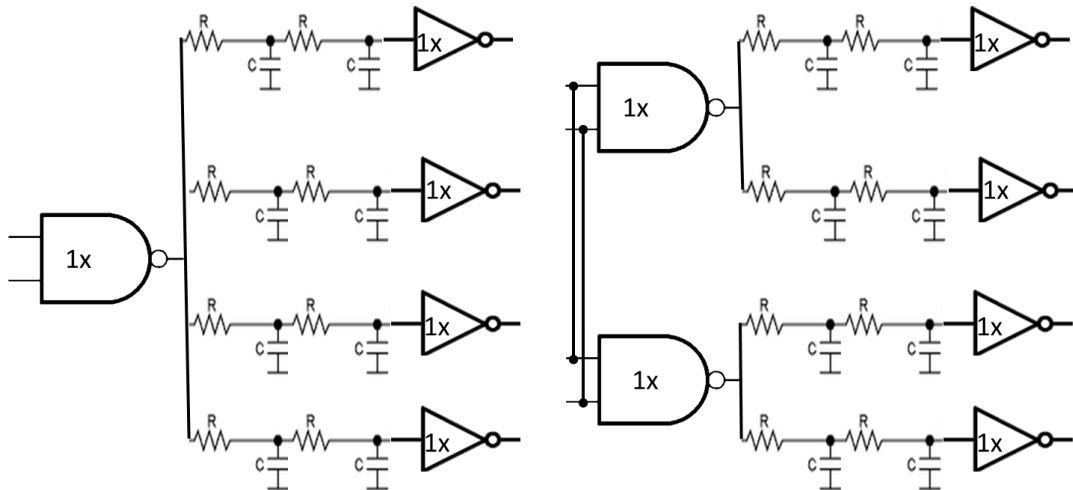


Figure 5-6 - Gate Cloning technique

The Gate Cloning has also been proposed for Soft Error mitigation as discussed by [60]. The standard cell is duplicated as shown in figure 5-7. In this case, the cell is cloned but the output is not split, so now

two cells drive the same node. This essentially increases the driving current at the output node, reducing the soft-error vulnerability.

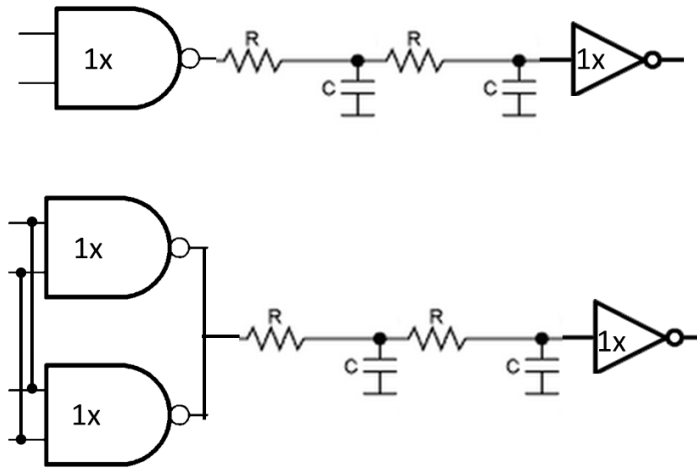


Figure 5-7 Gate Cloning for Hardening

Buffering [61] is another technique used to enable cells to drive big loads, as shown in figure 5-8. By adding an even number of inverters with increasing size, it's possible to make the design drive long interconnectivity lines. The number of buffers added to drive the interconnectivity delay will depend on the strategy used by the EDA tool. Due to the poor scaling of the interconnectivity delay when compared with standard cells, the number of buffers required in modern circuits grows fast. Care must be taken to ensure that these additional gates are taken into consideration for soft-error cross-section estimation.

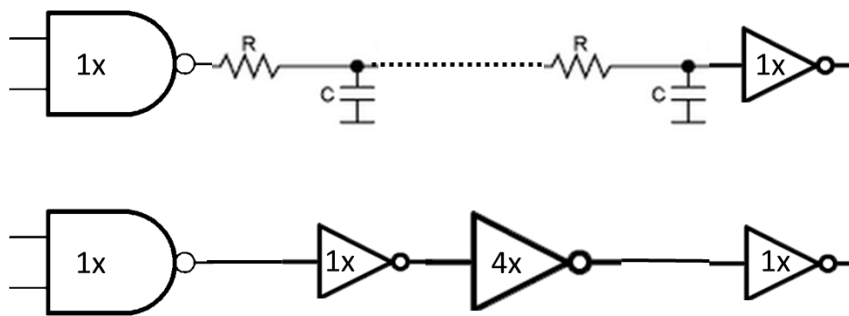


Figure 5-8 - Buffering the interconnectivity

Lastly, the Gate Relocation [62] works by swapping cells with its neighbors to achieve timing closure and is shown in figure 5-9. In the figure, only the interconnections of the NOR2 and INV cells involved in the delay problem are shown, with color blue indicating an acceptable delay and red a timing violation.

On the left side of the figure, the initial cell configuration, with a timing violation in the inverter output shown by a red interconnect line. This violation is caused by the long wire length needed to make the connection. This problem is fixed by shifting the inverter cell closer to the NOR gate. The relocation technique works by searching for a new cell position that solves the timing problem of the cell, and not creating additional timing violations. This technique is used by the detailed placement algorithms to perform local optimizations after the global placement is completed.

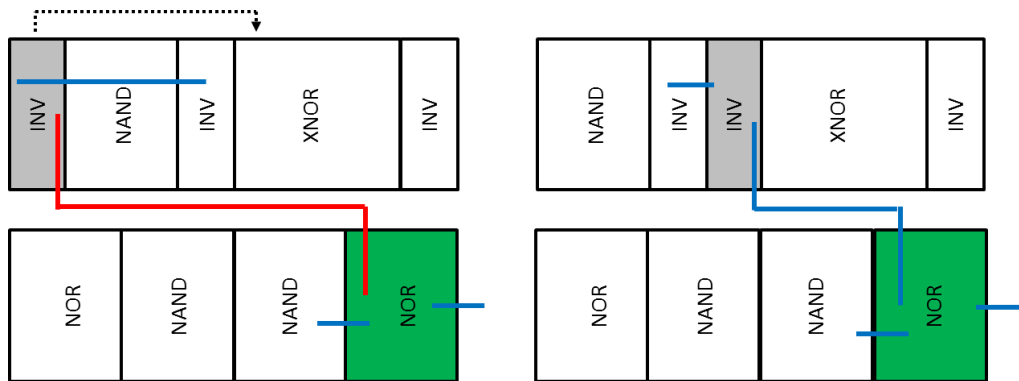


Figure 5-9 Gate Relocation (Cell swap)

It's the objective of this work to evaluate how the techniques Gate Sizing, Buffering and Gate Cloning influence the circuit soft-error cross-section. In the next section, the implementation of the automatic placement is discussed.

5.3 Automatic Placement

Now that techniques used to achieve the Timing Closure were discussed, focus is taken to the core step of the Physical Synthesis, the Automatic Placement. The placement requires the netlist generated by the Logical Synthesis and the geometry information from the standard cell containing both the dimensions of the cell and internal pin locations. The placement, then seeks to map these cells to physical positions in an IC area by minimizing constraints, like area and power. It turns out that this problem is solved by formulating this as an optimization problem, solved using a wide range of techniques, from

Monte Carlo methods to Genetic and Analytical algorithms as discussed by [53]. Some placement optimization algorithms are listed in table 5-1.

Table 5-1 – Comparison between placement algorithms [39]

Algorithm	Result quality	Speed
Simulated Annealing	Near optimal	Very slow
Genetic algorithm	Near optimal	Very slow
Force Directed	Medium ~ good	Slow ~ medium
Numerical optimization	Medium ~ good	Slow ~ medium
Quadratic Programming	Medium ~ good	Fast
Min-cut	Good	Medium
Clustering	Poor	Fast

From the techniques described in table 5-1, the focus of the EDA industry is currently between both Force Directed (FD) and Quadratic Programming (QP) based algorithms [63]. It is important to acknowledge that this is a very competitive field that is constantly changing with new heuristics and techniques constantly emerging. It is worth noticing that the placement flow might be divided into two steps. First a Global Placement is performed, where the optimization seeks the global position for the cells in the IC area (not necessarily with minimum wire length). At this moment the relative position between each cell is more important. Second a Detailed Placement uses the output of the first step to locally optimize the cell positions to minimize the wire length. Note that this sub-division in steps is implementation dependent. For this project, two different placement techniques are used to evaluate their performance in the Physical Synthesis. These algorithms may be replaced by new algorithms as needed and as available. The algorithms used are Simulated Annealing [64] based placement and a Quadratic Placement algorithm developed specifically for this project.

The placement problem is now mathematically formulated. The netlist provided by the Logical Synthesis is modeled as a graph of vertexes V and edges E . Standard cells are represented by vertexes and interconnectivities by edges. The cell position coordinates x, y are denoted by V_x, V_y and the pin positions are given by V_x^{pin} and V_y^{pin} . The pin position is given by the cell position and the pin offset extracted from the cell geometry, and represented as $V_x^{pin} = V_x + pin_offset$. The edge cost function is

implementation dependent, but is usually given by the Half-Perimeter Wire Length (HPWL) distance, calculated with Eqn. 5.3. The placement algorithm seeks the set V that minimizes the cost function, given by Eqn. 5.4. HPWL for each edge in the set E is used to estimate these functions.

$$G = \{V, E\} \quad (5.1)$$

$$E = \{E_0, \dots, E_K\}, e = \{V_i, V_j\}, V = \{V_0, \dots, V_N\} \quad (5.2)$$

$$HPWL_k(V_i, V_j) = (V_i^{pin_x} - V_j^{pin_x})^2 + (V_i^{pin_y} - V_j^{pin_y})^2 \quad (5.3)$$

$$HPWL_N = \sum_{k=0}^N HPWL_k(V_i, V_j) \quad (5.4)$$

For the SA placement to be discussed in the next section, the objective function that evaluates the cost of the edge, is given by Eqn. 5.3 and 5.4. The Eqn. 5.3 can also be referred to as Manhattan distance [65]. The HPWL is a good predictor for delay analysis, and is also correlated with congestion and several other circuit hazards related to routing. The placement algorithm implemented in this work seeks the timing closure by minimizing HPWL. The next sections describe the algorithm developed to solve the placement problem. These solutions were coded using C++11.

5.4 Simulated Annealing Placement

The SA, also known as Monte Carlo Simulated Annealing, is classified as a combinatorial heuristics and it was proposed for the placement solution by [66]. With proper setup, the algorithm is proven to reach near optimum solutions, but with an extremely high computational cost [67]. The heuristics simulate the heating and annealing process, and its algorithms pseudocode is shown in figure 5-10, based on the implementation from [51].

```

PROCEDURE simulated annealing;
  initialize; (temperature\loopcount\trials)
  configuration=generaterandomconfig(configuration);
  while (loopcount<maxloop && temperature>0){
    while (trials<maxtrials){
      new_configuration=perturb(configuration);
      delta=evaluate(new_configuration,configuration);
      if (delta<0) configuration=new_configuration;
      else if (accept(delta,tempeature)>random(0,1))
configuration=new_configuration;
    }
    --temperature; ++loop_count;
  }
}

```

Figure 5-10 - Simulated Annealing Pseudocode [66]

In the implementation shown in figure 5-10, the temperature is defined as an integer to control the acceptance probability of configurations during the placement. The algorithm starts with a very high temperature that is gradually reduced to not allow bad configurations be accepted by the heuristic. By applying random displacements in the set V , the cell positions (V_x, V_y) are changed randomly. This is done by the *perturb()* function. At the end of every iteration of the internal loop, the algorithm verifies the cost of the configuration, using the *evaluate()* function. The evaluate function uses the HPWL estimator (5.4) to compute the cost of the configuration and returns *delta*, the difference between the new configuration and the previous configuration. If the cost is reduced, the perturbation is accepted. If not, a second verification is performed. In case of HPWL increase, the algorithm is allowed to accept the change if the *accept()* functions returns a number higher than a randomly generated number (this random feature allows SA jumps through local minima's). When the system is on high temperature, the acceptance probability of bad moves is very high, but with the temperature decrease, the probability of the loop accepting a bad perturbation decreases.

$$accept(delta, temperature) = \exp\left(-\frac{delta}{temperature}\right) \quad (5.6)$$

The acceptance function is shown in Eqn. 5.6 [66] and it was used to provide a smooth reduction in the acceptance probability of bad perturbations with the temperature decrease. Once the temperature reaches

zero, the placement is legalized and accepted as the final solution. The SA placement implemented in this work is classified as a detailed placement algorithm. In the next section the QP placement is discussed.

5.5 Quadratic Placement

To replace the SA, due to its slow execution performance especially with big netlists, a placement based on the state-of-the-art algorithm used by the EDA industry was developed. The Quadratic Programming falls into the category of Analytical Placements and has emerged in the last few decades to become one of the key methods to solve the placement optimization problem. In modern Physical Synthesis tools, the placement problem is usually divided into two steps, the global placement and the detailed placement. The developed placement flow is shown in figure 5-11, with the global placement in the left, followed by the detailed placement on the right.

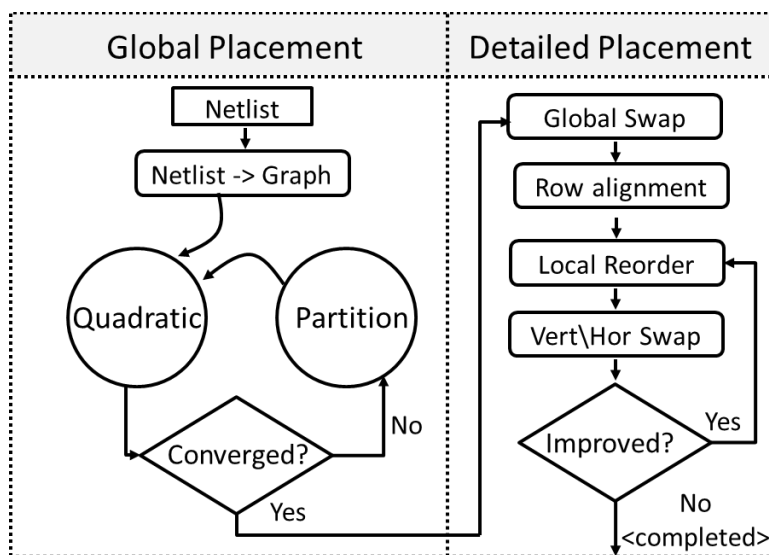


Figure 5-11 Developed placement flow

Once the Detailed Placement is completed, the electrical correction step is used to achieve timing closure. After that, a legalization is performed to ensure the resulting placement is legalized (does not have cell overlaps). The global and detailed placements are discussed next.

5.5.1 Global Placement

The global placement, as shown in figure 5-11, uses Quadratic Programming to solve the optimization problem discussed in section 5.3. In this work the Quadratic Placement uses the region partition algorithm to reduce the overlap between cells and spread of cells in the placement area [68]. In the QP implemented in this work, the length of the netlist is given by the quadratic cost function (shown in Eqn. 5.7) where the $w_{x,pq}$ and $w_{y,pq}$ are the Bound2Bound weight factor [69] used to take into account pin density when estimating the optimum wire length, and is given by Eqn. 5.8. The equations ahead are based on the implementation of both [68] and [69].

$$\Gamma = \frac{1}{2} \sum_{e=(V_i, V_j) \in E}^K w_{x,ij} (V_i^{pin} + V_j^{pin})^2 + w_{y,ij} (V_i^{pin} + V_j^{pin})^2 \quad (5.7)$$

$$w_{x,ij} = \begin{cases} 0, & \text{if } i \text{ and } j \text{ are inner pins} \\ \frac{2}{p-1} \cdot \frac{1}{|V_i^{pin} - V_j^{pin}|}, & \text{else} \end{cases} \quad (5.8)$$

To minimize the quadratic function Γ , Eqn. (5.7) is transformed in a matrix vector notation as shown in Eqn. 5.10, where \mathbf{x} and \mathbf{y} are the vector position of the M movable blocks (Eqn. 5.9), and vectors \mathbf{d}_x and \mathbf{d}_y are used to reflect the connections between fixed modules and movable blocks [68]. The matrices \mathbf{C}_x and \mathbf{C}_y are the connectivity matrices of dimension $M \times M$ and represent the connectivity between movable blocks. The generation of these matrixes is performed using the method proposed by [69].

$$\mathbf{x} = (Vx_0, \dots, Vx_M)^T \quad \mathbf{y} = (Vy_0, \dots, Vy_M)^T \quad (5.9)$$

$$\Gamma = \frac{1}{2} \mathbf{x}^T \mathbf{C}_x \mathbf{x} + \mathbf{x}^T \mathbf{d}_x + \frac{1}{2} \mathbf{y}^T \mathbf{C}_y \mathbf{y} + \mathbf{y}^T \mathbf{d}_y \quad (5.10)$$

The properties of the quadratic function Γ will depend on the characteristics of the connectivity matrices, which are built in such a way so that Γ is a positive semidefinite matrix if there are no modules fixed in \mathbf{C}_x and \mathbf{C}_y , or Γ is positive definite if some modules in the connectivity matrices are fixed. In both cases Γ is convex and the minimum can be obtained by setting its first derivative to zero. Notice also that Γ is separable, resulting in the quadratic functions (given by Eqn. 5.11 and 5.12) for the x and y coordinates respectively.

$$\Gamma_x = \frac{1}{2} \mathbf{x}^T \mathbf{C}_x \mathbf{x} + \mathbf{x}^T \mathbf{d}_x \quad (5.11)$$

$$\Gamma_y = \frac{1}{2} \mathbf{y}^T \mathbf{C}_y \mathbf{y} + \mathbf{y}^T \mathbf{d}_y \quad (5.12)$$

The system of linear equations (Eqn. 5.11) is represented by

$$\begin{pmatrix} \vdots \\ \Gamma_{x,i} \\ \vdots \\ \Gamma_{x,j} \\ \vdots \end{pmatrix} = \frac{1}{2} \begin{pmatrix} \cdots & x_i & \cdots & x_j & \cdots \end{pmatrix} \begin{pmatrix} \vdots & \vdots & \vdots & \vdots & \vdots \\ \cdots & C_{x,ii} & \cdots & C_{x,ij} & \cdots \\ \vdots & \vdots & \ddots & \vdots & \vdots \\ \cdots & C_{x,ji} & \cdots & C_{x,jj} & \cdots \\ \vdots & \vdots & \vdots & \vdots & \vdots \end{pmatrix} \begin{pmatrix} \vdots \\ x_i \\ \vdots \\ x_j \\ \vdots \end{pmatrix} + \begin{pmatrix} \cdots & x_i & \cdots & x_j & \cdots \end{pmatrix} \begin{pmatrix} \vdots \\ d_{x,i} \\ \vdots \\ d_{x,j} \\ \vdots \end{pmatrix}$$

The solution of the quadratic function is also subjected to a constraint used to restrict the freedom of movable blocks around the placement area. The partition step showed in figure 5-11 is used to split the placement area to both reduce the cell overlap and to distribute the cells during the optimization. The constraint utilized in this placement was based on the region partition proposed by [68], and is given by Eqn. 5.13. In the equation, the matrix \mathbf{A}^l indicates the module position in the region partition l and have its entry calculated using equation (Eqn. 5.14). The vector \mathbf{u} holds the center of the respective region l .

$$\mathbf{A}^{(l)} \mathbf{x} = \mathbf{u}^{(l)} \quad (5.13)$$

$$a_{pu} = \begin{cases} \frac{F_u}{\sum_u F_y}, & \text{if } u \in p \\ 0, & \text{otherwise} \end{cases} \quad (5.14)$$

In Eqn. 5.14, F_u is the total unit area occupied by module u and p is the region. The constraint, given by Eqn. 5.13, has the effect of restricting the modules in the matrix $\mathbf{A}^{(l)}$ to the respective regions indicated by the vector $\mathbf{u}^{(l)}$. The central coordinates of each region partition act like a gravity pulling modules away from the initial configuration, reducing the amount of overlap between the cells. The matrix $\mathbf{A}^{(l)}$ has M columns and L lines for each region partition. As a consequence to the fact that a module can't be in more than one region, each column in the matrix contains only one nonzero entry, as indicated by Eqn. 5.14. With both constraint and objective function stated, the minimization problem can be formulated as Eqn. 5.15.

$$\min_{\mathbf{x} \in M} = \left\{ \frac{1}{2} \mathbf{x}^T \mathbf{C}_x \mathbf{x} + \mathbf{x}^T \mathbf{d}_x \mid \mathbf{A}^{(l)} \mathbf{x} = \mathbf{u}^{(l)} \right\} \quad (5.15)$$

As discussed before, the Γ is positive-definite and convex. To solve Eqn. 5.15, the library QuadProg++ [70] is used in this work. The library solves the QP by implementing the *Goldfarb-Idnani active-set dual* method presented by [71] to solve strictly convex problems. With the resulting cell configuration, the placement evaluates the placement convergence seeking high density overlaps. If there are still areas in the placement with high density cell overlaps, the region is further partitioned. In every high overlapping area, the algorithm makes a new region partition and call QP iteratively. The addition of a new partition increases the size of $\mathbf{A}^{(l)}$ by one. Once the convergence is reached, the detailed placement process is started.

5.5.2 Detailed Placement

Once the Global Placement is completed, an optimization step is performed to fix bad cell positions generated by the region-partition algorithm discussed in the last section. During the global placement, when a region partition is defined, cells are pulled to the region central location and will not be able to return to the previous partition. In some cases, groups of connected cells are split in different regions when they shouldn't be split. To achieve a better cell organization, the Global Swap step is used

in the detailed placement to move cells freely in the placement area. During the detailed placement process, the region partition is ignored.

The Global Swap is performed once, by iterating over the M movable cells seeking for the optimum cell position (x_{opt}, y_{opt}) given the current placement condition. The estimated optimum cell position [72] is given by the median of the x and y coordinates of each cell neighbors (given by Eqn. 5.15-5.16).

$$x_{opt} = \text{median}(x_0, x_1, \dots, x_h) \quad (5.15)$$

$$y_{opt} = \text{median}(y_0, y_1, \dots, y_h) \quad (5.16)$$

After the x_{opt}, y_{opt} are calculated, the HPWL method is used to determine if there is an improvement over the wire length. If the netlist is improved, the cell position is updated and the algorithm continues until all cells are verified. With the Global Swap completed, the Row Alignment is started.

The resulting placement now has cells spread around the placement area and overlaps between cells exist. To remove the overlap and put the cells in a row organization (also adding routing channel), the algorithm uses the method proposed by [68] to align the modules without perturbing the relative organization between the cells. The Row Alignment step consists of first ordering the cells by the y coordination. Once the order along the y -direction is completed, an estimation is calculated to determine the number of cells per row to define sub-sets to be placed in each row. Once this sub-set is defined, the cells are now ordered by their x -coordinates and placed in their positions within each row. In this thesis, the cell ordering is performed using the Merge Sort algorithm [73], with worst case performance of $O(n \log n)$. At the end of this step, the cells are organized in rows separated by routing channels (with channel height defined by the user).

With the cells organized in rows, a local optimization step is performed to relocate bad placement of individual cells and to improve the netlist wire length. The algorithm starts a loop that stops when no

improvement is obtained to the netlist after the Local Reorder and Vertical\Horizontal Swaps are performed This step in the detailed placement is based on optimizations proposed by [72, 74]

The Local Reorder [74], shown in figure 5-12, works by iterating over each row and assigning each cell in the row to a segment. The size of the segment (how many cells are included in each segment) is user defined and may affect the computation complexity and performance. The Local Reorder algorithm permutes the relative position of cells within each segment. If the new permutation improves the HPWL, the new cell ordering is accepted and the algorithm proceeds to the next segment of cells until all segments are reordered.

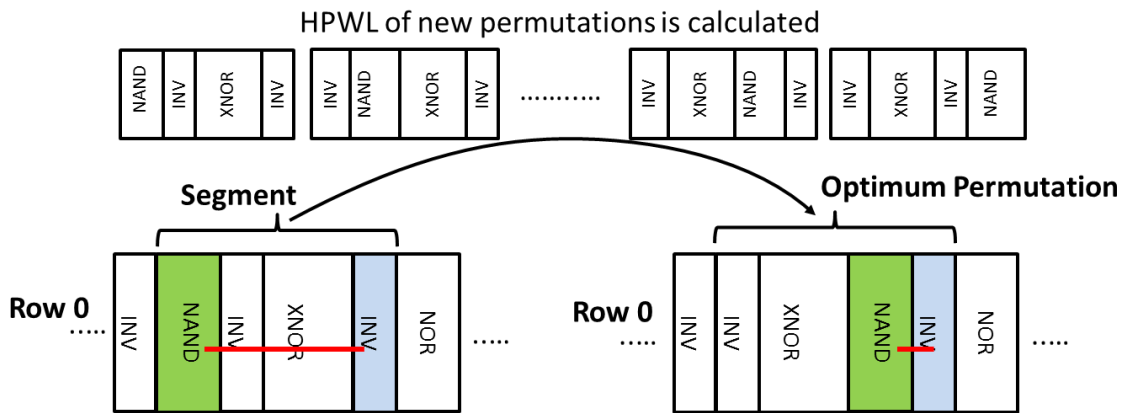


Figure 5-12 Placement Local Reordering

With the Local Reorder completed, the Vertical and Horizontal Swap algorithms [74] are used to replace individual cells as shown in figure 5-13 and 5-14 respectively. The Vertical Swap, shown in figure 5-13, exchanges cells placed in different rows. To identify the direction of the swap (bottom/top row), the y_{opt} is calculated for the cell being evaluated and the target cell to be swapped. If the swap improves both HPWL, the operation is approved and cells change the assigned rows.

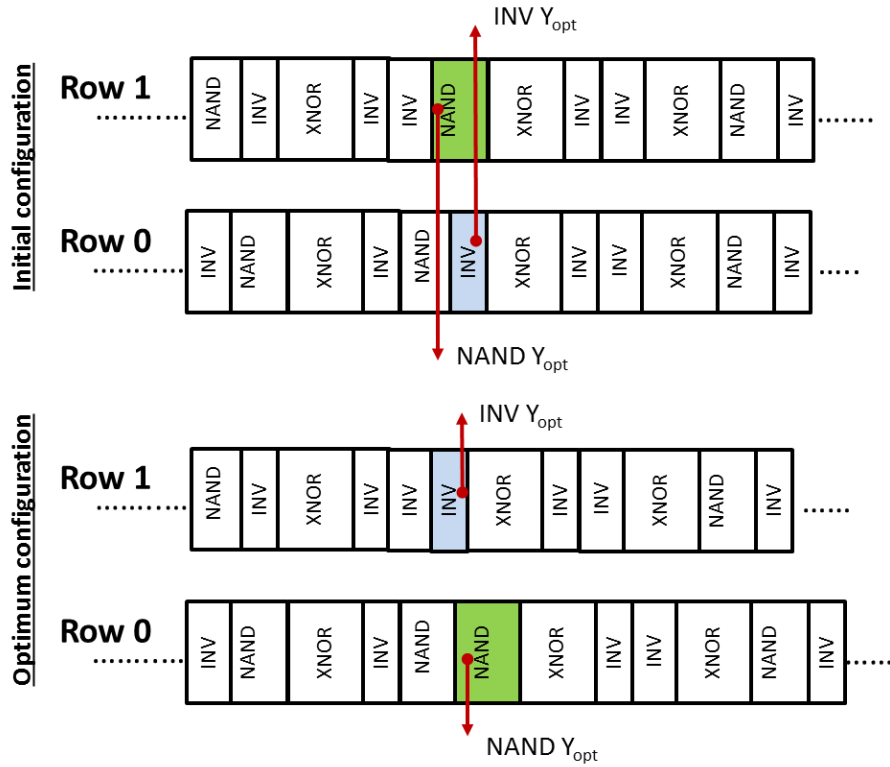


Figure 5-13 Vertical cell swap

The Horizontal Swap algorithm, showed in figure 5-14, works the same way along the horizontal direction by sliding the cell within the respective row seeking HPWL improvement. The cell stops moving along the row when the HPWL value does not improve.

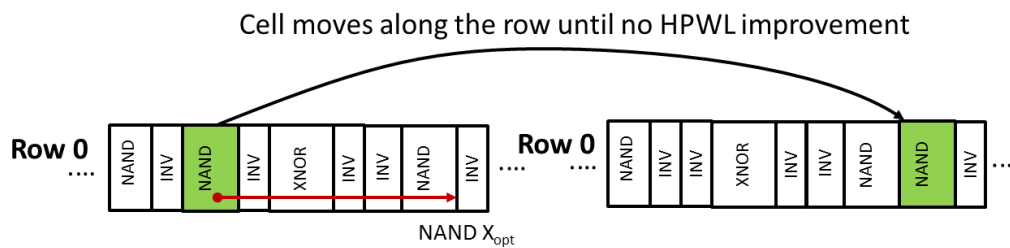


Figure 5-14 Horizontal Cell Swap

The result of the last 3 optimization algorithms is evaluated and compared with previous optimization iterations. Once the iteration does not improve the HPWL, the placed netlist is considered optimum and the Time Analysis is started.

5.6 Time Analysis

The calculation of the Electrical De-rating and the evaluation of the timing closure for the electrical correction require the estimation of the time delay associated with both the standard cell and the interconnectivity. To estimate the standard cell delay, first the output load capacitance has to be estimated; a task that can be performed once the interconnectivity RC characteristics is estimated. Once this load capacitance is known, a look-up table that models the cell delay, taking into account its input and output conditions, is used to estimate the standard cell delay. This information is available inside the Liberty Library file [75] contained within the Standard Cell Library.

The interconnectivity wire length is estimated using the Manhattan distance, a metric considered to be accurate for the wire length estimation. Given the wire length, the interconnectivity delay is calculated using the Elmore Delay [76] technique shown in Fig. 5-15. The Elmore delay has been used by the EDA industry to estimate the wire length delay for a few decades due to its simplicity and low computation requirements [77]. In the figure, the interconnectivity is modeled by a network of Resistances (R) and Capacitances (C) estimated from the wire characteristics and a load capacitance (C_l) from the next logic stage.

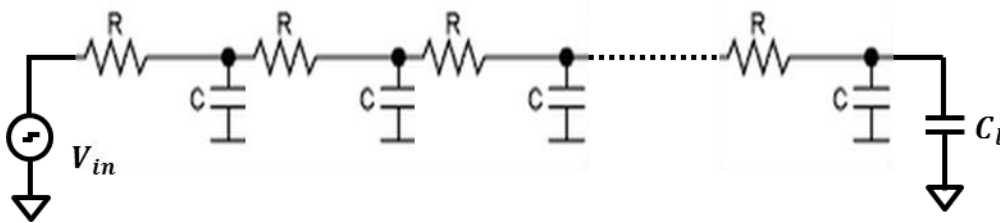


Figure 5-15 RC delay network used to model the interconnectivity [76]

The calculation of the Elmore delay is given by Eqn. 5.17 and is the summation of the path resistances (R) multiplied by each delay segment load (L) and the gate load capacitance of the connected standard cell.

$$T_D = \sum_{i=1}^N C_i \sum_{j=1}^i R_j \quad (5.17)$$

The interconnection resistance can be calculated using Eqn. 5.18 where ρ is the metal resistivity, L is the metal length, t and w are metal thickness and width, respectively.

$$R = \rho \left(\frac{L}{t * w} \right) \quad (5.18)$$

The interconnect capacitance can be split into a line-to-ground capacitance, line-to-line capacitance and crossover capacitance. The analysis of the line-to-line capacitance and crossover capacitance require the routing information, since the relative position between each metal line is required for the computation. Due to this reason, this work limits the interconnection capacitance to the computation of the Line-to-Ground capacitance [78] only as shown in Fig. 5-16.

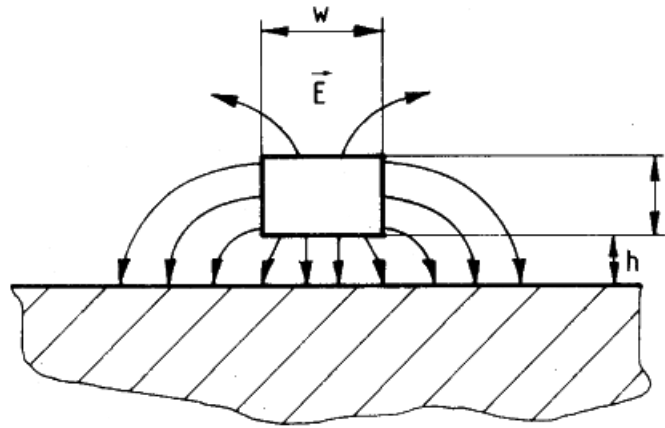


Figure 5-16 Line-to-Ground capacitance – Interconnection line facing the substrate [78]

The line-to-ground capacitance (C) is given by Eqn. 5.19, where w is the wire length width, h is the height between the metal layer and the substrate, and t is the metal thickness. The constant ϵ is the relative permittivity given by $\epsilon = \epsilon_0 \cdot k$, where the vacuum permittivity $\epsilon_0 = 8.85 \times 10^{-14} F/cm$ and k is the dielectric constant, for $SiO_2 = 3.9$.

$$C = \varepsilon \left[\frac{w}{h} + 0.77 + 1.06 \left(\frac{w}{h} \right)^{0.25} + 1.06 \left(\frac{t}{h} \right)^{0.5} \right] \quad (5.19)$$

This model was validated by comparing the delay estimations with electrical simulation results using *Cadence Virtuoso* [44]. In Fig. 5-17, a section of the network used to analyze the network delay is shown. In the figure, 3 different interconnects are modeled between two inverters from the 40-nm technology node. Each *RC* component in the figure was calculated to model a wire segment of $1\mu\text{m}$ of length, based on data from [77].

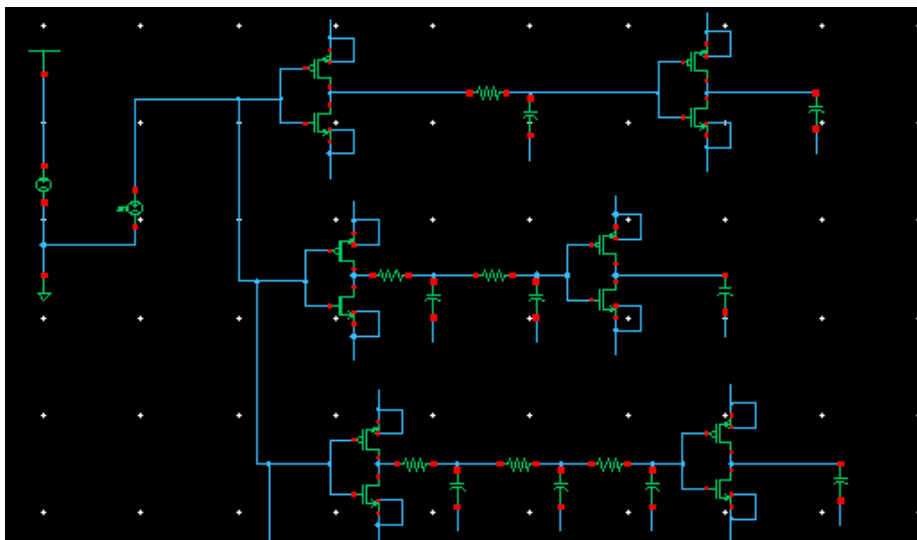


Figure 5-17 Snapshot of part of the RC network used in simulations

The simulation results along with the Elmore delay estimations are shown in Fig. 5-18. These results indicate that the model does serve as a very good first-order estimation for the delay. It must be noted here that the error in Elmore delay increases significantly as wire length increases. Care must be taken to estimate delays using more accurate methods when wire length exceeds a certain user-defined threshold (process dependent).

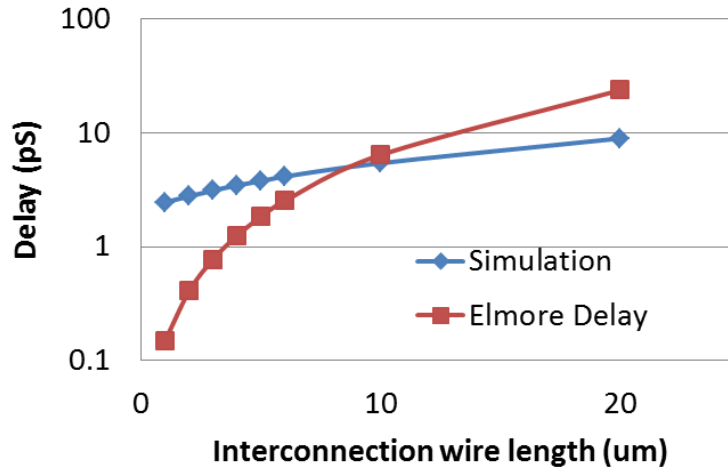


Figure 5-18 Elmore model vs simulation

This data is used carefully in this thesis since it has a direct impact in both the Electrical Correction step and the Electrical De-rating estimation. Note that the Manhattan Distance is used to estimate the wire length and the RC network used to estimate the wire delay assuming a single wire section. The final circuit interconnection is better represented as a spanning tree, rather than a single interconnection for improved accuracy. Usually, this would lead to additional RC elements in the calculations. Thus, the algorithms and the constraints discussed above do yield a lower bound on the RC delay for the network. Also, note that the capacitance model only considers line-to-ground capacitance, not taking account line-to-line and crossover capacitances. Adding these capacitances will increase the overall load capacitance for a given node. It is not possible to estimate these capacitances without performing the detailed routing.

These two characteristics result in a lower bound of the estimation, a desired property for this model. The interest is in the lower bound due to Electrical De-rating calculation. As discussed in chapter 4, the RC delay is able to filter short SET's, and by using the lower bound estimation one can guarantee that SET pulses masked by the methodology would not be able to propagate to the output. The impact of this under-estimation on the electrical correction is to increase the number corrections performed by the

model. However, this does not affect the trend to be observed in the analysis of the SER cross-section analysis.

5.7 Electrical Correction & Legalization

With the Time Analysis completed, the Electrical Correction techniques can be applied to solve netlist violations. The placed netlist is now analyzed to verify if the Timing Closure is achieved. In this work, the Timing Closure is formulated using Eqn. 5.20 and 5.21 from [79]. The Eqn. 5.20 models the Setup Constraint, and refers to long paths in the netlist. It specifies the amount of time the input data should be steady before the clock edge of each Flip Flop. In the equation T_{cycle} is the clock period, $t_{critical_path}$ is the longest path in the circuit, t_{setup} the clock setup time and t_{skew} is the clock skew.

$$T_{cycle} \geq t_{critical_path} + t_{setup} + t_{skew} \quad (5.20)$$

The Eqn. 5.21 models the Hold Constraint and refers to short paths used to indicate the amount of time the input data should be stable after the clock edge. In the Eqn. 5-21, t_{comb_delay} is the path delay from the cell to FF, and t_{hold} is the Flip Flop hold time.

$$t_{comb_delay} \geq t_{hold} + t_{skew} \quad (5.21)$$

The Time Analysis section discussed before is used to provide the timing information required in this section. The worst and the best case delay paths associated with each node in the graph is stored to evaluate conditions (given by Eqn. 5.20 and 5.21. If time closure is not achieved (equalities specified in Eqn. 5-20 and 5-21 are not satisfied), the electrical correction techniques, such as Gate Sizing, Gate Cloning and Buffering are used to adjust the netlist timing. Because we are also interested in analyzing the effects of each of these individual techniques, algorithms used in this work allow the user to block the use of each of these electrical correction techniques. Notice also that the input frequency provided by the user will have a strong influence on the Timing Closure and will directly impact the number of operations to be performed in this step.

In this work, the Gate Sizing was applied in the critical paths where there is a violation for Eqn. 5.20. The algorithm seeks the smaller cell size to fix the delay constraint. First, the magnitude of the time violation is calculated by measuring the difference between the $t_{critical_path} + t_{setup} + t_{skew}$ and T_{cycle} . Once the time violation is measured, the time improvement per cell in the critical path is estimated dividing the time violation by the number of cells in the critical path. The algorithm iterates over the critical path, sizing cells according to the necessary cell delay improvement for the given critical path.

The Gate Clone is applied by analyzing both the cell fan-out and wire length delay. If the cell has a high fan-out (>4) and high load capacitance, beyond the acceptable threshold indicated by the standard cell library, the cell is cloned as described in section 5.2. The addition of Buffers is done by analyzing the wire length delay, previously calculated by the Timing Analysis. If the load of the interconnection is beyond the acceptable cell load, and the wire length is the source of the high load, the interconnection is split in half, using Buffers. In some situations, like a high load occasioned due to a small cell driving a big cell (for example, INV_1X driving INV_32X), the choice for buffering and gate sizing to solve the violation will depend on the magnitude of the violation. When possible, the algorithm will prioritize Gate sizing to avoid additional interconnections in the netlist. Once the timing closure is achieved, the Legalization is performed to verify if there is any overlap in the placement. Any overlap violation will be solved by sliding the cells in the respective row. This concludes the discussion over the Physical Synthesis algorithms used in this thesis. Next, a brief comparison between the performance of the SA and QP placements implemented is discussed.

5.8 Simulated Annealing vs Quadratic Programming Placement

Benchmarks from ISCAS'85 [80] and the Open Cell Library (OCL) from Nangate 45nm library [81] were used for this evaluation. The table 5-2 shows the results for these benchmarks. These algorithms were executed on a laptop, equipped with an Intel core i5, 2.4 Ghz, 6 GB RAM, 64 bit word. The solutions run in an Ubuntu operating system.

Table 5-2 Simulated Annealing vs Quadratic Programming

Circuit	Number of Cells	Simulated Annealing			Quadratic Programming		
		Wire Length (um)	Run Time (s)	Area (um ²)	Wire Length (um)	Run Time (s)	Area (um ²)
C17	6	17.9	81.33	4.78	17.06	0.01	4.78
C17a	10	26	113.3	8	23.2	0.018	8
C432	168	1744.7	7925.4	248	1403	3.09	248
C499	210	2877.8	13130.2	407	2679	4.27	407

In Fig. 5-19, the total post placement wire length of the circuits placed by the SA and QP are compared. It clearly shows that the QP solution quality matches, and in certain cases, has a better performance than the SA solution. Notice that the QP detailed placement step has a strong influence in the final performance of the algorithm by performing a significant number of local optimizations.

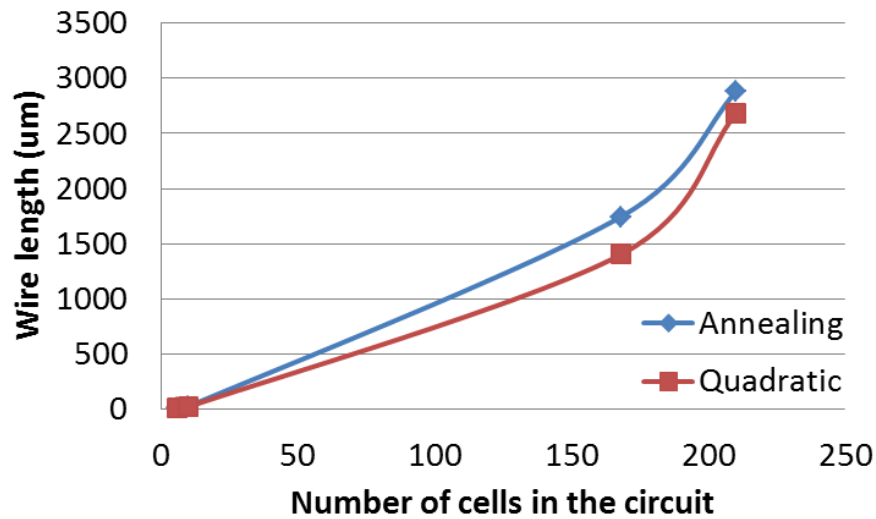


Figure 5-19 Total post placement circuit wire length of Simulated Annealing vs Quadratic Programming Placement

In Fig. 5-20, the total run time execution of both placements is shown. Note the significant performance improvement for QP over the SA algorithm. On average, the QP algorithm is about 5,016 times faster than the SA algorithm, making this the right solution to be used in this framework.

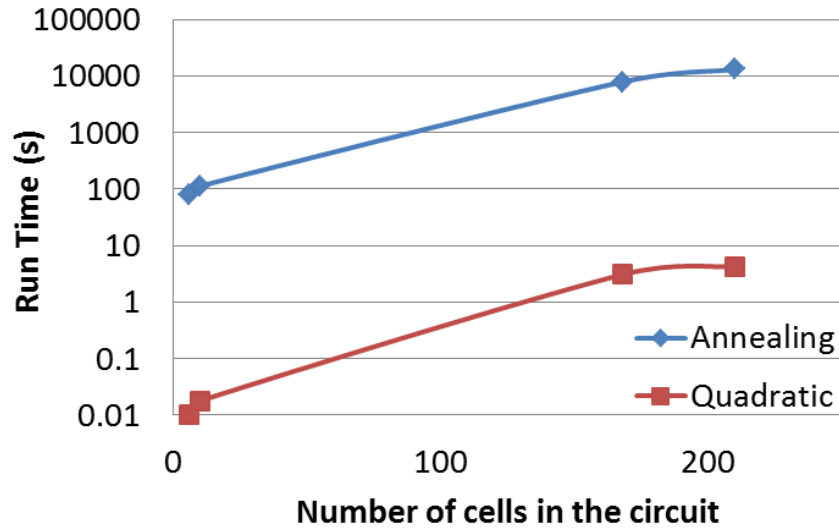


Figure 5-20 Total placement execution run time of Simulated Annealing vs Quadratic Programming Placement

These results validate the solution quality of the QP implementation due to expected near optimum solution quality compared to SA implementation. With the good run-time performance demonstrated by the QP placement, this becomes the placement algorithm of choice to have the soft-error models integrated into the flow.

5.9 Conclusions

In this chapter, the Physical Synthesis flow was presented. The algorithms and required steps to produce the Physical design of the circuit were discussed. The use of electrical correction techniques, like Gate Sizing, Cloning, Buffering and Relocation, were presented as methodologies to achieve timing closure and a quick overview of their usage to soft-error mitigation was presented.

A Physical Synthesis flow using Simulated Annealing algorithm and Quadratic placement algorithm was developed. Results from both algorithms were compared to show the improved performance and good placement quality produced by the QP algorithm. In the next chapter, the Soft Error analysis integration into the Physical Synthesis is discussed along with both electrical correction and hardening techniques evaluated.

CHAPTER 6

Soft Error Aware Physical Synthesis

As discussed in last chapter, the Physical Synthesis process is a sequential and time consuming process. Because of the size of the input netlists, ICs with billions of transistors are very common now-a-days, it is important for the Physical Synthesis process to be computational efficient. The radiation event modeling discussed in chapter 3 (Collected Charge) and in chapter 4 (SET Pulse width and Soft Error cross-section estimation) was specially designed to be integrated into the Physical Synthesis. The synthesis flow, along with the soft error evaluation modules, are shown in figure 6-1. In the figure, the soft error modules are painted with yellow color and the conventional Physical Synthesis flow with blue and green color.

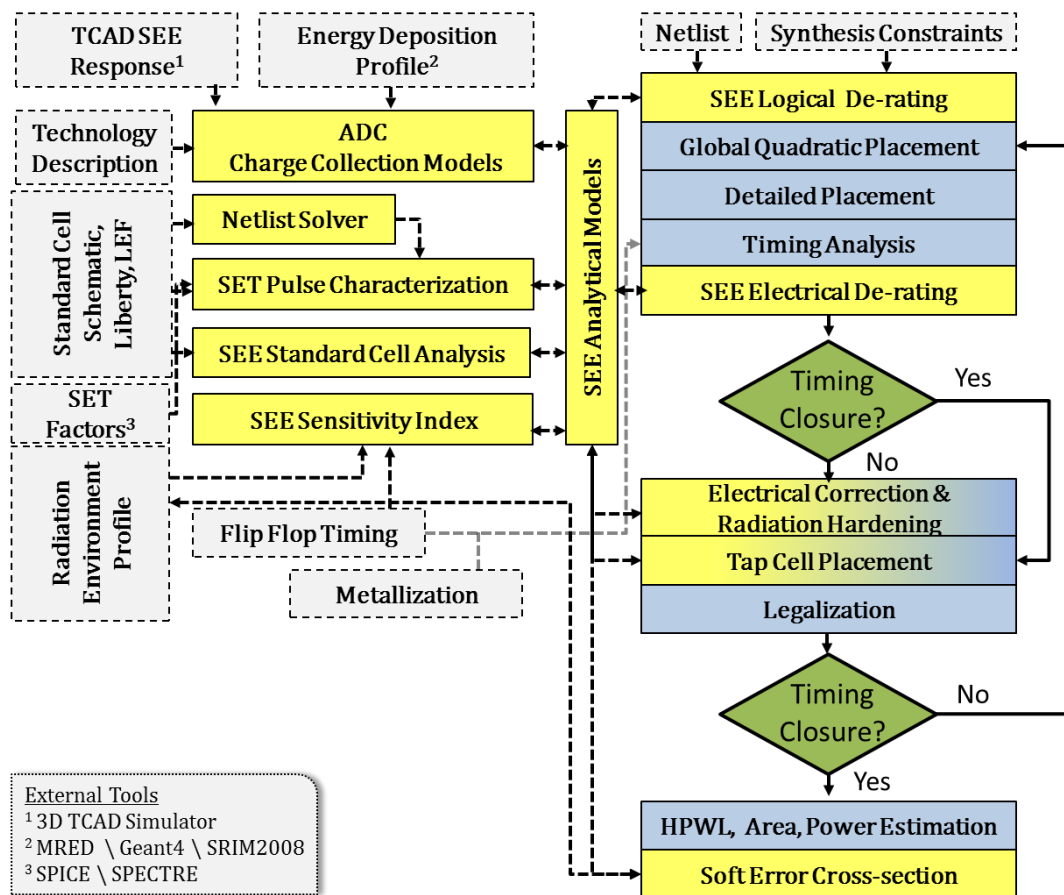


Figure 6-1 – Soft Error Aware Physical Synthesis Flow

In the figure the SEE Analytical Models is used as an interface to provide access to the soft error models developed in Chapter 3 and 4. The modules on the left side of the diagram receives Pre-Characterized data, therefore no simulation is performed during the synthesis. The soft error analysis is performed during the synthesis at different steps. .

The first soft error analysis step in Figure 6-1 is the SEE Logical De-rating step. In this step the evaluation of the Logical Masking is executed, since it can be performed before the placement of individual cells due to the no dependency in the electrical characteristics of the circuit. A single vector of primary input and output probabilities are required. The second analysis step of SEE Electrical De-rating is performed once the timing analysis has provided the HPWL and the delay characteristics of both standard cells and interconnections. In the Electrical Correction & Radiation Hardening step, the soft error analysis is again performed seeking for reduction in the soft error sensitivity. During the Tap Cell Placement step, the soft error analysis is performed if the Tap Cell Hardening was selected by the user.

In the final step, the Soft Error cross-section is analyzed by performing the Monte Carlo soft error analysis presented in chapter 4. Once the flow is completed the final placement is provided with detailed information from the HPWL and Soft Error report. In next section a quick overview of the Visual Interface of the framework is presented. Netlists from ISCAS'85 benchmark [80] and the Nangate OCL 45nm [81] were used in this thesis. For an overview over both ISCAS and the OCL used in the thesis, consult attachment 1.

6.1 Framework – Data Visualization

As discussed in the last chapter the evaluation of the SEE event is a challenging task, especially when dealing with circuits. Up to this date, most reports of the SEE analysis are done by analyzing the soft error sensitivity indexes, and statistical data from each circuit node sensitivity. In this thesis, the visualization of the SEE event is improved by providing to the Reliability Engineer detailed graphical information of

the event, to allow for a better understanding of the metrics and its circuit dependencies. For instance, the user interface (UI) developed in this work allows the real time visualization of the whole Physical Synthesis flow, along with the visualization of the SEE De-rating factors of each standard cell in a graphical interface. The SET Pulse width associated with each SEE and cell is also saved, and can be consulted by the engineer. The framework allows the user to perform the analysis with or without the UI. The UI runs in a different thread and it was implemented using GLUT OpenGL [82]. In the framework logic standard cells are represented as 2 dimensional structures (rectangles), but the environment is 3 dimensional. In figure 6-2 a picture taken from the tool UI, showing the circuit c880 (8-bit ALU) from ISCAS [80], placed using the QP placement is shown. In this figure, the circuit area boundary is indicated by the outer rectangle with white border. The primary input PADS are placed in the left corner and are painted with green color. The primary output PADS are placed in the right corner and painted with red color. The standard cells are painted with purple color and the TAP Cells, when used are painted with white color. The interconnections (wires) are represented with gray lines and can be visualized with a straight line or the equivalent HPWL (two orthogonal lines) used to estimate the wire length. Notice that the placement has enforced the use of routing channels, thus each row is separated by a spacing defined by the user.

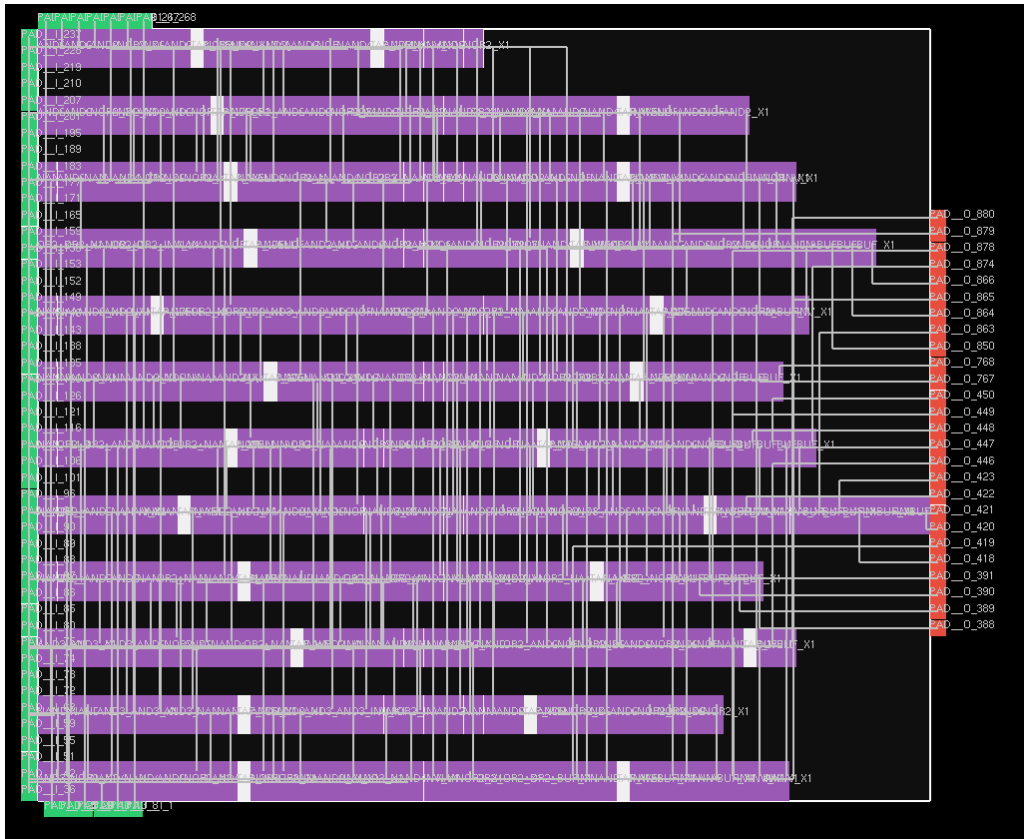


Figure 6-2 –Placed Netlist- ISCAS C880

Next the Logical De-rating can be visualized for the same circuit, and is shown in figure 6-3. In this image the standard cells are now colored according to their propagation probability. The higher the propagation probability, the more likely is the probability that the SET generated in the given cell will propagate to the output PAD. The intensity of the propagation probability is mapped from high using the red color, to low propagation probability using green color. Notice in the figure that the highest concentration of cells with high propagation probability are placed in the right corner, close to the output PADS. This is a good indication of placement quality since it is desired to have these cells near the outputs.

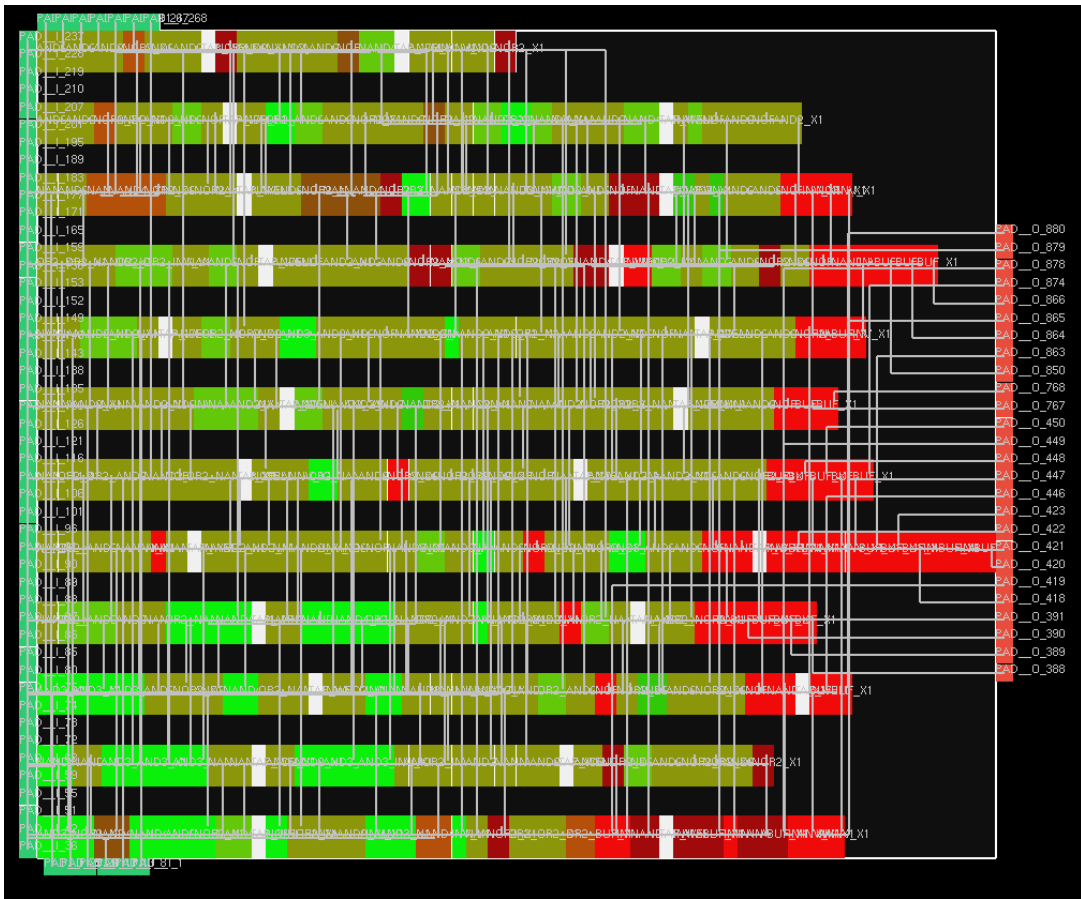


Figure 6-4 Netlist Minimum Propagation delay

The SET pulse width can be visualized in figure 6-5. In the figure cells with longer SET pulse width are painted with red color, and cells with shorter SET pulse width are painted with green color. Notice that the SET pulse width has no connection with the logical depth. Also important to observe is that standard cells closer to the Tap Cells have shorter SET pulse widths, a property that was used in this case to reduce the overall SET pulse width generated in the circuit.

netlist without any additional standard cell and modification will be referred to as the Standard netlist. The final circuit requires the changes performed in the Electrical Correction step to ensure the proper circuit operation. As the objective of this thesis is to analyze the impact of the electrical correction techniques, a very high frequency operation is set to force the use of these techniques. For the results in this chapter, a frequency of 2 GHz was used to constrain the placement. The LET spectrum was considered to have a Gaussian shape, with a constant sigma of about $0.7 \text{ MeV}\cdot\text{cm}^2/\text{mg}$, and the expectation was $5.5 \text{ MeV}\cdot\text{cm}^2/\text{mg}$ and $14 \text{ MeV}\cdot\text{cm}^2/\text{mg}$ for low and high LET analysis, respectively. To analyze the impact of each electrical correction technique, these circuits were placed using each electrical technique individually and combining all of them. In table 6-1 the area penalty of Gate Sizing, Cloning and Buffering are shown. The last column on table 6-1 indicates the total penalty of the Electrical Correction step (Considering gate sizing, cloning and buffering been used).

Table 6-1 Electrical correction techniques impact to the circuit area

Circuit	Cells	Gate Sizing		Gate Cloning		Buffering		Electrical Correction - Total	
		Sized cells	Δ Area (%)	Cloned cells	Δ Area (%)	Buffers Added	Δ Area (%)	Cells Changed	Δ Area (%)
C432	168	62	51.67	7	5.20	5	3.72	72	60.59
C499	210	16	10.67	8	2.93	24	8.80	48	22.39
C880	383	32	11.70	8	2.57	36	11.58	72	25.85
C1355	554	148	37.40	8	1.53	45	8.61	196	47.55
C1908	932	245	36.80	22	2.06	77	7.21	329	46.08
Avg	449.4	100.6	29.65	10.60	2.86	37.40	7.98	143.40	40.49

With each circuit in table 6-1 placed, its possible to analyze the soft-error cross-section trend for each of these techniques. In figure 6-6 the effect of the Gate Sizing on the soft-error cross section is evaluated for an LET of $5.5 \text{ MeV}\cdot\text{cm}^2/\text{mg}$ and an LET of $14 \text{ MeV}\cdot\text{cm}^2/\text{mg}$. In the figure the reduction of about 15.9% in soft-error cross-section for the Gate Sizing circuit is noticeable for the low LET. A 13.9% increase of the soft error cross section is observed for the high LET case. The reduction of the cross-section with the gate sizing is expected for low LET particles, due to the increased cell current drive. But this current drive increase is not effective once higher LET particles are prevalent. It is also important to notice that the cells selected to be sized were chosen based on the delay, and thus are not optimized for soft error reduction.

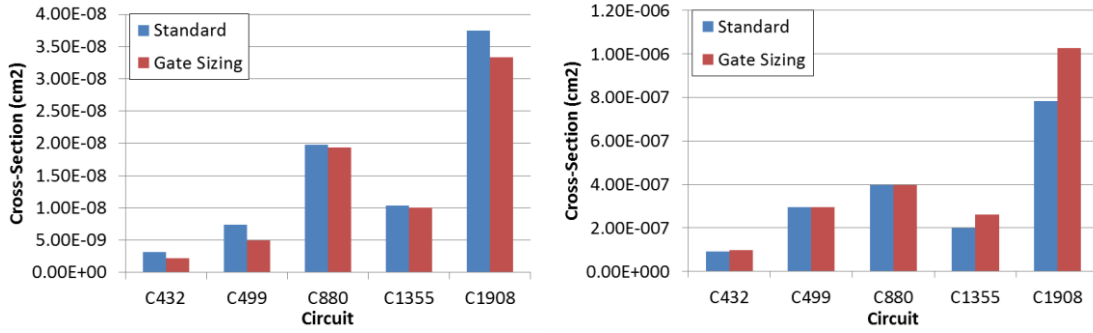


Figure 6-6 Electrical Correction – Gate Sizing – LET= 5.5 MeV*cm²/mg and 14 MeV*cm²/mg

The Gate Clone soft error cross-section is shown in figure 6-7. Notice from table 6-1 that few cloning operations were required for these circuits, so variations in the cross-section will be small. For the low LET particles the cross-section increased about 10.8%, mostly due to the impact of the gate clone for the C432 circuit. Not considering this particular circuit the soft error cross-section increase is about 2% for all other circuits. For high LET particles the soft error cross-section increases to by about 5%. The gate clone should result in an increased cross-section due to the fact that these nodes have no drive current increase, and thus not capable of reducing the soft error cross-section.

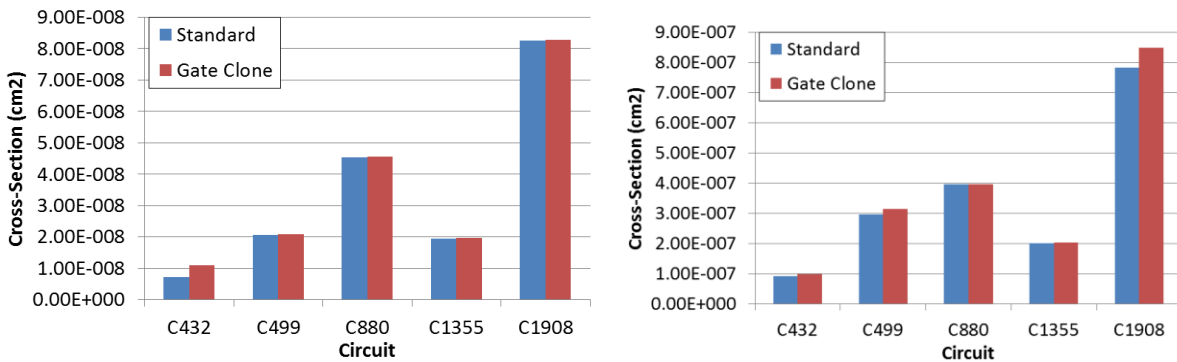


Figure 6-7 Electrical Correction – Gate Clone – LET= 5.5 MeV*cm²/mg and 14 MeV*cm²/mg

The addition of Buffers to the circuits, along with Gate Sizing is the most used electrical correction techniques to solve delay violations. In figure 6-8, the buffering impact on the cross-section is shown, with a 9% increase in cross-section for the low LET particles and a 23.7% cross-section increase for high LET particles. It is important to notice that for the buffering case, these cells were about 8% of

the circuit, and have a direct impact on the cross-section, with about 9% cross-section increase for low LET particles.

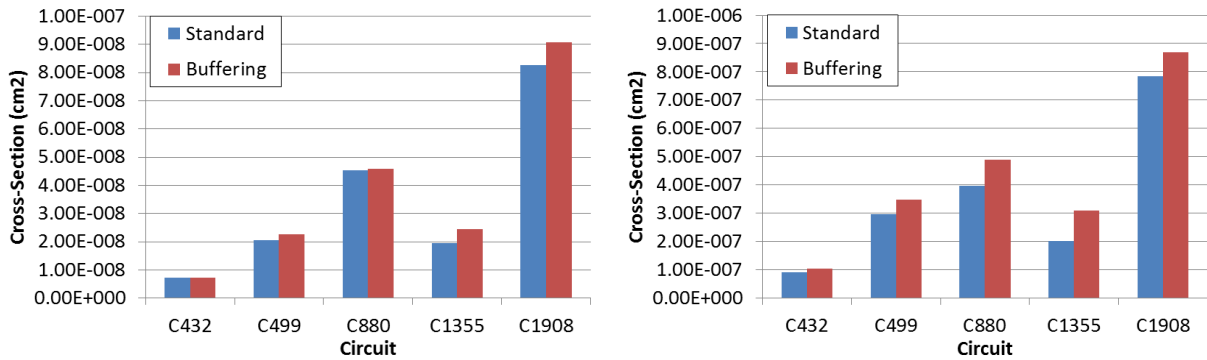


Figure 6-8 Electrical Correction – Buffering – LET= 5.5 MeV*cm²/mg and 14 MeV*cm²/mg

Now the techniques discussed before are combined, as they would be in a synthesis step. In figure 6-9 the soft error cross-section is shown for both low and high LET case. For the low LET case the soft-error cross section increases about 15%. For the high LET case the increase of the soft-error cross-section is consistent for all circuits, with an average increase of 45%.

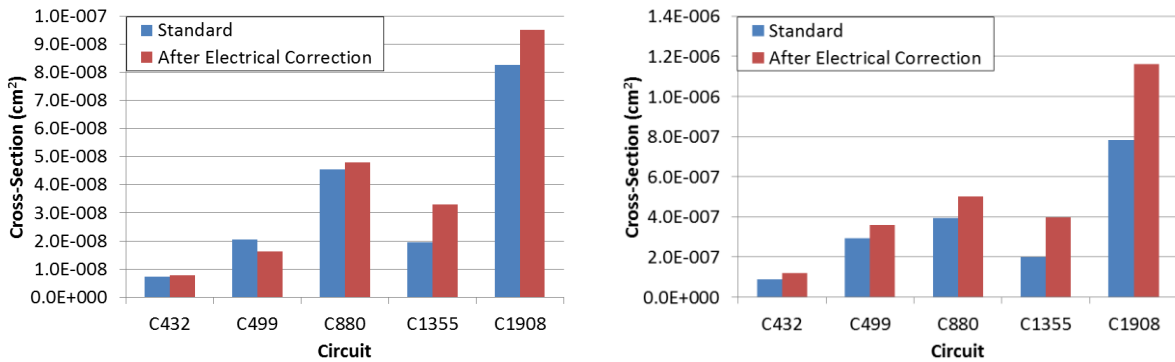


Figure 6-9 Electrical Correction – LET= 5.5 MeV*cm²/mg and 14 MeV*cm²/mg

Electrical correction techniques do have a direct impact to the circuit soft-error cross-section and should be properly evaluated in the design flow. The use of buffers and gate clone will increase the soft-error cross-section if the additional cells are vulnerable to the given LET particle, therefore these operations should be used carefully if the soft error rate is a concern. The specification of the radiation environment (fluence and energy spectrum) are very important to determinate the expected circuit soft-error cross-section. The circuit soft error response will strongly depend on the radiation environment

characteristics, since the trend observed by the electrical correction techniques used in the physical synthesis will be directly influenced by the radiation environment. In the next section, the gate sizing and the gate cloning will be applied to the circuit to reduce the soft error cross-section

6.3 Gate Sizing and Cloning for Hardening Soft Error Cross-Section

In this section the Gate Sizing and Gate Cloning are used to reduce the soft error cross-section. Notice that Gate Cloning in this case is applied as a hardening technique as discussed in chapter 5, section 2. In table 6-2 the placed netlists generated for this analysis are shown with their impact on the circuit area.

Table 6-2 Radiation hardening techniques impact to the circuit area

Circuit	Post Placement		Hardening Sizing			Hardening Cloning		
	Cells	HPWL (um)	Sized cells	Δ Area (%)	Δ HPWL (%)	Cloned cells	Δ Area (%)	Δ HPWL (%)
C432	178	2.51E+03	19	12.3	1.0	33	15.3	4.2
C499	242	3.97E+03	42	17.6	3.3	50	15.0	44.4
C880	423	5.88E+03	57	20.4	9.9	76	19.4	53.0
C1355	602	8.16E+03	110	20.0	10.4	110	14.3	51.4
C1908	1016	1.58E+04	186	16.7	23.2	186	11.9	23.2
C2670	1434	4.13E+04	156	22.2	9.7	240	20.4	41.4
C3540	1859	4.11E+04	132	7.3	15.1	340	10.2	71.2
C5315	3116	9.81E+04	276	11.5	6.7	466	11.6	48.5
C6288	2552	5.46E+04	0	0.0	0.0	26	0.7	4.5
C7552	4162	9.87E+04	631	26.4	14.6	596	12.1	41.9
Avg	1558.4	37009	160.9	15.4	9.40	212.3	13.1	38.4

For these analysis the algorithm searched for cells with average SET pulse width greater than $FF_{Setup} + FF_{Hold}$ and smaller then $2 * (T_{Setup} + T_{Hold})$. These cells were selected as good candidates to be sized and cloned because cells generating SETs with width below $T_{Setup} + T_{Hold}$ will be masked, but if the SET pulse generated by the cell is too long, the Gate Sizing and Gate Cloning will likely not be able to mitigate the SEE event. In figure 6-10 a comparison among the soft error cross-sections of the circuits from table 6-2 is shown. The gate sizing does show the best results with a 58.6% soft-error cross-section reduction to the standard placed circuit, while gate clone Sizing shows an average 31.4% soft-error cross-section reduction for the low LET case. The use of gate cloning has a considerable impact to the HPWL, with an average 38.4% wire length increase when compared with the standard placement. The addition of a new

cell to the placement row, in several situations increases the wire length associated with the duplicated cell. The increase of the RC associated with the output of the duplicated cell, can potentially produce longer SET pulses, associated with that specific standard cell, as shown in figure 4-6. In figure 6-10 the low soft error cross-section of c6288 circuit, a 16x16 multiplier, is attributed to the high logical de-rating of the circuit.

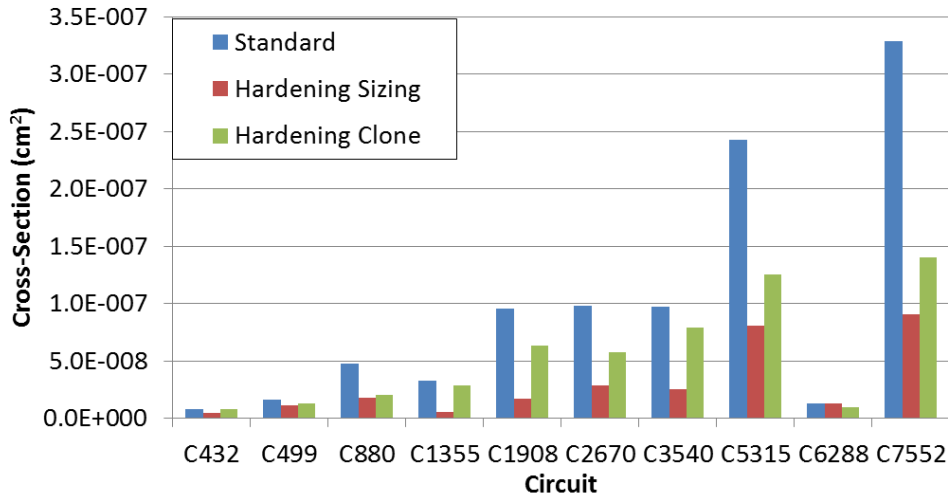


Figure 6-10 Hardening techniques Soft error cross-section - 5.5 MeV*cm²/mg

The soft-error cross-section for the high LET case is shown in figure 6-11 for a smaller number of circuits (the analysis of additional circuits will not change the expected trend). In this case the soft-error cross-section shows an average 24.7% increase for the sized circuit, and a 3.6% reduction for the cloned circuit. The high LET case reduce the effectiveness of the hardening techniques, a trend also observed when analyzing the electrical correction techniques.

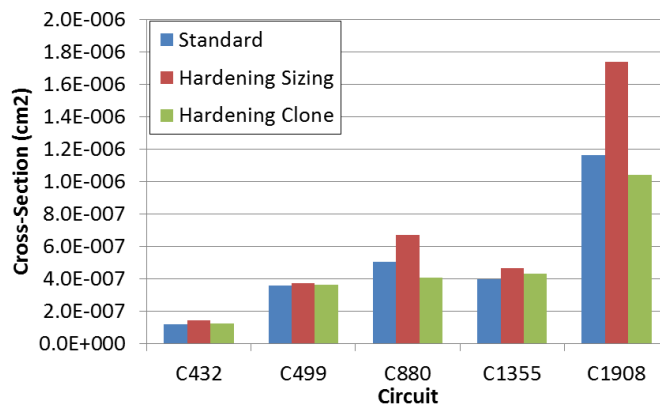


Figure 6-11 Hardening techniques Soft error cross-section -14 MeV*cm²/mg

These results indicate that for circuit hardening, for low LET particles the gate sizing is the preferable technique for reducing the soft-error cross-section reduction. In terms of area penalty Gate Sizing increases the circuit area about 17% and gate cloning increases the circuit area about 15%. The major problem with the use of Gate Cloning is the high increase of wire length (average 38% for ISCAS benchmark), having a direct impact on the circuit routability.

6.4 Tap Cell Placement Aware of Soft Errors

With the basic electrical correction and circuit hardening techniques discussed, the impact of Tap Well contacts discussed in chapter 2, section 1 to the soft error cross-section is evaluated. In this section first circuits with standard cells containing well taps (well\substrate contact) are compared with tap less cells. In figure 6-12 a comparison between cells with tap and tap less is shown, with tap less cells showing a soft error cross-section increase of about 210%. For the tap less circuits, a Tap Cell containing the well\substrate tap was added in each placement row every *40 um*. For the case shown in figure 6-12 the taps were added in regular intervals (a tap cell at the end of every row is also enforced). An increase in the soft error cross-section with tap less cells is expected due the increasing distance from the parasitic bipolar device to the well contact discussed in chapter 2, resulting in a higher resistance between the bipolar device base and the well/substrate contact.

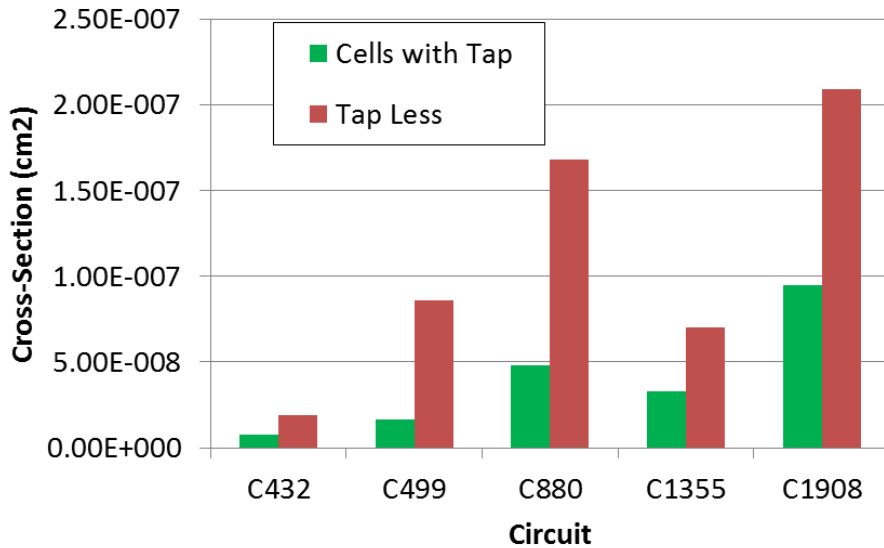


Figure 6-12 Soft-error cross-section comparison between cells with tap and tap less cells – 5.5 MeV*cm²/mg.

With the increase in soft-error cross-section resulting from the use of tap less standard cells, a placement algorithm is proposed in this thesis to reduce the impact of tap less cells to the soft error cross-section. Considering that the Standard Tap Cell algorithm places cells at regular intervals according to DRC rules, it is proposed in this thesis to optimize the tap cell position according the SET pulse width of the cells in the given placement row. In this work, the Tap Cell position is given by both a minimum required distance from the standard cells to the Tap Cell (as done in the standard tap placement algorithm) and the tap cell position that minimizes the SET pulse width in a given placement row section.

The algorithm, now referred to as the Tap Hardening algorithm simply divide the placement row into segments according to the minimum required distance from the standard cell to the tap cell (in this case 40um). Now within this segment the algorithm iterates over each cell position and calculate the total SET pulse width of all cells in the row segment. The final tap cell position will be the one with minimum total SET pulse width. For this analysis to ensure a fair comparison between the standard tap placement and the hardening tap placement, the number of Tap Cells used for each technique was constant, therefore in each placement row the number of tap cells used by the hardening technique and the standard tap placement are the same. In table 6-3 the impact of the tap cell placement techniques to the HPWL is shown, with the Tap Hardening technique causing a very small increase to the HPWL.

Table 6-3 Impact of tap placement algorithms to the HPWL

	Standard Tap HPWL	Tap Hardening HPWL
C432	2513	2527
C499	3967	4000
C880	5517	5543
C1355	9719	9753
C1908	15812	15880

In figure 6-13 (a) the circuit c1355 (32 bit Single-Error-Correcting circuit) is shown placed with the Standard Tap placement algorithm, and in figure 6-13 (b) with the Tap Hardening algorithm. For the standard tap placement, the Tap Cells with the well/substrate contacts are placed in the corners. The Tap Hardening places the tap cells in different row positions, according to the cells SET pulse widths.

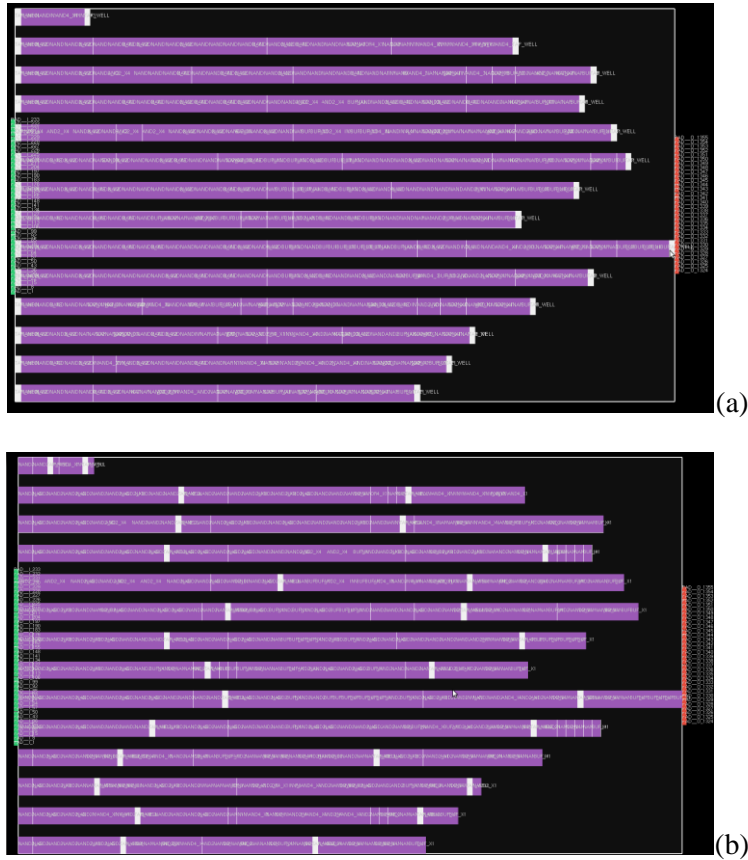


Figure 6-13 Standard Tap Placement vs Hardening Tap Placement C1355

In figure 6-14 the resulting placed circuits using the Standard Tap and Hardening algorithm are compared for a low LET particle. On average the use of the tap hardening technique reduces the soft-error

cross-section by about 28%, with no significant impact on the HPWL and no area penalty, since no cells are added to the circuit.

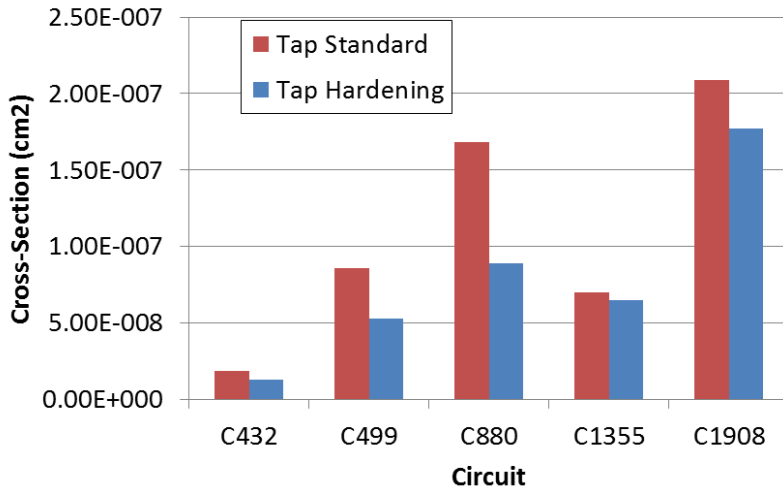


Figure 6-14 Standard Tap Placement vs Hardening Tap Placement 5 MeV*cm²/mg

For high LET particles the soft error cross-section is shown in figure 6-15. The effectiveness of the technique to reduce the SET pulse width for high LET particles is limited, resulting in a soft error cross-section reduction of about 11%.

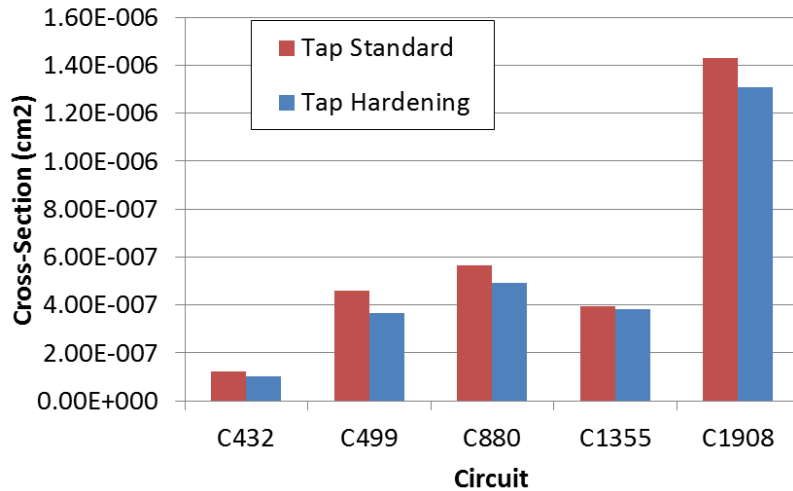


Figure 6-15 Standard Tap Placement vs Hardening Tap Placement 14 MeV*cm²/mg

To analyze the root cause of this soft-error cross-section reduction, SET pulse width estimated by the tool are analyzed. In figure 6-16 the SET pulse width distribution of circuit c880 is shown for both 5.5 MeV*cm²/mg and 14 MeV*cm²/mg particle energy distribution. It is noticeable in both distributions that

the Standard Tap placement has longer SET pulse widths when compared with the Tap Hardening. The same trend is observed for both low and high LET particles.

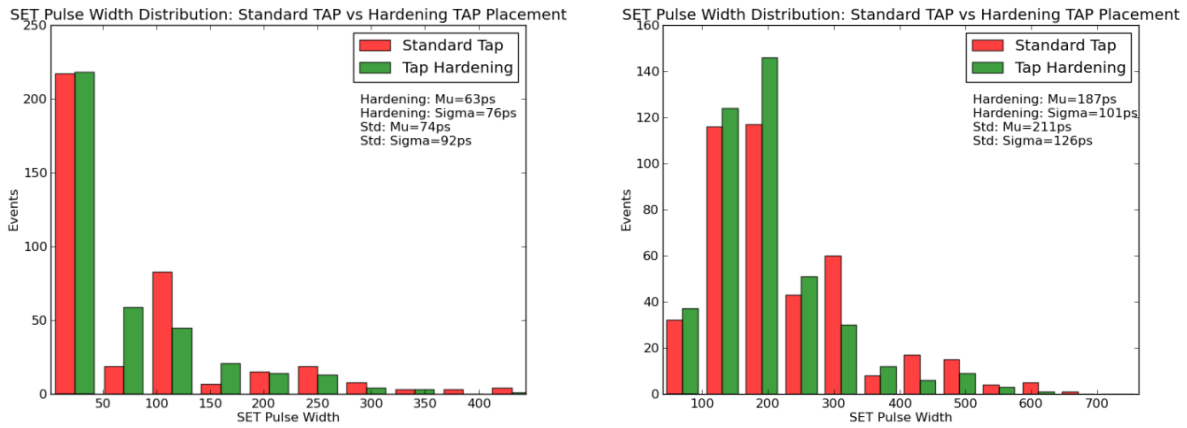


Figure 6-16 SET pulse width distribution – C880 - (a) LET=5.5 MeV*cm²/mg. (b) LET=14 MeV*cm²/mg.

The SET pulse width distribution of circuit c1355 is shown in figure 6-17. The trend observed with this circuit is the same as observed for the c880 circuit. The Tap Hardening placement does produce circuits with shorter SET pulse widths when compared with the circuits generated by the Standard Tap placement.

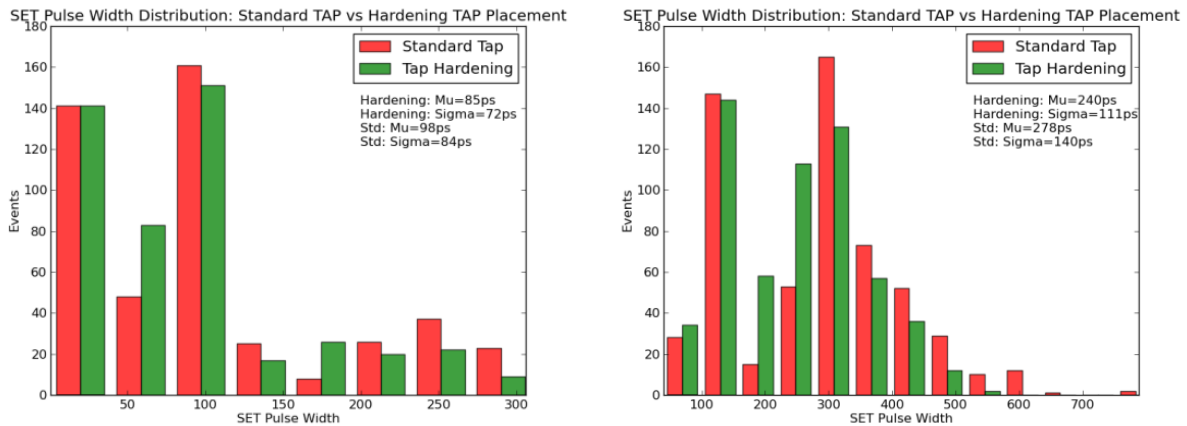


Figure 6-17 SET pulse width distribution – C1355 - (a) LET=5.5 MeV*cm²/mg. (b) LET=14 MeV*cm²/mg.

The SET pulse width can also be observed in the framework UI, as shown in figure 6-18 for circuit c880. Notice that standard cells closer to the Tap Cells (white rectangle) have shorter SET pulse widths (indicated with green).

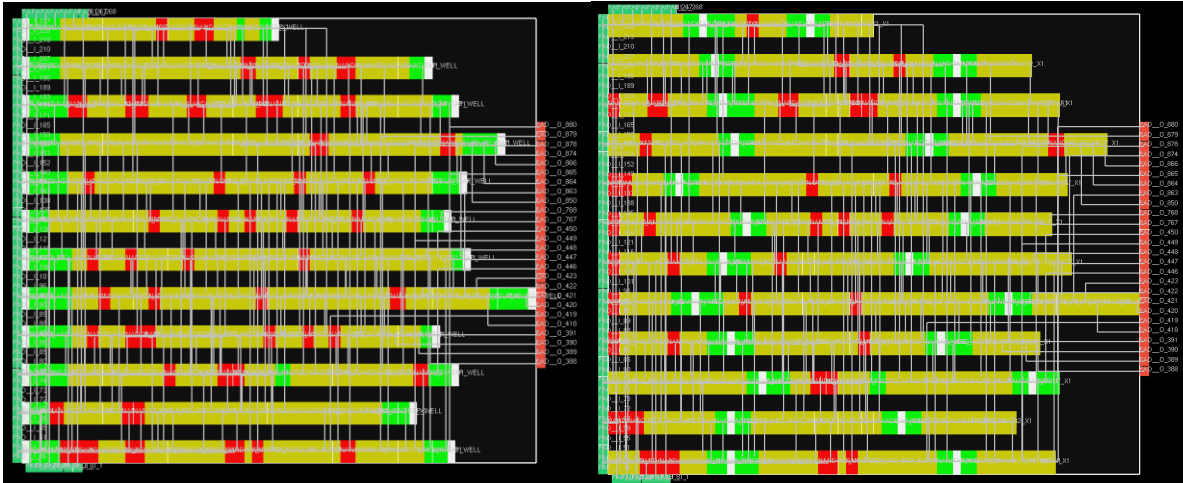


Figure 6-18 SET pulse width – c880 – Standard Tap vs Tap Hardening

The adoption of an optimization algorithm to find a better placement position for Tap Cells to reduce cells SET pulse width has a clear impact on the circuit soft-error cross-section. The possibility of selective increase in the number of Tap Cells to further reduce the soft-error cross-section should be seen as an extension of this technique, since the tap cells have a direct impact on the SET pulse width.

6.5 Conclusions

In this chapter the Physical Synthesis integration with the Soft Error models discussed in early chapters was presented. The impact of electrical correction techniques and hardening to the circuit soft error cross-section was evaluated. It was shown that for low LET particles, both buffering and gate cloning can potentially increase the circuit soft-error cross-section when the electrical correction is performed. For hardening gate sizing showed better improvement when compared with gate cloning.

The impact of the Tap Cell placement to the circuit soft-error cross-section was evaluated. A novel technique was proposed to optimize the Tap Cell position to reduce the overall SET pulse width of cells in the placement row.

CHAPTER 7

Conclusions

In this work a computationally feasible approach to model the Soft Error event in the Physical Synthesis process is presented. By using a combination of analytical models and lookup tables soft error event were modeled for different technologies, using a flexible and robust framework.

This thesis first, expands the ADC model used to estimate collected charge for the multiple transistor charge collection case. By using the ADC, detailed characteristics of the SEE event can be dynamically computed by the framework, without much computational impact. Results were validated with TCAD data, showing good model accuracy.

A methodology to estimate the SET pulse width was developed to model the SEE response of multiple standard cells, with limited pre-characterization. By using only 3 lookup tables and the standard cell circuit analysis, the SET pulse width can be accurately estimated. Results were validated using SPICE simulations and experimental SET pulse width data from a 45nm Test Chip. To measure the circuit SEE response, a Monte Carlo analysis using probability matrices was developed. The model is able to estimate the circuit soft-error cross-section with a good accuracy and it was validated using data from a 45nm Test Chip.

With the Soft Error analysis methodology implemented, the Physical Synthesis flow was modeled. Two placement algorithms based on Monte Carlo Simulated Annealing and Quadratic Programming were implemented and evaluated in this work. The Physical synthesis flow was partially implemented covering the Timing Analysis, Electrical Correction, Detailed Placement and Legalization algorithms.

The electrical correction techniques, Gate Sizing, Gate Cloning and Buffering had their impact to the circuit soft error cross-section evaluated. It was found out that the electrical correction is going to increase the circuit soft-error cross-section if the cells added to the circuit are vulnerable to the analyzed radiation

environment. The use of buffers are particular likely to increase the soft-error cross-section if the particle is able to upset these cells. The Gate Sizing and Gate Cloning as circuit hardening techniques were also evaluated. Results have shown a better efficiency of Gate Sizing to reduce the circuit soft error cross-section when compared to the gate cloning. The high wire length increase occasioned due to the Gate Cloning is another drawback to the use of this technique for hardening circuits against soft errors.

The use of the circuit hardening techniques like Gate Sizing and Gate Cloning should be used with care, due to the limitation of these techniques to mitigate low LET particles. The radiation environment should be carefully evaluated, before consider the use of these circuit operations as hardening techniques. As demonstrated in this thesis, the methodology proposed is able to detect these trends and provide valuable information about the circuit soft error response.

An algorithm to reduce the circuit soft error cross-section by optimizing the Tap Cell placement was proposed. The algorithm was able to generate circuits with consistent reduced soft-error cross-section when compared with circuits generated by standard tap cell placement algorithms. The better placement of the Tap Cells has a consistent impact on the soft-error cross-section for both low and high LET particles.

A design flow to allow the integration of the soft error analysis into the Physical Synthesis without disrupting the standard ASIC design flow used by the industry was developed. By adopting two metrics to evaluate the soft error event (soft error sensitivity index and the cross section calculation) in different steps of the design flow, the soft error evaluation can be effectively integrated into the Physical Synthesis.

Lastly, the capabilities of the resulting framework developed in this thesis to analyze circuits and provide detailed information from the soft error event was shown. The framework enables the traceability of the SEE events to identify the root-cause of events and circuit characteristics impacting the standard cell response.

Future work includes the modeling of collected charge and charge sharing on *FinFET* transistors and the evaluation of this analysis to estimate the soft-error cross-section of sequential circuits.

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Open Cell Library and Benchmarks

For the proper evaluation of the standard cell response to soft errors, the use of a standard cell library designed by a commercial company becomes important. In this thesis the Nangate FreePDK 45nm Open Cell Library (OCL) [81] was used for characterizing the standard cells.

A.1 Open Cell Library

The OCL 45nm library uses a constant cell height of 1.63 μ m with the length of both NMOS and PMOS devices equal to 50nm. The maximum width for the NMOS transistor is 415nm and for the PMOS 630nm. The standard cell INV_X1, is shown in figure A-1 for dimensional reference. Notice that the increase of the gate size does not change the cell height but cell width, with the addition of transistors in parallel.

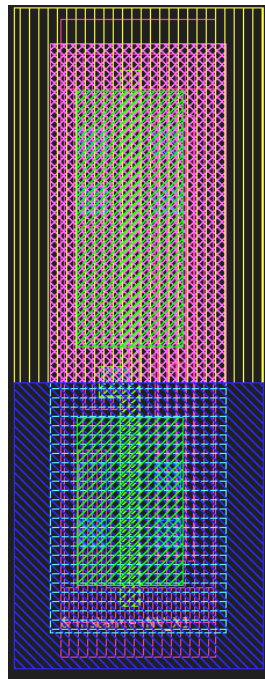


Figure A-0-1 - INV_X1 Standard Cell - OCL 45nm [81]

The use of the OCL was especially important for the SET Pulse Width estimation and Timing Analysis. For the SET Pulse Width estimation provided the standard cell schematic files for simulation of both standard and complex gates. For the Timing Analysis the OCL provided the delay and electrical characteristics of the target cells. This information is made available by the libraries contained in the OCL package. A special package is the Liberty Library containing detailed characterization of timing, power and noise of the standard cells. In table A-1 the snapshot of the full-cycle delay table is shown for some standard cells used in this work.

Table A--1 The full cycle delay of standard cells, over different load capacitances. Delay provided in ns.

Cell\Load	0.36fF	15.14fF	30.28fF	60.57fF
.....
AND2_X1	0.13	0.2	0.26	0.37
AND3_X1	0.15	0.23	0.29	0.4
AND4_X1	0.17	0.25	0.31	0.42
AOI222_X1	0.16	0.34	0.4	0.4549
AOI22_X1	0.11	0.27	0.38	0.4349
.....
.....
INV_X1	0.03	0.18	0.27	0.41
MUX2_X1	0.17	0.24	0.3	0.41
NAND2_X1	0.06	0.21	0.31	0.46
NAND3_X1	0.08	0.23	0.34	0.51
NAND4_X1	0.1	0.25	0.37	0.55
.....

The metallization information from the interconnection was obtained from [77]. In some cases values were interpolated to complete the table A-2. In A-2 the first column indicates the technology node, followed by the minimum metal 1 interconnection width, thickness, distance to the substrate and the sheet resistance. Distance units in micro meters. This information is particularly useful during the Time Analysis when computing the Elmore Delay,

Table A-2 Metallization characteristics

KIND	NODE	M1_MINW	M1_T	M1_SUB_H	M1_SHEETR
BULK	130	0.13	0.62	0.6	0.1
BULK	90	0.09	0.45	0.4	0.16
BULK	65	0.065	0.303	0.27	0.25
BULK	40	0.045	0.202	0.18	0.38
BULK	20	0.02	0.12	0.108	0.4

Flip flop information required for both Time De-rating calculation and the Timing Closure estimation are shown in table A-3. The delay data from Setup Time and Hold Time refer to the full cycle delay, and is given in nano seconds. The Flip Flop Leakage is given in nano Watts.

Table A--3 Flip Flop Technology Information

KIND	NODE	FF_SETUP_TIME	FF_HOLD_TIME	FF_LEAKAGE
BULK	130	0.1	0.11	175.21
BULK	90	0.09	0.1	144.14
BULK	65	0.058	0.065	110.1
BULK	40	0.035	0.04	79.11
BULK	20	0.02	0.015	49.44

The analysis over the benchmarks to be presented in the next section have indicated the standard cells to be used in this work, and shown in table A-4. Information regarding the number of transistors, cell width and cell height are shown for reference.

Table A--4 - OCL Cells used by the benchmarks

[45nm]				
Standard Cells	Transistors	W(um)	H(um)	Area(um2)
INV_X1	2	0.61	1.63	0.9943
INV_X2	4	0.8	1.63	1.304
INV_X4	8	1.18	1.63	1.9234
INV_X8	16	1.94	1.63	3.1622
INV_X16	32	3.46	1.63	5.6398
BUF_X1	4	0.8	1.63	1.304
NAND2_X1	4	0.8	1.63	1.304
NAND3_X1	6	0.99	1.63	1.6137
NAND4_X1	8	1.18	1.63	1.9234
NAND2_X2	8	1.18	1.63	1.9234
NAND3_X2	12	1.56	1.63	2.5428
NAND4_X2	16	1.94	1.63	3.1622
AND2_X1	6	0.99	1.63	1.6137
AND3_X1	8	1.18	1.63	1.9234
AND4_X1	10	1.37	1.63	2.2331
NOR2_X1	4	0.8	1.63	1.304
NOR3_X1	6	0.99	1.63	1.6137
NOR4_X1	8	1.18	1.63	1.9234
NOR2_X2	8	1.18	1.63	1.9234
NOR3_X2	12	1.56	1.63	2.5428
NOR4_X2	16	1.94	1.63	3.1622
OR2_X1	6	0.99	1.63	1.6137
OR3_X1	8	1.18	1.63	1.9234

OR4_X1	10	1.37	1.63	2.2331
XOR2_X1	10	1.37	1.63	2.2331
OAI21_X1	6	0.99	1.63	1.6137
OAI21_X2	12	1.56	1.63	2.5428
OAI22_X1	8	1.18	1.63	1.9234
OAI22_X2	16	1.94	1.63	3.1622
AOI21_X1	6	0.99	1.63	1.6137
AOI21_X2	12	1.56	1.63	2.5428
AOI22_X1	8	1.18	1.63	1.9234
AOI22_X2	16	1.94	1.63	3.1622

Next section the benchmarks to be used by this thesis are summarized.

A.2 Benchmarks

To represent common circuit designs used by the industry the benchmarks from ISCAS 85 were used, due to its broad range of circuits design. A summary of some ISCAS 85 netlists is shown in figure A-2.

Circuit	Function	No. of input lines	No. of output lines	No. of logic gates	No. of major functional blocks
c432	27-channel interrupt controller	36	7	160	5
c499	32-bit SEC circuit	41	32	202	2
c880	8-bit ALU	60	26	383	7
c1355	32-bit SEC circuit	41	32	546	2
c1908	16-bit SEC/DED circuit	33	25	880	6
c2670	12-bit ALU and controller	233	140	1,193	7
c3540	8-bit ALU	50	22	1,669	11
c5315	9-bit ALU	178	123	2,307	10
c6288	16 × 16 multiplier	32	32	2,406	240
c7552	32-bit adder/comparator	207	108	3,512	8

Figure A--2- ISCAS'85 Benchmark details [83]

As seen in figure A-2 the circuits evaluated in this work differ significantly in both functional purpose and architectural characteristics. The number of cells varies from 160 to 3,512 standard cells with the number of primary inputs ranging from 36 to 233 ports. The number of outputs range from 7 to 140 ports.

Publications

List of Publications

Journals

1. T.R. Assis, K. Ni, J. Kauppila, B. L. Bhuvu, R.D. Schrimpf, L.W. Massengill, S. Wen, R. Wong, C. Slayman. **Estimation of Single-Event Induced Collected Charge for Multiple Transistors Using Analytical Expressions.** IEEE Transactions on Nuclear Science. 2015.
2. J.S. Kauppila, W.H. Kay, T.D. Haeffner, D.R. Rauch, T.R. Assis, N.N Mahatme, J. Gaspard, B.L. Bhuvu, M.L. Alles, W.T. Holman, L.W. Massengill. **“Single-Event Upset Characterization Across Temperature and Supply Voltage for a 20nm Bulk Planar CMOS Technology”** IEEE Transactions on Nuclear Science. 2015
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Conferences

1. T.R. Assis, K. Ni, J. Kauppila, B. L. Bhuvu, R.D. Schrimpf, L.W. Massengill, S. Wen, R. Wong, C. Slayman. **Estimation of Single-Event Induced Collected Charge for Multiple Transistors Using Analytical Expressions.** IEEE Nuclear and Space Radiation Effects Conference 2015.
2. T.R. Assis, N. J. Gaspard, N. N. Mahatme, R. A. Reed, R. A. Weller, M. H. Mendenhall, B. L. Bhuvu, L. W. Massengill, E. Blackmore, M. Trinczek, S. A. Wender. **Neutrons and Protons Radiation Induced Single Event Effects in an Advanced Technology Node.** Single Event Effects (SEE) Symposium, May. 2014.
3. Mahatme, N.N.; Gaspard, N.J.; Assis, T.; Jagannathan, S.; Chatterjee, I.; Loveless, T.D.; Bhuvu, B.L.; Massengill, L.W.; Wen, S.J.; Wong, R., **“Impact of technology scaling on the combinational logic soft error rate,”** *Reliability Physics Symposium, 2014 IEEE International* , vol., no., pp.5F.2.1,5F.2.6, 1-5 June 2014
4. Narasimham, B.; Chandrasekharan, K.; Wang, J.K.; Djaja, G.; Gaspard, N.J.; Mahatme, N.N.; Assis, T.R.; Bhuvu, B.L., **“High-speed pulsed-hysteresis-latch design for improved SER performance in 20 nm bulk CMOS process,”***Reliability Physics Symposium, 2014 IEEE International* , vol., no., pp.5F.4.1,5F.4.5, 1-5 June 2014
5. Adrian Evans (Cisco), Michael Nicolaidis (IMAG-FR), Shi-Jie Wen(Cisco), Thiago Assis(Vanderbilt) B. Bhuvu (Vanderbilt). **Clustering Techniques and Statistical Fault Injection for Selective Mitigation of SEUs in Flip-Flops.** International Symposium on Quality Electronic Design (ISQED) 2013

Industry Awards\Competitions

1. Cisco Systems - 1st Place Supply Chain Operation Case Competition – 2014
2. Cisco Systems - 1st Place Supply Chain Operation Case Competition – 2013