# Efficient Binary Decision Diagram Manipulation by Reducing the Number of Intermediate Nodes 

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#### Abstract

The complexity of hardware systems has increased significantly in recent decades. Due to increasing user requirements, there is a need to develop more efficient data structures and algorithms to guarantee the correct behavior of such systems. A Reduced Ordered Binary Decision Diagram (BDD) is a suitable data structure as it represents all Boolean functions canonically given a variable order as well as provides algorithms for efficient manipulation. However, BDDs also have challenges: practicability depends on their minimization and there is a large memory consumption for some complex functions. To address these issues, this work investigates the number of emerged intermediate nodes that are not used in the final BDD result and presents a novel approach for efficient BDD manipulation by reducing the number of such nodes. Experiments on BDD benchmarks show that peak BDD node sizes can be significantly reduced, leading to accelerated BDD manipulation.

Index Terms-Boolean functions, binary decision diagrams, software packages, formal verification, model checking


## I. Introduction

Moore's law describes that the number of transistors in Integrated Circuits (ICs) doubles every two years [1]. Due to technological progress, billions of transistors are nowadays present in Very Large Scale Integration (VLSI) circuits that can be found, i. a., in smartphones used by today's society. As ICs become more complex, VLSI design cannot be created without Computer-Aided Design (CAD), making it an essential part of the hardware design process known as VLSI CAD [2].

In order to meet time-to-market constraints and to guarantee the quality of VLSI CAD, continuous algorithmic improvements in the field of verification are necessary. Model checking is an important approach to assess the correctness of hardware systems through state exploration and property checking [3]. Traditionally, data structures have been implemented that explicitly consider system states [4]. Thus, only automata with at most $10^{6}$ reachable states could be processed [5]. However, since real models consist of billions of states, they cannot be considered in a reasonable amount of time [6]. Reduced Ordered Binary Decision Diagrams (BDDs) [7] are suitable for this application as they can compactly encode Boolean functions and allow efficient algorithms, such as reachability analysis, leading to a breakthrough in this technique [8].

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Therefore, extensive research has been conducted to improve verification techniques by developing components that are combined in BDD packages [9]. Components include, i. a., BDD manipulation by the If-Then-Else (ITE) algorithm, hashbased Computed Tables (CTs) for caching operations, Unique Tables (UTs) with collisions resolved by chaining to identify all existing nodes, and Garbage Collection (GC) methods to remove unused (dead) nodes [10].
Although a BDD is an efficient data structure for Boolean functions, there are existing challenges: practicability depends on their minimization and there is a large memory consumption for some complex functions like multipliers [11], [12]. Firstly, the final BDD size of many functions like adders depends on the variable order [13], [14]. Secondly, when logical operations are performed repeatedly during BDD manipulation, the number of UT nodes temporarily used for final BDD construction, called Intermediate Nodes (Inodes), can dramatically increase, leading to memory overflow, so that subsequent operations such as reachability analysis cannot be performed [15], [16], [17], [18], [19], [20].
To address these issues, in this paper we investigate the emergence of Inodes and their number during BDD manipulations. Based on these results, we propose a novel approach for efficient BDD manipulation by reducing the number of Inodes. Experiments confirm that using this approach can significantly reduce the greatest number of nodes in use at any point during the process lifetime, leading to accelerated BDD manipulation that is on average about $20 \%$ faster compared to the single use of the ITE algorithm and related work.

In summary, the main contributions are as follows:

1) Investigation of the number of Inode emergences using Boolean functions;
2) Development of an approach to reduce Inodes and accelerate BDD manipulation;
3) Approach evaluation and comparison with related work.

This paper is organized as follows: We summarize the Boolean function fundamentals and related work in Section II. Section III investigates Inode emergences and describes the proposed approach based on observations. In Section IV, the efficiency of this approach is evaluated and experimental results are discussed. Finally, Section V concludes the paper.

## II. BACKGROUND

This section introduces important fundamentals in an attempt to keep this work self-contained. While Section II-A describes concepts for understanding Boolean functions, Section II-B briefly discusses related proposals for increasing the efficiency of memory use during function manipulations.

## A. Preliminaries

In ICs, signals can be symbolized by variables $x_{1}, \ldots, x_{n}$ taking logical values from $\mathbb{B}:=\{0,1\}$. In statement logic, $0(1) \in \mathbb{B}$ is interpreted as false (true). Thus, outputs whose values are specified by inputs can be described by mathematical mappings like Boolean functions.

Definition 1. A mapping $f: \mathbb{B}^{n} \rightarrow \mathbb{B}^{m}$ is called a Boolean function, where $n, m \in \mathbb{N} . \mathcal{B}_{n, m}:=\left\{f \mid f: \mathbb{B}^{n} \rightarrow \mathbb{B}^{m}\right\}$ describes the set of Boolean functions, where $\mathcal{B}_{n}:=\mathcal{B}_{n, 1}$.

The Boolean calculus [21], the basis for today's computer systems, allows computations with Boolean functions as well as their manipulation and defines algebraic structures.
Definition 2. The quadruple $\left(\mathcal{B}_{n},+, \cdot,{ }^{-}\right)$with

$$
\begin{aligned}
f+g \in \mathcal{B}_{n} & :=(f+g)(\alpha)=f(\alpha) \vee g(\alpha) \forall \alpha \in \mathbb{B}^{n} \\
f \cdot g \in \mathcal{B}_{n} & :=(f \cdot g)(\alpha)=f(\alpha) \wedge g(\alpha) \forall \alpha \in \mathbb{B}^{n} \\
\bar{f} \in \mathcal{B}_{n} & :=\bar{f}(\alpha)=1 \Longleftrightarrow f(\alpha)=0 \forall \alpha \in \mathbb{B}^{n}
\end{aligned}
$$

is called the Boolean algebra of functions.
Based on Definition 2, properties can be derived including but not limited to commutative, absorption, resolution, and annulment laws [22]. Properties come in pairs, i. e. the dual of a Boolean Expression (BE) - a common representation of Boolean functions - is obtained by interchanging + with $\cdot$ and 0 with 1 [23].

Definition 3. Let $X_{n}=\left\{x_{1}, x_{2}, \ldots, x_{n}\right\}$ be a variable set and $\Sigma=X_{n} \cup\{0,1,+, \cdot,-,()$,$\} be an alphabet. The set B E\left(X_{n}\right)$ over $X_{n}$ is the subset of $\Sigma^{*}$ that is defined inductively:

Identity elements 0 and 1 , and variables are BEs.
If $g$ and $h$ are BEs, then the disjunction $g+h$, conjunction $g \cdot h(g h)$, and negation $\bar{g}$ are also BEs.
Nothing else is a BE.
If an IC is modular, a technique such as model checking can first compute single representations in order to subsequently combine them [12], enabled by Shannon expansion [24].
Definition 4. Let $f \in \mathcal{B}_{n}$ be a $n$-ary function. The partitioning $f$ to $x_{i}$ with $f_{x_{i}=1}\left(\alpha_{1}, \alpha_{2}, \ldots, \alpha_{i-1}, 1, \alpha_{i+1}, \alpha_{i+2}, \ldots, \alpha_{n}\right)$ and $f_{x_{i}=0}\left(\alpha_{1}, \alpha_{2}, \ldots, \alpha_{i-1}, 0, \alpha_{i+1}, \alpha_{i+2}, \ldots, \alpha_{n}\right) \forall \alpha \in \mathbb{B}^{n}$ is called the Shannon expansion

$$
f=x_{i} \cdot \underbrace{f_{x_{i}=1}}_{\text {positive cofactor }}+\bar{x}_{i} \cdot \underbrace{f_{x_{i}=0} .}_{\text {negative cofactor }}
$$

If variables are successively decomposed using Definition 4 respecting a total order $\pi$ and avoiding redundancies/isomorphisms by exploiting laws of Boolean algebra, a $B D D$ results.

(a) $x_{1}+x_{2}$

(b) $x_{1} \cdot x_{2}$

(c) $\bar{x}_{1}$

Fig. 1: BDDs representing the Boolean basis functions

Definition 5. A $B D D$ is a directed acyclic graph $G=(V, E)$ over variables $X_{n}:=\left\{x_{1}, x_{2}, \ldots, x_{n}\right\}$ and a value set $\mathbb{B}$. Each node is assigned such a label where a Boolean function $f$ is interpreted as follows:

If $v$ is labeled with $b \in \mathbb{B}$, then the leaf represents the constant function.
If $v$ is an inner node, it is labeled with $x_{i} \in X_{n}$, where the variable is decomposed by $x_{i} \cdot f_{v}+\bar{x}_{i} \cdot f_{\bar{v}}=\left(x_{i}, f_{v}, f_{\bar{v}}\right)$ respecting a total order $\pi: x_{1}<x_{2}<\ldots<x_{n}$, where $f_{v}$ is the high child and $f_{\bar{v}}$ is the low child referenced by the parent $v$. The edge set $E$ contains all such references.
If $\left(f_{v}\right)_{x_{i}} \neq\left(f_{v}\right)_{\bar{x}_{i}} \forall v \in V$ and no distinct nodes $v, w \in V$ exist which are labeled with the same variable and whose children are identical, then $G$ is called reduced.

Example 1. BDDs for the Boolean functions $+, \cdot,-\in \mathcal{B}_{2}$ are shown in Fig. 1: 1) disjunction (Fig. 1a), 2) conjunction (Fig. 1b), and 3) negation (Fig. 1c).

Remark. The referencing is typically drawn using solid edges (1-edges) and dashed edges ( 0 -edges).

Applying Definition 4 successively top-down to build a BDD requires repeatedly performing an equivalence test to check whether subfunctions are already represented [2]. A more efficient way is to transform Definition 4 using Boolean algebra laws and combine nodes via

$$
f \otimes g=x_{i} \cdot\left(f_{x_{i}=1} \otimes g_{x_{i}=1}\right)+\bar{x}_{i} \cdot\left(f_{x_{i}=0} \otimes g_{x_{i}=0}\right)
$$

where $\otimes \in \mathcal{B}_{2}$. These operations can be traced back to

$$
\operatorname{ITE}(f, g, h)=f \cdot g+\bar{f} \cdot h
$$

that computes If $f$ Then $g$ Else $h$ such as $f+h=\operatorname{ITE}(f, 1, h)$, which is compatible with Definition 4 because of

$$
f \cdot g+\bar{f} \cdot h=\left(x_{i}, \operatorname{ITE}\left(f_{x_{i}}, g_{x_{i}}, h_{x_{i}}\right), \operatorname{ITE}\left(f_{\bar{x}_{i}}, g_{\bar{x}_{i}}, h_{\bar{x}_{i}}\right)\right)
$$

The resulting triple corresponds to Definition 5 and formulates the conventional divide-and-conquer Algorithm 1 [25] that constructs BDDs by a sequence of logical operations starting from single nodes. At first, terminal cases are checked in Line 1. Using a CT, Lines $2-4$ check if an operand combination has already been computed. Otherwise, two cofactors are computed in Lines 5-7 decomposing according to the previously determined order of variables. Lines 8-10 check for isomorphism. In Line 11, canonicity is ensured by either finding or adding the computed triple into the UT $u t$.

Algorithm 1: Conventional BDD manipulation operator ITE

```
Input: BDDs \(f, g, h\)
Output: Constructed BDD based on \(f, g, h\)
    . . .
        \(\triangleright\) terminal cases
    if ct.has_entry \((f, g, h)\) then
        return \(\operatorname{ct}(f, g, h)\)
    end if
    \(x \leftarrow\) top variable of \(f, g, h\)
    \(t \leftarrow \operatorname{ITE}\left(f_{x_{i}}, g_{x_{i}}, h_{x_{i}}\right)\)
    \(e \leftarrow \operatorname{ITE}\left(f_{\bar{x}_{i}}, g_{\bar{x}_{i}}, h_{\bar{x}_{i}}\right)\)
    if \(t=e\) then
        return \(t\)
    end if
    \(r \leftarrow u t . f i n d \_o r \_a d d(x, t, e)\)
    ct.insert \((f, g, h, r)\)
    return \(r\)
```

| Node | Label | Then | Else |
| :---: | :---: | :---: | :---: |
| $v_{0}$ | 0 | - | - |
| $v_{1}$ | 1 | - | - |
| $v_{2}$ | $x_{2}$ | $v_{1}$ | $v_{0}$ |
| $v_{3}$ | $x_{1}$ | $v_{1}$ | $v_{2}$ |
| $v_{4}$ | $x_{3}$ | $v_{1}$ | $v_{0}$ |
| $v_{5}$ | $x_{2}$ | $v_{1}$ | $v_{4}$ |
| $v_{6}$ | $x_{1}$ | $v_{1}$ | $v_{5}$ |

(a) UT

(b) $f$

(c) $h$

(d) $\operatorname{ITE}(f, 1, h)$

Fig. 2: Performing ITE, where $f=x_{1}+x_{2}$ and $h=x_{3}$

Depending on the current number of dead nodes or if there is not enough space to add new nodes, a GC is performed to periodically free unused memory with subsequent table expansions if necessary. Finally, Line 12 stores the result in the CT followed by the return of the constructed BDD.

Example 2. Let $\rho \in \mathcal{B}_{3}$ with $\rho\left(x_{1}, x_{2}, x_{3}\right)=x_{1}+x_{2}+x_{3}$, where $\pi: x_{1}<x_{2}<x_{3}$. Fig. 2 illustrates the UT (Fig. 2a) with $\operatorname{BDD}$ nodes after performing $\operatorname{ITE}(f, 1, h)$ on the basis of $f=x_{1}+x_{2}$ (Fig. 2b), $g=1$, and $h=x_{3}$ (Fig. 2c), where Fig. 2d represents the constructed final BDD. It can be seen that during the BDD manipulation there are no more references to the nodes $v_{2}$ and $v_{3}$, i.e. they are dead.

Based on these components, Algorithm 1 can be carried out in a time almost linear to the number of BDD nodes assuming an ideal UT and CT, i.e. checking and storing nodes in constant time.

## B. Related Work

In practical applications, logical operations such as Algorithm 1 are performed successively [5]. Beyond the theoretical point of view w.r.t. complexity described in the last section, this can dramatically increase the UT size which leads to a failure (or a significant slowdown) of the computation caused by a memory overflow.

Although optimizations for choosing a proper variable order have been developed in recent years to address this issue [26], the reduction of intermediate computations has not been investigated to this extent: While algorithms such as [15] create BDD nodes only when such are represented in the final result, there exist approaches like [16] that get final BDDs top-down and bottom-up. Below is a brief explanation of these methods.

The so-called Multi-way method [15] exploits implicit don't cares in a BE including a cube cofactorization by recursively applying Definition 4 to address wasteful intermediate computations. For example, when computing a BE like $f g+h$, each of $f$ and $g$ can be minimized using $h$ as a don't care set before the product is formed. That is, for any subexpression where $h$ is 1 , the final result will also be 1 on that subexpression regardless of the logical value of $f$ or $g$. Thus, the creation of Inodes is prevented and only the final BDD is constructed.
The second method, called XTop [16], first performs a topological analysis top-down to find good decomposition points for the IC. To this end, a cut-set is computed by reducing the number of compositions and dependent variables w.r.t. the outputs, where a cut-set is a set of gates such that any path from an IC input to an IC output has to cross through one of the gates in the IC. Based on selected decomposition points, single BDDs are then constructed bottom-up using Algorithm 1, which are finally composed.
Although the aforementioned methods have proven successful for various problems, such as formal design verification and graph-theoretic problems, they come with some drawbacks: While with Multi-way from [15] the explosion in the number of BE operations exists, XTop from [16] has problems with essential algorithms like dynamic variable ordering. Therefore, the main goal of this work is to overcome these limitations.

## III. Reducing the Number of Intermediate Nodes

In this section, we describe the core development from our investigation to the approach of reducing the number of Intermediate Nodes (Inodes). First, the emergence of Inodes and their number are investigated in Section III-A. Second, in Section III-B, we present our approach to allow systematic reduction of the number of Inodes.

## A. Investigation of the number of emerged Inodes

In BDD packages, complicated BEs in the form of BDDs are constructed by logically combining single BDD nodes using Algorithm 1. If combining is performed conventionally, this can lead to unnecessary memory and runtime overhead due to intermediate results that are not in the final BDD.

Considering Example 2, for the construction of $x_{1}+x_{2}+x_{3}$ first $f=x_{1}+x_{2}$ is built, which is afterwards combined with $h=x_{3}$ via logical disjunction. However, the problem with this combination is that two Inodes $\left(v_{2}, v_{3}\right)$ emerge since they are not used in the final BDD (Fig. 2d).
Remark. In the case of dead nodes, it is not clear during BDD manipulation whether they are finally intermediate, since they can be reactivated and thus also be in the final result [11]. Every Inode is dead, but not every dead node is intermediate.


Fig. 3: ITE constructions for $x_{1}+x_{2}+x_{3}$, where $f=x_{1}+x_{2}$, $h=x_{3}, f^{\prime}=x_{3}+x_{2}$, and $h^{\prime}=x_{1}$

Via an ITE walkthrough, it can be observed why these Inodes emerge and how they can be prevented, which is illustrated in Fig. 3: Respecting the order $\pi: x_{1}<x_{2}<x_{3}$, it is shown in Fig. 3a that $h$ must be "transported" to the intended level 3 in the BDD, causing references to change, and $v_{2}$ and $v_{3}$ to become intermediate. If, taking commutativity into account as shown in Fig. 3b, $f^{\prime}=x_{3}+x_{2}$ is constructed first and then combined with $h^{\prime}=x_{1}$, no Inode results since there is no recursive descent and $h^{\prime}$ can be "docked" directly. This observation applies analogously to logical conjunction $\operatorname{ITE}(f, h, 0)$ because of the principle of duality mentioned in Section II-A, i.e. only the edge redirections need to be swapped. For negation it is sufficient to swap only the 1 -edges and 0 -edges.
Remark. Binary operators, by themselves, do not perform intermediate computations nor create unnecessary nodes. Thus, it makes no difference for this investigation whether ITE or related algorithms such as $A P P L Y$ [2] construct BDDs.

Generally, the order in which the operands are processed can drastically affect the memory and thus the runtime since the ITE recursion terminates in a specific branch depending on the combination for an operand: for example, for conjunction the value 0 applies, for disjunction if an operand becomes 1.

Another weakness of a conventional BDD manipulation becomes apparent by possible simplifications using the Boolean algebra laws (Section II-A).

Example 3. Let $\rho \in \mathcal{B}_{3}, \rho\left(x_{1}, x_{2}, x_{3}\right)=x_{1} x_{3}+x_{2}+x_{1}$, and $\pi: x_{1}<x_{2}<x_{3}$. Anticipating the absorption law, $\rho$ can be simplified to $x_{2}+x_{1}$. Fig. 4 shows performing $\operatorname{ITE}(f, 1, h)$ based on $f=x_{1} x_{3}+x_{2}$ (Fig. 4a) and $h=x_{1}$ (Fig. 4b). During the recursion steps, it is detected that the node $v_{3}$ already contained in the UT can be absorbed making $v_{5}$ and $v_{6}$ isomorphic and thus mergeable as illustrated in Fig. 4c.

The findings from Example 3 can easily be transferred to other laws such as the resolution laws. An extreme case occurs when, e.g., a BE like $f g+h$ is constructed where $h=1$. Then the construction of $f g$ is wasted since the final BDD is only the 1-leaf. This obviously becomes worse as the BE grows longer. Analogous to the operand order, redundancies can also have negative effects on memory and runtime.

(a) $f$

(b) $h$

(c) $\operatorname{ITE}(f, 1, h)$

Fig. 4: Absorption detection during $\operatorname{ITE}(f, 1, h)$ manipulation, where $f=x_{1} x_{3}+x_{2}$ and $h=x_{1}$

Algorithm 2: Preprocessing approach Sortify to reduce the number of Inodes during BDD manipulation

```
Input: BE \(f\)
Output: Sorted simplified BE based on \(f\)
    \#pragma omp parallel
    \#pragma omp master
    \(\operatorname{sort}\left(f_{\text {begin }}, f_{\text {end }}\right)\)
    simplify \((f)\)
```

In summary, the operand order should already be observed before the BDD manipulation and redundancies should be discovered as early as possible in order to reduce the number of Inodes.

## B. Proposed Approach To Reduce the Number of Inodes

BDDs are constructed on the basis of combining single nodes during BDD manipulation. Due to our observations from the last section, (unnecessary) Inodes emerge during BDD manipulation for two main reasons: 1) poorly selected operand order and 2) redundancies.

As explained in Section II-B, intermediate computations are completely prevented in [15]. However, there is usually a high rebirth rate using well-known techniques such as model checking [11], i.e. dead nodes that may be intermediate are often reactivated before a GC which can prevent deep recursive descents due to caching. In general, it is difficult to predict the "value" of a created node [19]. Therefore, Inodes are permitted up to a certain threshold value in [16]. However, due to the direct integration into the BDD manipulation there are problems with the dynamic variable ordering.
We propose a parallel preprocessing approach called Sortify, using w.l.o.g. OpenMP [27], as a heuristic to reduce the number of Inodes during BDD manipulation as shown in Algorithm 2. To this end, we are oriented to Quicksort [28], one of the fastest sorting algorithms. To address issue 1), a BE is sorted using a predefined sort key invoked by a master thread of a parallel region (Lines $1-3$ ). To address issue 2), the sorted BE is searched neighbor by neighbor to simplify (Line 4).

Algorithm 3: Sorting method sort for Boolean subexpressions

```
Input: BE iterators \(f_{\text {begin }}, f_{\text {end }}\)
Output: Sorted BE \(f\)
    \(\ldots \quad \triangleright\) terminal cases
    \(p \leftarrow \operatorname{median}\left(f_{\text {begin }}, f_{\text {begin }}+\left(f_{\text {end }}-f_{\text {begin }}\right) / 2, f_{\text {end }}\right)\)
    \(f_{\text {left }} \leftarrow f_{\text {begin }}\)
    \(f_{\text {right }} \leftarrow f_{\text {end }}\)
    \(i \leftarrow f_{\text {begin }}+1\)
    while \(i \leq f_{\text {right }}\) do
        if \(k_{\pi}(i, p)\) then
            \(\operatorname{swap}\left(f_{\text {left }}, i\right)\)
            \(++f_{\text {left }}\)
            \(++i\)
        else if \(k_{\pi}(p, i)\) then
                \(\operatorname{swap}\left(f_{\text {right }}, i\right)\)
                \(--f_{\text {right }}\)
        else \(++i\)
        end if
    end while
    if \(\delta\left(f_{\text {begin }}, f_{\text {end }}\right) \geq c\) then
        \#pragma omp taskgroup
        \{
        \#pragma omp task
        if \(\delta\left(f_{\text {begin }}, f_{\text {left }}\right)>0\) then
            \(\operatorname{sort}\left(f_{\text {begin }}, f_{\text {left }}-1\right)\)
        end if
        \#pragma omp task
        if \(\delta\left(f_{\text {right }}, f_{\text {end }}\right)>0\) then
                \(\operatorname{sort}\left(f_{\text {right }}+1, f_{\text {end }}\right)\)
        end if
        \}
    else \(\quad \triangleright\) serial sorting
    end if
```

The main idea (Algorithm 3) is to divide a BE into three subsequent partitions consisting of subexpressions that are less, equal to, or greater than a preselected subexpression $p$ of the entire BE picked by the median of three based on the first, middle, and last subexpression (Lines 1-2). Primarily, there is now a single pass through each subexpression, from left to right (Lines 3-6). Each subexpression is compared to $p$, with the sort key $k_{\pi}$ following the variable order $\pi$. There are three main cases that are handled: If a subexpression is less than $p$, this subexpression is swapped to the left partition (Lines 7-10), otherwise if a subexpression is greater than $p$, it is swapped to the right partition (Lines 11-13), else nothing is swapped (Lines 14-16). Afterwards, the left and right partitions are recursively sorted in the same way. Since the sorting of the left and right partitions is independent, it can be parallelized by two concurrent tasks. To reduce the number of parallel recursive tasks that are scheduled, the cutoff $c(100,000$ according to [27]) is introduced. Before concurrency it is checked if the current number of subexpressions $\delta$ to be sorted reaches $c$ (Line 17). If this is the case, then the recursive descent is performed in parallel (Lines 18-28), otherwise serial (Lines 29-30).

To simplify the BE, its order is now exploited and, similar to finding prime implicants of the Quine McCluskey method [29], the subexpressions are compared neighbor by neighbor in one "round" to see if Boolean algebra laws can be applied.

Compared to Quicksort, Sortify allows to reduce the bestcase complexity from linearithmic $\mathcal{O}(n \log n)$ to linear $\mathcal{O}(n)$, where $n$ is the number of subexpressions. This is achieved by counteracting unbalanced partitions via the selection procedure and avoiding unnecessary recursive calls using three partitions. Since there is no logic minimization, but the subexpressions are only compared neighbor by neighbor after sorting, this effort is negligible. In addition, Sortify is parallelized and cache-coherent greatly affecting the CPU's cache pipeline.

## IV. Experimental Results

This section summarizes the experiments conducted in order to empirically analyze our approach. While Section IV-A describes the setup used for the evaluations. Section IV-B presents the impact of our approach compared to related work.

## A. Experimental Setup

Our preprocessing approach was implemented in $\mathrm{C}++20$ as w.l.o.g. EDDY [30] was used as BDD package for performance evaluation. To allow a fair comparison, the methods discussed in Section II-B were directly integrated into EDDY. For representative purposes, IWLS-93 benchmark instances were taken from [31]. The initial UT (CT) size was set to $2^{20}\left(2^{18}\right)$ due to the complexity of the instances. Using these instances, peak node usage $N$ - the greatest number of nodes in use at any point during the process lifetime was measured during the BDD manipulation and effects on CPU time $T$ (in sec) were recorded, both reported as node ratio $N R$ and time ratio $T R$. The used variable order follows the order of appearance in the respective file. All evaluations were carried out on a Fedora 28 machine with an Intel Xeon E3-1270 v3 CPU with 3.5 GHz and 32 GB of main memory. For each instance, 10 runs were performed and the average was calculated. The Time Out (TO) was set to 10 min , whereas the Memory Out (MO) was configured to a node limit of 1 M .

## B. Performance Evaluation

The experimental results can be seen in Table I and confirm that our approach meets the objectives of this work. Sortify is able to significantly reduce the number of Inodes for all considered benchmark instances thereby accelerating BDD manipulation due to a lower number of recursive descents, GC calls, UT/CT expansions, and hash collisions. As a preprocessing method for $I T E$, it results in peak node usage being reduced by about $29 \%$ compared to single usage (Conventional). This accelerates the runtime by about $23 \%$. While minimally more nodes are needed compared to Multi-way, Sortify is about 19 times faster due to better cache performance which is a successful compromise. In addition, Sortify is stable in solving these instances: While XTop, e.g., has a higher peak node usage than Conventional to construct the BDD for c5315 due to poorly selected decomposition points, Sortify always uses a lower number of nodes at any point during the process lifetime.

TABLE I: Experimental comparison of Sortify and related work in terms of peak node usage and CPU time

| Instance Name | Conventional |  | Multi-way |  | XTop |  | Sortify |  | Conv./Sortify |  | Multi-way/Sortify |  | XTop/Sortify |  |
| :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: | :---: |
|  | $N$ | $T$ | $N$ | $T$ | $N$ | $T$ | $N$ | $T$ | $N R$ | TR | $N R$ | $T R$ | $N R$ | $T R$ |
| des | 21,449 | 0.52 | 13,455 | 5.88 | 21,316 | 0.52 | 15,209 | 0.45 | 1.41 | 1.16 | 0.88 | 13.07 | 1.40 | 1.16 |
| rot | 14,038 | 0.43 | 9,781 | 4.60 | 13,805 | 0.35 | 10,228 | 0.16 | 1.37 | 2.69 | 0.96 | 28.75 | 1.35 | 2.19 |
| c880 | 15,452 | 0.47 | 10,757 | 2.42 | 13,674 | 0.28 | 11,497 | 0.19 | 1.34 | 2.47 | 0.94 | 12.74 | 1.19 | 1.47 |
| c1355 | 46,238 | 0.65 | 32,768 | 3.98 | 40,233 | 0.52 | 34,105 | 0.37 | 1.36 | 1.76 | 0.96 | 10.76 | 1.18 | 1.41 |
| c1908 | 24,854 | 0.54 | 18,201 | 2.93 | 24,032 | 0.38 | 22,193 | 0.24 | 1.12 | 2.25 | 0.82 | 12.21 | 1.08 | 1.58 |
| c2670 | 322,988 | 4.47 | 167,241 | 86.87 | 259,734 | 4.13 | 184,179 | 3.61 | 1.75 | 1.24 | 0.91 | 24.06 | 1.41 | 1.14 |
| c3540 | 405,544 | 5.39 | 292,834 | 97.84 | 393,867 | 5.20 | 318,223 | 4.52 | 1.27 | 1.19 | 0.92 | 21.65 | 1.24 | 1.15 |
| c5315 | 39,983 | 0.64 | 30,992 | 13.96 | 40,577 | 0.76 | 33,109 | 0.55 | 1.21 | 1.16 | 0.94 | 25.38 | 1.23 | 1.38 |
| c6288-12 | 256,878 | 2.28 | , | TO | 208,887 | 2.13 | 176,092 | 1.83 | 1.46 | 1.25 | - | - | 1.19 | 1.16 |
| c6288-13 | 732,667 | 9.27 | - | TO | 611,420 | 8.81 | 428,977 | 7.02 | 1.71 | 1.32 | - | - | 1.43 | 1.25 |
| c6288-14 | MO | - | - | TO | MO | - | 926,563 | 31.97 | - | - | - | - | - | - |
| Total (1-8) | 890,546 | 13.11 | 576,029 | 218.48 | 807,238 | 12.14 | 628,743 | 10.09 | 1.40 | 1.65 | 0.92 | 18.58 | 1.27 | 1.39 |
| $N$ | Peak node usage |  |  | $T$ | CPU time in sec |  | $N R$ |  | Node ratio |  | $T R$ |  | Time ratio |  |

## V. Conclusion

This paper focused on investigating the number of Inode emergences using Boolean functions in order to reduce them and accelerate BDD manipulation. By observing the variable order and detecting redundancy, our developed approach can significantly reduce the number of Inodes and is on average about $20 \%$ faster compared to the single use of ITE and related work as demonstrated by experiments.

In addition to the consideration of further benchmark instances and study of various parameters, future research will be directed towards integrating our approach directly into the BDD manipulation, e.g., to detect redundancies structurally in order to try to further improve the manipulation.

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