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Efficient ROBDD based computation of common

decomposition functions of multi-output boolean functions *

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Abstract

One of the crucial problems multi-level logic synthesis techniques for multi-output boolean functions $f = (f_1, \ldots, f_m) : \{0, 1\}^n \to \{0, 1\}^m$ have to deal with is finding sublogic which can be shared by different outputs, i.e., finding boolean functions $\alpha = (\alpha_1, \ldots, \alpha_h) : \{0, 1\}^p \to \{0, 1\}^h$ which can be used as common sublogic of good realizations of f_1, \ldots, f_m .

In this paper we present an efficient ROBDD based implementation of this COMMON DECOMPOSITION FUNCTIONS PROBLEM (CDF). The key concept of our method is the exploitation of "equivalences" of the functions f_1, \ldots, f_m which considerably reduces the running time of the tool.

Formally, CDF is defined as follows: Given m boolean functions $f_1, \ldots, f_m : \{0, 1\}^n \to \{0, 1\}$, and two natural numbers p and h, find h boolean functions $\alpha_1, \ldots, \alpha_h :$ $\{0, 1\}^p \to \{0, 1\}$ such that $\forall 1 \leq k \leq m$ there is a decomposition of f_k of the form

$$f_k(x_1, \dots, x_n) = g^{(k)}(\alpha_1(x_1, \dots, x_p), \dots, \alpha_h(x_1, \dots, x_p), \alpha_{h+1}^{(k)}(x_1, \dots, x_p), \dots, \alpha_{r_k}^{(k)}(x_1, \dots, x_p), x_{p+1}, \dots, x_n)$$

using a minimal number r_k of single-output boolean decomposition functions.

1 Introduction

The long term goal for logic synthesis is the automatic transformation from a behavioral description of a boolean function to near-optimal netlists, whether the goal is minimum delay, minimum area, or some combination. Most of the approaches attacking the multi-level logic synthesis problem use gate count as optimization criterion. A survey can be found in [4]. Alternatively, some recent papers [10, 11, 13, 15] propose an approach different from the one addressed above. This approach to multi-level logic synthesis which originates from Ashenhurst [1], Curtis [8], Hotz [9], and Karp [12] is based on minimizing communication complexity. The methods used to reduce communication complexity employ functional decomposition, i.e., given a Paul Molitor

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boolean function $f: \{0,1\}^n \to \{0,1\}$ they are looking for (multi-output) boolean functions α , β , and g, such that $f(x_1,\ldots,x_n) = g(\alpha(x_1,\ldots,x_p),\beta(x_{p+1},\ldots,x_n))$ holds for all $(x_1,\ldots,x_n) \in \{0,1\}^n$ (see Figure 1). If one



Figure 1: Decomposition of a boolean function f: $\{0,1\}^n \rightarrow \{0,1\}$

wants to apply functional decomposition recursively to the decomposition functions α and β , a generalization to multi-output boolean functions is required. Of course, the approaches of the papers above can also be applied to multi-output boolean functions $f = (f_1, \ldots, f_m)$: $\{0,1\}^n \rightarrow \{0,1\}^m$ either by considering multi-output boolean functions as single-output multi-value functions $f': \{0, 1\}^n \to \{0, \dots, 2^m - 1\}$ defined by $f'(x_1, \dots, x_n) = \sum_{i=1}^m f_i(x_1, \dots, x_n) \cdot 2^{i-1}$ or by decomposing each singleoutput boolean function f_i independently of the other single-output functions f_j $(j \neq i)$ and only then testing whether they use some identical decomposition functions by chance. The first method has the drawback that it decomposes each single-output function f_i with respect to the same input partition which can result in poor realizations. It does not take into consideration that there possibly exist single-output boolean functions f_i and f_j such that there does not exist one input partition good for both, f_i and f_j . Furthermore, even in case that there is an input partition good for every single-output function of f, it is unlikely

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that there is a decomposition where function g has much less inputs than f (if m is large enough), i.e., g is not much easier to synthesize than f. The drawback of the second method is clear. In the final analysis, it does not use the potential of reusing subcircuits for different outputs of f.

In [14] the authors presented a multi-level synthesis method for multi-output boolean functions based on communication complexity which avoids both drawbacks. The method can be divided into two steps. In the first step, output partitioning is performed, i.e., $\{f_1, \ldots, f_m\}$ is partitioned into disjoint sets Y_1, \ldots, Y_u . Single-output functions f_i and f_j of the same set Y_k will be decomposed with respect to the same input partition. The partitioning is executed such that for every Y_k there is an input partition which is 'near-optimal' for every $f_i \in Y_k$. In the second step, the decomposition functions of the singleoutput functions of each class Y_k are constructed giving special attention to generate these functions in such a way that many of them can be used in the decomposition of different elements of Y_k . Benchmarking results of circuits taken from 1991 MCNC benchmark set have shown the technique to be effective with respect to layout size and signal delay.

The paper in hand presents a much more efficient version of our algorithms from [14]. During the computation of common decomposition functions we efficiently make use of Reduced Ordered Binary Decision Diagrams (ROBDD), which are compact representations for many of the boolean functions encountered in typical applications [6]. In this paper we show that it is possible to carry out all necessary steps based on ROBDD's. In particular we show that the computation of common decomposition functions for the decomposition of several single-output functions, which is the basis of step 2 of the technique above [14], can be performed efficiently based on ROBDD's. (The crucial point of this method is the exploitation of "equivalences" of the functions f_1, \ldots, f_m which results in the computation of connected components (in the graph-theoretical sense)).

Benchmarking results show the new method to be efficient with respect to layout size, signal delay and running time.

We start by giving some basic definitions (section 2) and summarize the algorithm for computing common decomposition functions (section 3). In section 4 we demonstrate how to apply ROBDDs to implement this algorithm. Experimental results close the paper (section 5).

2 Basic definitions

A multi-output boolean function ϕ with *n* inputs is represented as a set $\{\phi_1, \ldots, \phi_m\}$ of boolean-valued output functions. We denote the set of completely defined boolean functions with *n* inputs and *m* outputs by $B_{n,m}$. Let B_n be an abbreviation for $B_{n,1}$. $\phi_{i,\ldots,j}$ $(i \leq j)$ denotes the tuple (ϕ_i, \ldots, ϕ_j) .

Definition 1 A decomposition of a multi-output boolean function $f \in B_{n,m}$ with respect to the input partition $\{X_1, X_2\}$ $(X_1 = \{x_1, \ldots, x_p\}, X_2 = \{x_{p+1}, \ldots, x_{p+q}\}, p+q = n)$ is a representation of f of the form

$$f_k(x_1, \dots, x_n) = g^{(k)}(\alpha_1^{(k)}(X_1), \dots, \alpha_{r'_k}^{(k)}(X_1), \\ \beta_1^{(k)}(X_2), \dots, \beta_{s'_k}^{(k)}(X_2))$$

 $(\forall k \in \{1, \dots, m\}), \text{ where } \alpha_i^{(k)} \in B_p \ (\forall i), \ \beta_j^{(k)} \in B_q \ (\forall j), \ and \ g^{(k)} \in B_{r'_k + s'_k}, \ \alpha_i^{(k)} \ and \ \beta_j^{(k)} \ are \ called \ decomposition \ functions \ of \ f_k \ g^{(k)} \ is \ called \ composition \ function \ of \ f_k \ \diamond$

With respect to a given input partition $\{X_1, X_2\}$, a single-output function f_k can be represented as a $2^p \times 2^q$ matrix $M(f_k)$, the decomposition matrix of f_k or the chart of f_k with respect to $\{X_1, X_2\}$. (For illustration see Figure 2.) Each row and column of $M(f_k)$ is associated with a distinct assignment of values to the inputs in X_1 and X_2 , respectively, such that $f_k(X_1, X_2) = M(f_k)[X_1, X_2]$ where $M(f_k)[X_1, X_2]$ represents the element of $M(f_k)$ which lies in the row associated with X_1 and the column associated with X_2 .

Note that $(\alpha_1^{(k)}, \ldots, \alpha_{r'_k}^{(k)})$ of definition 1 encodes the rows of chart $M(f_k)$. Of course, the following property has to hold.

Encoding Property If the row pattern of row $(v_1, \ldots, v_p) \in \{0, 1\}^p$ differs from the row pattern of row $(v'_1, \ldots, v'_p) \in \{0, 1\}^p$, then $(\alpha_1^{(k)}, \ldots, \alpha_{r'_k}^{(k)})$ has to assign different codes to (v_1, \ldots, v_p) and (v'_1, \ldots, v'_p) .

The minimum number of communication wires required between the subcircuit which encodes the rows of $M(f_k)$ and the composition function $g^{(k)}$ is $\lceil \log p_1^{(k)} \rceil$ where $p_1^{(k)}$ is the number of distinct row patterns in $M(f_k)$. r_k will denote value $\lceil \log p_1^{(k)} \rceil$ in the following. *

Definition 2 A decomposition of $f_k : \{0,1\}^n \to \{0,1\}$ is optimal (for X_1 with respect to a given input partition $A = \{X_1, X_2\}$) if it uses only r_k (= $\lceil \log p_1^{(k)} \rceil$) decomposition functions $\alpha_1^{(k)}, \ldots, \alpha_{r_k}^{(k)}$ with domain X_1 .

To compute decomposition functions (with domain X_1) of a multi-output function f which are used by different single-output functions f_k , we have to consider the following problem which will be denoted by CDF. (CDF can be posed for X_2 in an analogous manner.)

Given: Let $f = \{f_1, \ldots, f_m\} \in B_{n,m}$ be a multi-output boolean function, $A = \{X_1, X_2\}$ with $X_1 = \{x_1, \ldots, x_p\}$ and $X_2 = \{x_{p+1}, \ldots, x_n\}$ be an input partition, and h be a natural number with $h \leq r_k$ (= $\lceil \log p_1^{(k)} \rceil$) ($\forall k$).

Find: *h* single-output boolean functions $\alpha_1, \ldots, \alpha_h \in B_p$, which can be used as decomposition functions of every

^{*}Likewise, the minimum number s_k of interconnections between $\beta^{(k)}$ and $g^{(k)}$ is $\lceil \log p_2^{(k)} \rceil$ where $p_2^{(k)}$ is the number of distinct column patterns in $M(f_k)$.

1									
/									
		X,	00	1		X,	00	1	
		x ₅ 4	00	1		x ₅ ⁴	00	1	
	f₁	5			f	5			
	1	x _n	01	1	2	x _n	01	1	
	X4XaXa				X4XaXa				
	000		row p	attern 1	000		row pattern 1		
	001		row p	attern 1	001		row	pattern 2	
	010		row p	attern 2	010	0 1 0 row pattern 3			
	011		row p	attern 2	0 1 1 row pattern 4			pattern 4	
	100		row p	attern 3	100	row pattern 2			
	101		row p	attern 3	101	0 1 row pattern 2			
	1 1 0		row p	attern 3	1 1 0		row	pattern 1	
	1 1 1		row p	attern 2	1 1 1		row	pattern 4	
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Figure 2: Charts $M(f_1)$ and $M(f_2)$ of the multi-output boolean function $\{f_1, f_2\}$ which will be used to illustrate the ideas of the paper. The input partition A is given by $(\{x_1, x_2, x_3\}, \{x_4, \ldots, x_n\})$. Each chart obviously consists of 8 rows. A row pattern is associated to each row. There are three (four) different row patterns in $M(f_1)$ $(M(f_2))$ denoted by the numbers 1, 2, and 3 (1 to 4). Thus, $r_1 =$ $r_2 = 2$ holds.

single-output function f_k for k = 1, ..., m such that there is an optimal decomposition of f_k of the form

$$f_k(x_1, \dots, x_n) = g^{(k)}(\alpha_1(X_1), \dots, \alpha_h(X_1), \\ \alpha_{h+1}^{(k)}(X_1), \dots, \alpha_{r_k}^{(k)}(X_1), X_2).$$

Of course, such h boolean functions need not to exist. We have proven the problem whether such functions $\alpha_1, \ldots, \alpha_h$ exist to be NP-complete. Nevertheless, we have to solve CDF. An algorithm which is applicable from the practical point of view (as shown by the benchmarking results) is presented in the next two sections.

3 An algorithm for CDF

In the following, let $f = \{f_1, \ldots, f_m\} \in B_{n,m}$ be a multioutput boolean function. Each single-output function f_k has to be decomposed with respect to the same given input partition $A = \{X_1, X_2\}$ (computed during output partitioning (see [14])) which will be fixed in the following.

3.1 Theoretical background

We start with a theoretical result working towards a solution to CDF. It gives a condition necessary and sufficient that h single-output functions $\alpha_1, \ldots, \alpha_h \in B_p$ are common decomposition functions of f_1, \ldots, f_m . It is a generalization of a lemma shown by Karp [12]. For this, we need the following notations. The rows of chart $M(f_k)$ induce a partition of $\{0, 1\}^p$ into equivalence classes $K_1^{(k)}, \ldots, K_{p_1^{(k)}}^{(k)}$ such that $v, v' \in \{0, 1\}^p$ belong to the same class $K_j^{(k)}$ if and only if the two corresponding row patterns of $M(f_k)$ are identical. We denote the corresponding equivalence relation by \equiv_k and the set of the equivalence classes $\{K_1^{(k)}, \ldots, K_{p_1^{(k)}}^{(k)}\}$ by $\{0, 1\}^p/_{\equiv_k}$. Let

 $\theta^{(k)}:\{0,1\}^p\to\{1,\ldots,p_1^{(k)}\}$ be the function which maps $v\in\{0,1\}^p$ to the index j of the class $K_j^{(k)}$ to which it belongs.

Furthermore, for given $\alpha_{1,...,h}$ [†] and all $a \in \{0, 1\}^h$, let $S_a^{(k)}$ be the set $\{\theta^{(k)}(v); \alpha_{1,...,h}(v) = a\}$ of those classes which contain a row mapped to a by $\alpha_{1,...,h}$. $(\alpha_{1,...,h}$ is not able to tell these rows apart (see the Encoding Property).) Note that $S_a^{(k)}$ and $S_{a'}^{(k)}$ need not to be disjoint for $a \neq a'$, and that the number $|S_a^{(k)}|$ of elements of $S_a^{(k)}$ equals the number of distinct row patterns of $M(f_k)$ mapped to a by $\alpha_{1,...,h}$. Thus, $\max\{|S_a^{(k)}|; a \in \{0,1\}^h\}$ denotes the 'inability to distinguish' of $\alpha_{1,...,h}$ with respect to f_k . $itd(A, f_k, \alpha_{1,...,h})$ will denote value $\max\{|S_a^{(k)}|; a \in \{0,1\}^h\}$ in the following.

Continued example: Let h be equal 2. Assume that $\forall (v_1, v_2, v_3) \in \{0, 1\}^3 \alpha_1(v_1, v_2, v_3) = v_2$ and $\alpha_2(v_1, v_2, v_3) = v_3$. It follows that $S_{00}^{(1)} = \{1, 3\}$ holds because $\alpha_{1,2}(000) = \alpha_{1,2}(100) = 00$ and $\theta^{(1)}(000) = 1$, $\theta^{(1)}(100) = 3$. Furthermore the following equalities hold: $S_{01}^{(1)} = \{1, 3\}, S_{10}^{(1)} = \{2, 3\}, \text{ and } S_{11}^{(1)} = \{2\}$. Thus the inability to distinguish $itd(A, f_1, \alpha_{1,2})$ of $\alpha_{1,2}$ with respect to f_1 is equal to 2.

Lemma 1 $\alpha_1, \ldots, \alpha_h \in B_p$ are common decomposition functions of f_1, \ldots, f_m with respect to A such that there is an optimal decomposition of f_k of the form

$$f_k(x_1, \dots, x_n) = g^{(k)}(\alpha_1(X_1), \dots, \alpha_h(X_1), \\ \alpha_{h+1}^{(k)}(X_1), \dots, \alpha_{r_k}^{(k)}(X_1), X_2)$$

 $(\forall k \in \{1, \ldots, m\})$ if and only if the inability to distinguish $itd(A, f_k, \alpha_{1,\ldots,h})$ of $\alpha_{1,\ldots,h}$ with respect to f_k is $\leq 2^{r_k - h}$ $(\forall k)$.

Proof: Since $(\alpha_{1,\ldots,h}, \alpha_{h+1,\ldots,r_{k}}^{(k)})$ has to assign different values to rows of chart $M(f_{k})$ with different row patterns (see the Encoding Property), $\alpha_{h+1,\ldots,r_{k}}^{(k)}$ has to assign different values to those rows which cannot be told apart by $\alpha_{1,\ldots,h}$. As $\alpha_{h+1,\ldots,r_{k}}^{(k)}$ can produce at most $2^{r_{k}-h}$ different values, the statement of the lemma follows.

3.2 The basic algorithm

CDF can be solved by computing $\alpha_{1,...,h}$ by a (simplified) branch and bound algorithm. The sets $S_a^{(k)}$, which determine the inability to distinguish $itd(A, f_k, \alpha_{1,...,h})$ with respect to f_k , are constructed step by step. In the initialization phase, $\alpha_{1,...,h}(v')$ is set to *undef* for all $v' \in \{0, 1\}^p$,

[†]Remember that $\alpha_{1,\ldots,h}$ denotes the tuple $(\alpha_1,\ldots,\alpha_h)$.

and $S_{a'}^{(k)}$ is set to the empty set for all a' and k. Each time we enter the main loop (step 4 of the algorithm; see Figure 3) there is a $v \in \{0,1\}^p$ and a vector $a \in \{0,1\}^h$ such that $\alpha_{1,\ldots,h}(v')$ is defined for all v' with int(v') < int(v),[‡] and there is no extension of the present function table with $\alpha_{1,\ldots,h}(v) = a'$ and int(a') < int(a) which does not violate the condition of lemma 1. In this step we test whether the condition of lemma 1 is violated if $\alpha_{1,\ldots,h}(v)$ is set to a. If the condition is violated, we have to backtrack if $int(a) = 2^h - 1$, i.e., $a = (1,\ldots,1)$. If $int(a) < 2^h - 1$, we enter the loop once again with a incremented by 1. The sets $S_a^{(k)}$ are updated in each step. (Thus, the numbers $itd(A, f_k, \alpha_{1,\ldots,h})$ of lemma 1 are implicitely updated, too.) For detailed information of the algorithm see the pseudo code shown in Figure 3.

Continued example: Assume h = 1, and remember that $r_1 = r_2 = 2$ holds. The CDF algorithm above computes $\alpha_1 : \{0,1\}^3 \rightarrow \{0,1\}$ defined by $\alpha_1(v) = 1 \iff v \in \{010,011,111\}$ as common decomposition function of f_1 and f_2 . The inability to distinguish with respect to f_k (k = 1, 2) is equal to 2 as $S_0^{(1)} = \{1,3\}, S_1^{(1)} = \{2\}$, and $S_0^{(2)} = \{1,2\}, S_1^{(2)} = \{3,4\}.$

4 A ROBDD based implementation

Assume the function $f = \{f_1, \ldots, f_m\} \in B_{n,m}$ we want to decompose is given by m ROBDDs bdd_1, \ldots, bdd_m and that the ordering of the variables is given by (x_1, \ldots, x_n) (for illustration see Figure 4). Then, the following observations result in an efficient implementation.

4.1 Some observations

The first observation, already made in [7, 13], is that for all $(v_1, \ldots, v_p) \in \{0, 1\}^p$ the row pattern belonging to row (v_1, \ldots, v_p) of $M(f_k)$ equals the function table of the cofactor $(f_k)_{x_1^{v_1}\cdots x_p^{v_p}}$ (with $x_i^0 = \overline{x_i}$ and $x_i^1 = x_i$). Thus the problem of determining the number $p_1^{(k)}$ of different row patterns of $M(f_k)$ is equivalent to the problem of computing the number of different cofactors $(f_k)_{x_1^{\nu_1}\cdots x_p^{\nu_p}}$. The ROBDD of the cofactor $(f_k)_{x_1^{\nu_1},\ldots,x_p^{\nu_p}}$ is given by the sub-bdd of bdd_k whose root is reached by starting at the root of bdd_k and then following the path corresponding to (v_1,\ldots,v_p) . The roots of these cofactors are called *linking* nodes (the shaded nodes in Figure 4). Since f_k is given by a ROBDD, the number of different linking nodes of bdd_k obviously equals the number of different cofactors. The computational complexity of determining the number of different linking nodes is at most linear in the size of bdd_k since it can be determined by traversing bdd_k in a depth first search manner.

```
1. Let S_{a'}^{(k)} = \emptyset (\forall 1 \le k \le m, a' \in \{0, 1\}^h),

\alpha_{1,...,h}(v') = undef (\forall v' \in \{0, 1\}^p),

v = (0, ..., 0) \in \{0, 1\}^p, and

a = (0, ..., 0) \in \{0, 1\}^h.
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2. Let
$$\alpha_{1,\dots,h}(v) = a, /* \alpha_{1,\dots,h}(0,\dots,0) = (0,\dots,0) */$$

 $S_a^{(k)} = \{\theta^{(k)}(v)\} \quad (\forall k).$

3. Increment v.

```
4. Let \alpha_{1,...,h}(v) = a.
    If (\forall k) \mid S_a^{(k)} \cup \{\theta^{(k)}(v)\} \mid \leq 2^{r_k - h}
     /* test whether the condition of lemma 1 is
         not violated */
    then let S_a^{(k)} = S_a^{(k)} \cup \{\theta^{(k)}(v)\} \ (\forall k).
              Increment v.
              Let a = (0, \ldots, 0) \in \{0, 1\}^h.
     else while \alpha_{1,...,h}(v) == (1,...,1)
              do let \alpha_{1,\ldots,h}(v) = undef.
                   Decrement v;
                   {}^{"}S^{(k)}_{\alpha_{1,...,h}(v)} = S^{(k)}_{\alpha_{1,...,h}(v)} \setminus \{\theta^{(k)}(v)\} {}^{"} \forall k.
              od
              a = \alpha_{1,\ldots,h}(v).
              Increment a.
              Let \alpha_{1,\ldots,h}(v) = undef.
    fi
5. If (\forall v \in \{0, 1\}^p) \ \alpha_{1, \dots, h}(v) \neq undef
     then return \alpha_1, \ldots, \alpha_h.
    else if v = (0, ..., 0)
              then return "There is no solution"
              else goto 4 fi
    fi
```

Figure 3: Pseudo code of the algorithm solving CDF. In step 2 of the algorithm, we can set $\alpha_{1,\ldots,h}(0,\ldots,0) = (0,\ldots,0)$ without loss of generality because if there exist h common decomposition functions then there also exist h common decomposition functions with $\alpha_{1,\ldots,h}(0,\ldots,0) = (0,\ldots,0)$. Furthermore, the operation $S_{\alpha_{1,\ldots,h}(v)}^{(k)} = S_{\alpha_{1,\ldots,h}(v)}^{(k)} \setminus \{\theta^{(k)}(v)\}$ in step 4 is somewhat more complex than common set difference: the index $\theta^{(k)}(v)$ is only removed from $S_{\alpha_{1,\ldots,h}(v)}^{(k)}$ if there is no $v' \neq v$ with $int(v') < int(v), \alpha_{1,\ldots,h}(v') = \alpha_{1,\ldots,h}(v)$, and $\theta^{(k)}(v') = \theta^{(k)}(v)$.

The second observation is that encoding the linking nodes of bdd_k with a code of length r_k (which is the logarithm of the number of linking nodes of bdd_k) results in an optimal decomposition of f_k . (Of course this simple approach does not lead to common decomposition functions.) For $1 \leq i \leq r_k$ the corresponding decomposition function $\alpha_i^{(k)}$ is given by substituting the linking nodes of bdd_k by the *i*th bits of the codewords belonging to the linking nodes. The composition function $g^{(k)}$ is given by

 $^{^{\}ddagger}int(y)$ denotes the natural number represented by the boolean vector y

[§]Because of $itd(A, f_k, \alpha_1) = 2$, there obviously exists a decomposition function $\alpha_2^{(k)} : \{0,1\}^3 \to \{0,1\}$ such that $\forall v, v' \in \{0,1\}^3$ $(\alpha_1(v), \alpha_2^{(k)}(v)) \neq (\alpha_1(v'), \alpha_2^{(k)}(v'))$ if the row patterns of chart $M(f_k)$ corresponding to v and v' are different (k = 1, 2).



Figure 4: Continued example: The ROBDDs of f_1 and f_2 . The left (right) outgoing edge of node x_i corresponds to the case $x_i = 0$ ($x_i = 1$).

substituting the part of bdd_k , which corresponds to the variables x_1, \ldots, x_p , by the corresponding code-tree. For illustration see Figure 5.

A decomposition constructed in this way leads to decomposition functions $\alpha_j^{(k)}$ of f_k which assign the same value to all those $(v_1, \ldots, v_p) \in \{0, 1\}^p$ for which the corresponding row patterns of $M(f_k)$ are identical. (Note that till now the decomposition functions $\alpha_1^{(k)}, \ldots, \alpha_{r_k}^{(k)}$ of a single-output function f_k are allowed to assign different values to rows with identical row patterns in the case that the total number of different row patterns is less than 2^{r_k} .)

In order to use this simple method of constructing the decomposition functions and composition functions, the branch and bound algorithm above is modified so that only equivalence preserving decomposition functions are considered. We will name the modified algorithm "ROBDD based branch and bound algorithm".

Definition 3 A decomposition function $\alpha_i \in B_p$ of a boolean function $f_k \in B_n$ is said to preserve equivalences if $\alpha_i(v) = \alpha_i(v')$ holds for every $v, v' \in \{0, 1\}^p$ with $v \equiv_k v' .\diamond$

Thus, common equivalence preserving decomposition functions $\alpha_1, \ldots, \alpha_h$ of f_1, \ldots, f_m have to assign the same value to v and $v' \in \{0, 1\}^p$ whenever there is a $k \in \{1, \ldots, m\}$ such that the rows of $M(f_k)$ corresponding to v and v' have identical row patterns. More formally, let

$$v \sim v' \iff (\exists 1 \le k \le m) \ v \equiv_k v',$$

then the corresponding equivalence relation partitions the rows, i.e. $\{0, 1\}^p$, into equivalence classes E_1, \ldots, E_l such that common equivalence preserving decomposition functions have to assign the same value to each $v \in E_i$. We will denote the set of these equivalence classes by $\{0, 1\}^p / \sim$.

Continued example: Figure 6 illustrates the definition of this equivalence relation. Consider a graph whose vertices are the 8 rows $000, \ldots, 111$. (Note that this graph will not be constructed by the algorithm presented in subsection 4.2.2.) There is an edge between two rows v and v' if the

corresponding row patterns in $M(f_1)$ or $M(f_2)$ are identical, i.e., if $v \equiv_1 v'$ or $v \equiv_2 v'$. (Edges resulting from f_1 are drawn in bold.) This results in a graph whose connected components determine the equivalence classes E_i . There are two classes, namely $E_1 = \{000, 001, 100, 101, 110\}$ and $E_2 = \{010, 011, 111\}$. E_1 is marked by shaded nodes. \diamond



Figure 6: Continued example: Illustration of the definition of the equivalence classes E_i .

Thus, the ROBDD based branch and bound algorithm assigns values to the subsets E_1, \ldots, E_l . Because l mostly is much smaller than 2^p , this approach considerably reduces the running time compared to the original branch and bound algorithm (see subsection 3.2). (The benchmarking results will also show that this reduction of the running time can be achieved without reducing the quality of the circuits constructed.) During the ROBDD based branch and bound algorithm, every time a value $a \in \{0, 1\}^h$ is assigned to an equivalence class E_i by $\alpha_{1,\ldots,h}$, the sets $S_a^{(k)}$ ($\forall k$), which contain the different row patterns of $M(f_k)$ mapped to a by $\alpha_{1,\ldots,h}$, have to be updated by $S_a^{(k)} = S_a^{(k)} \cup SET_i^{(k)}$. $SET_i^{(k)}$ denotes the set $\{j; K_j^{(k)} \subseteq E_i\}$, i.e., the set which consists of the indices (with respect to $\theta^{(k)}$) of the different row patterns of $M(f_k)$ belonging to E_i . Note that the sets $SET_i^{(k)}$ and $SET_i^{(k)}$ are disjoint if $i \neq j$.

Continued example:
$$SET_1^{(1)} = \{1,3\}, SET_2^{(1)} = \{2\}, SET_1^{(2)} = \{1,2\}, \text{ and } SET_2^{(2)} = \{3,4\}.$$

Of course, the ROBDD based branch and bound algorithm requires a preprocessing and postprocessing phase described in the following.

4.2 **Preprocessing steps**

In the first preprocessing step we have to efficiently determine the minimum number r_k of decomposition functions required for a decomposition of f_k ($\forall k$). In the second preprocessing step we construct ROBDDs representing the equivalence classes $K_j^{(k)} \in \{0,1\}^p /_{\equiv_k}$. Then, we determine the ROBDDs of the equivalence classes $\{0,1\}^p /_{\sim}$ corresponding to $\{f_1,\ldots,f_m\}$ and thus also the corresponding sets $SET_i^{(k)}$ the algorithm requires.



Figure 5: Continued example: Optimal decomposition of f_2 by encoding the first linking node by 00, the second by 01, the third by 10 and the fourth by 11. (To make things clear the OBDDs shown are not reduced.)

Note that all these computation steps have to be done without constructing the charts $M(f_k)$. Since we have already explained in section 4.1 how to efficiently determine the minimum number r_k of decomposition functions required for a decomposition of a function f_k , it remains to show how to efficiently compute the equivalence classes.

4.2.1 Computation of the equivalence classes with respect to \equiv_k

As already mentioned, identical row patterns of $M(f_k)$ correspond to the same linking node of bdd_k . Thus every equivalence class $K_j^{(k)}$ is associated to exactly one linking node $n_j^{(k)}$ and vice versa. $K_j^{(k)}$ is given by the set of the paths from the root of bdd_k to linking node $n_j^{(k)}$. Thus, substituting the sub-bdd of bdd_k with root $n_j^{(k)}$ by constant 1 and connecting every edge which leaves the cone of $n_j^{(k)}$ to constant 0 results in a ROBDD, whose ON-set is given by $K_j^{(k)}$. We call this ROBDD $bdd_j^{(k)}$. The cone of node $n_j^{(k)}$ is defined to be the set of those nodes x of bdd_k such that there is a path from x to $n_j^{(k)}$. For illustration see Figure 7.

4.2.2 Computation of the equivalence classes with respect to \sim

We implicitely construct a graph G = (V, E) where the set V of vertices is given by the ROBDDS $bdd_j^{(k)}$ representing the equivalence classes $K_j^{(k)}$. At the end, there is an undirected edge $\{bdd_{j_1}^{(k_1)}, bdd_{j_2}^{(k_2)}\}$ if and only if $bdd_{j_1}^{(k_1)} \wedge bdd_{j_2}^{(k_2)} \neq \emptyset$, i.e., iff their ON-sets are not disjoint (see Figure 8). Obviously, there is a one-to-one relation between the set of the connected components (in the graph-theoretical sense) of G and the set of the equivalence classes $\{0, 1\}^p/\sim$. For every class E_i , there is a connected component CC_i of G such that the logical-or of the ROBDDS $bdd_j^{(k)}$ (for any fixed k) corresponding to vertices of CC_i results in a representation of E_i and vice versa.

Continued example: E_1 is represented by the ROBDD $bdd_1^{(1)} \lor bdd_3^{(1)}$ which is the same ROBDD as $bdd_1^{(2)} \lor bdd_2^{(2)}$. E_2 is represented by $bdd_2^{(1)}$ which is the same ROBDD as $bdd_3^{(2)} \lor bdd_4^{(2)}$.



Figure 8: Continued example: Computation of the equivalence classes $\{0, 1\}^p / \sim$. The connected components of graph G represent the classes E_i .

Note that the algorithm described below does not have to test each pair of ROBDDS $bdd_{j_1}^{(k_1)}$, $bdd_{j_2}^{(k_2)}$ whether their ON-sets are disjoint. Virtually, it performs depth first search on graph G. In each step we compute a connected component which contains a node $bdd_z^{(1)}$ not yet touched by calling procedure $search(bdd_z^{(1)}, 1)$. This procedure recursively constructs ROBDDS $cc^{(1)}, \ldots, cc^{(m)}$. At each moment $cc^{(k)}$ equals the ROBDD representing the logical-or of all the nodes $bdd_j^{(k)}$ of the present connected component which have already been touched. At the end of procedure call $search(bdd_z^{(1)}, 1)$, the equation $cc^{(1)} = \ldots = cc^{(m)}$ holds, and $cc^{(1)}$ represents the connected component computed. The exact implementation of search is shown in Figure 9. Note that, if the ROBDD notcovered is non-



Figure 7: Continued example: Illustration of how to efficiently compute the equivalence classes $\{0,1\}^p/_{\equiv_k}$. (Note that the OBDDs shown still have to be reduced.)

Figure 9: Pseudo code of the algorithm computing the equivalence classes $\{0,1\}^p/\sim$

empty during a step, there is a row v belonging to the present connected component which is not in the ON-set of $cc^{(j)}$ yet, so that the ROBDD $bdd_u^{(j)}$ which describes the set of the rows which have identical row patterns as v in $M(f_j)$ has to be joined to $cc^{(j)}$.

Procedure *search* is called exactly once for every node of G. During the execution of the body of procedure search(b, k) (without the recursive calls) there are one apply operation (see [2, 6]) performing the logical-or of two ROBDDs and at most m - 1 + degree(b) apply operations performing logical-and of two ROBDDs, where degree(b) denotes the degree of vertex b with respect to G. This results in a number of apply operations which is linear in the size of G as $m - 1 \leq degree(b)$ holds. The running time of the remaining operations is linear in the size of the relevant ROBDD.

4.3 **Postprocessing steps**

After the execution of the preprocessing steps, the ROBDD based branch and bound algorithm encodes the equivalence classes as already described.

4.3.1 Computation of the ROBDDs of the decomposition functions

Assume that a single-output function α_i has been found by the ROBDD based algorithm which can be used as decomposition function of f_k . Note that the algorithm does not explicitly assign values to every $v \in \{0, 1\}^p$ but only to the equivalence classes $\{0, 1\}^p/\sim$. As the ROBDD of these equivalence classes are known, we only have to connect by logical-or those ROBDDs whose corresponding equivalence classes are mapped to value 1 by α_i in order to obtain the ROBDD representing α_i .

Continued example: Assume h = 1. The ROBDD based branch and bound algorithm constructs $\alpha_1 : \{0, 1\}^3 \rightarrow \{0, 1\}$ defined by $\alpha_1(E_1) = 0$ and $\alpha_1(E_2) = 1$ as common decomposition function of f_1 and f_2 . Thus the ROBDD of α_1 is given by the ROBDD specifying E_2 , and equals function $\alpha_1^{(2)}$ of Figure 5. \diamondsuit

4.3.2 Computation of the ROBDDs of the composition functions

Once that r_k boolean valued functions $\alpha_1^{(k)}, \ldots, \alpha_{r_k}^{(k)}$ which can be used to decompose function f_k are determined, we have to compute the ROBDD of the corresponding composition function $g^{(k)}$. This is done as (informally) described in section 4.1 (second observation). For illustration see Figure 5 once again. The ROBDD of $g^{(k)}$ can be constructed using the linking nodes $n_j^{(k)}$ and combining these cofactors through the codetree with *if-then-else*-operations of the ROBDD-package [2].

5 Benchmarking results

We have synthesized several examples of the 1991 MCNC multi-level logic benchmark set in order to compare factor, factorII [10, 11], which are communication based multi-level synthesis tools developed at PennState University, and misII [5], sis 1.1 [16] to our tool which uses the CDF algorithm described above as basis.* We will call the ROBDD based implementation of our tool mulopII. The former implementation working on charts will be called mulop.

As running time considerations motivate this paper, we compared the running times of our ROBDD based implementation *mulopII* to those of our former version *mulop*. Table 1 shows gate counts[†] and running times for some of the MCNC multi-level logic benchmark circuits. The running times are measured on a SPARCstation 10/30 (64 MByte RAM). The experiments show that the ROBDD based implementation is much faster than the former version. For these examples mulop has running times up to about 50 CPU minutes while the running time of mulopII is at most a few seconds. In particular, these experiments prove our synthesis tool to be applicable in terms of running time. Although the running times of mulopII are much smaller than those of our former version mulop, in almost all cases the numbers of gates of the computed circuits are not larger.

We compared *mulopII* to *factor*, *factorII*. We ran the experiments with the technology mapping used in [11]. Since the quality of the layouts synthesized by *factorII* approximately equals the quality of the layouts synthesized by *factor* (see [11]), Table 2 only reports the results of the comparison of *mulopII* and *factorII*. (The results concerning *factorII* are those which are published in [11].) Compared to *factorII*, our approach generates realizations with a smaller (or equal) number of gates for almost all circuits. As layout considerations motivate the approach of minimizing communication complexity, we compared layout sizes in the following. A comparison to *factorII* and *factorII* and *factorII* and *factorII* and *factorII* and *factorII* approach generates realizations with a smaller (or equal) number of gates for almost all circuits. As layout considerations motivate the approach of minimizing communication complexity, we compared layout sizes in the following. A comparison to *factorII* with respect to layout size was not possible because *factorII* and

	Num	ıber of	N			
Circuit	inputs	outputs	factorII	mulopH	ratio	
9symm1	9	1	75	52	1.44	
$\mathrm{cm}138\mathrm{a}$	6	8	21	21	1.00	
cm151a	12	2	37	40	0.93	
$\mathrm{cm}162\mathrm{a}$	14	5	80	41	1.95	
$\mathrm{cm}163\mathrm{a}$	16	5	47	36	1.31	
cm82a	5	3	18	14	1.29	
cmb	16	4	33	34	0.97	
decod	5	16	31	32	0.97	
$f51\mathrm{m}$	8	8	107	54	1.98	
$\mathbf{x}2$	10	7	65	51	1.27	
z4m1	7	4	25	21	1.19	

Table 2: Comparison between our tool *mulopII* and *factorII* with respect to the number of gates used. The technology file used is taken from the paper of Hwang et al. Note that it is different from the technology file in the other comparisons.

the layout tool used in the paper of Hwang et al. was not at our disposal. We used the standard cell place and route package wolfe which is integrated in octtools. Table 3 shows the comparison between mulopII, misII and sis 1.1 with respect to layout size. The technology library used consists of the set of the 2-input gates.[‡] For almost two thirds of the benchmark set, our approach dominates (or is as good as) that of misII and sis with respect to layout size. The signal delays of our realizations for more than two thirds of the circuits considered are better (or equal) than those of the realizations synthesized by misII and sis. However, on the other hand the results confirm the observation already made in [10, 11] that some circuits, e.g., cm151a, are not suited for being decomposed with respect to disjoint input partitions. The most dramatic improvement has been obtained for circuit 9symm1 which is a symmetric function. This confirms the approach of searching equivalence preserving decomposition functions. §

6 Conclusion

We have presented a ROBDD based technique of computing common decomposition functions of multi-output boolean functions. This algorithm has been integrated in our multi-level synthesis tool which has been presented in [14] where more details of how the CDF algorithm is integrated can be found. The benchmarking results show that most of the circuits constructed by our synthesis tool are very efficient. They also prove it to be applicable in terms of running time.

^{*} The (maximal) value of parameter h of CDF is determined by logarithmic search. More details of how the CDF algorithm is integrated in the tool can be found in [14].

[†]The library consists of the 2-input gates from *stdcell2_2.genlib* available in *octtools*.

[‡]For the technology file itself see $stdcell2_2.genlib$ available in octtools.

[§]Let $f : \{0,1\}^m \to \{0,1\}^n$ be a boolean function which is symmetric in some variables. Then, each equivalence preserving decomposition function of f is symmetric in these variables, too.

	N	o. of gates		Running time			
Circuit	mulop	mulopII	ratio	mulop	mulop II	ratio	
9symm1	40	45	0.89	1.40 sec	1.23 sec	1.14	
C17	6	7	0.86	0.32 sec	$0.15 \mathrm{sec}$	2.13	
$\mathrm{cm}138\mathrm{a}$	20	18	1.11	1.01 sec	0.18 sec	5.61	
cm151a	48	41	1.17	4.16 sec	1.09 sec	3.82	
$\mathrm{cm}152\mathrm{a}$	34	27	1.26	2.15 sec	0.50 sec	4.30	
$\mathrm{cm}162\mathrm{a}$	46	44	1.05	$350.65 \mathrm{sec}$	3.32 sec	105.62	
$\mathrm{cm}163\mathrm{a}$	38	34	1.12	2923.31 sec	2.35 sec	1243.96	
cm82a	13	13	1.00	0.38 sec	0.21 sec	1.81	
cm85a	42	42	1.00	7.46 sec	3.73 sec	2.00	
cmb	24	29	0.83	1836.13 sec	2.52 sec	728.62	
decod	31	28	1.11	26.15 sec	2.56 sec	10.21	
f51m	64	56	1.14	3.14 sec	1.83 sec	1.71	
majority	9	9	1.00	0.44 sec	0.08 sec	5.50	
parity	15	15	1.00	111.06 sec	$1.37 \sec$	81.07	
z4m1	20	20	1.00	0.66 sec	0.76 sec	0.87	

Table 1: Comparison between the (prototype) ROBDD based implementation of our synthesis tool *mulopII* and the former version *mulop* working on decomposition charts. The technology file consists of the 2-input gates from *stdcell2_2.genlib* available in *octtools*.

	Layout size			rat	io	Signal delay			ratio	
Circuit	misII	sis	mulopII	misII	sis	misII	sis	mulopII	misII	sis
9symm1	917928	1194336	201400	4.56	5.93	26.0	27.6	13.6	1.91	2.03
C17	28800	28800	31744	0.91	0.91	4.2	4.2	4.2	1.00	1.00
cm138a	101528	103896	87480	1.16	1.19	5.8	5.8	6.8	0.85	0.85
cm151a	102528	95312	177712	0.58	0.54	12.6	12.6	16.4	0.77	0.77
cm152a	90360	85536	106704	0.85	0.80	10.0	10.0	13.2	0.76	0.76
cm162a	149736	131976	192000	0.78	0.69	13.8	12.0	13.2	1.05	0.91
cm163a	153272	144008	164416	0.93	0.88	11.0	13.0	10.4	1.06	1.25
cm82a	83104	74784	61600	1.35	1.21	8.2	7.2	7.0	1.17	1.03
cm85a	171584	165456	180000	0.95	0.92	10.2	10.2	11.0	0.93	0.93
cmb	198616	204792	123496	1.61	1.66	14.8	9.4	6.8	2.18	1.38
decod	133496	140448	119496	1.12	1.18	6.2	6.2	5.0	1.24	1.24
f51m	536016	561184	251392	2.13	2.23	51.0	51.0	18.4	2.77	2.77
majority	42200	42200	39168	1.08	1.00	7.8	7.8	6.6	1.18	1.18
parity	96976	99408	96976	1.00	1.03	6.2	5.0	5.0	1.24	1.00
z4m1	176160	156288	103896	1.70	1.50	18.0	16.2	9.8	1.74	1.65
Σ	2982K	$3228\mathrm{K}$	1937 K	1.54	1.67	205.8	198.2	147.4	1.40	1.34

Table 3: Comparison between mulopII, misII, and sis1.1 with respect to layout size, and signal delay.

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