Abstract—High-level synthesis consists of many interdependent tasks such as scheduling, allocation and binding. All tasks in high-level synthesis are NP-complete and the design objectives are in conflict for nature, most of the already proposed approaches are not efficient in the exploration of the design space and not effective in the identification of different trade-offs. For these reasons, genetic algorithms can be considered as good candidates to tackle such difficult explorations. A new algorithm that named Learning Automata Genetic Algorithm (LAGA) is used in this paper to perform scheduling and allocation concurrently. This algorithm is based on the Genetic Algorithm, the difference is that the Learning Process is added to the Genetic Algorithm. This strategy can complete the scheduling and the allocation effectively in the high-level synthesis under certain time and resource constraints. This algorithm is implemented in C language and is tested finally on a number of DSP benchmarks, and the test results then are compared with those obtained from four other different techniques which are commonly used in high-level synthesis. The experimental results show that the high-level synthesis using the LAGA algorithm is very effective, especially under the area constraint.

Index Terms—Genetic algorithms, Learning Automata, High-level synthesis (HLS), Scheduling, design space exploration

I. INTRODUCTION

There is a growing consensus among VLSI designers that one of the most effective methods to handle the complexity of today’s system-on-chip (SoC) designs is to use computer-aided design (CAD) techniques. CAD techniques start with an abstract behavioral or algorithmic description of a circuit and automatically synthesize a structural description of a digital circuit that realizes the behavior. The behavioral description consists of computational operations (additions, multiplications, comparisons, logical operations) and control operations (conditional statements, loops, and procedure calls) [1].
High-level synthesis (HLS) is the process of translating a behavioral description into a hardware implementation at register transfer level [3]. The design specification is usually written as a behavioral description, in a language such as C. The behavioral description is first compiled into an internal representation (such as data flow graphs - DFGs), which are then mapped to the functional units that are selected from the resource library to meet design goals (such as power, area, and performance). This process of transforming a behavioral description into a synthesizable structural description affords a methodology of automatically synthesizing a realizable digital circuit from an abstract algorithmic specification of the design, thus considerably reducing the design cycle time. VLSI designs are multiobjective by nature, since they have to trade-off several conflicting design objectives such as chip area, circuit delays, and power dissipation. The shorter design time using behavioral synthesis allows one to examine many alternative circuit realizations during the design process [4]. Often, the structural specification is divided into a data path comprised of the functional and storage units and a control unit that coordinates the flow of data between the data path elements [5]. Due to the division, high-level synthesis is traditionally divided into data path synthesis and controller synthesis[6]. The primarily focus will be on data path synthesis.

Datapath synthesis can be modeled as the process of searching a complex multidimensional space represented by the set of possible schedules, allocations, and bindings that can realize a given behavioral specification.

As modern VLSI and SoC designs become more complex, a major problem is the extremely large number of possible schedule and allocation combinations that must be examined in order to select a design that meets constraints and is optimal [7]. This process, called design space exploration, is further compounded by the need for shortening design times due to time-to-market pressures. Since an exhaustive search could be prohibitive and an ad hoc design exploration could be inefficient, designers often select a conservative architecture after some experimentation, which often results in a suboptimal design. Given this scenario, there is an acute need for techniques that automate the efficient exploration the large space in a reasonable time, during high-level synthesis of datapaths [8]. Searching a complex space of problem solutions often involves a tradeoff between two apparently conflicting objectives: exploiting the best solutions currently available and robustly exploring the design space. Genetic algorithms (GAS) manage this tradeoff in an intelligent way. GAS have recently been applied successfully to optimization problems in diverse fields, such as standard cell placement [9], searching and machine learning [10] and data path synthesis [11].

The Genetic Algorithm begins with a randomly selected population, and through recurrence of the production of the generation, looks for the best chromosome [12]. The aim of the Genetic Algorithm is to find the best chromosome. The position of genes in each chromosome, in the Genetic Algorithm is random. If we should select the appropriate position of genes, it would be possible to appropriate the nearly optimal answer in fewer generations [13]. The Genetic Algorithm, in fact chooses the best chromosome from among the existing ones, and the positions of the genes of chromosomes are totally random [14]. If it were possible to find the optimal place of the genes of chromosomes, we would be able to find the ideal answer in fewer generations. Through utilizing the advantage of both methods, the proposed algorithm tries to achieve the optimal answer in fewer generations.

In the LAGA algorithm each chromosome is equal to an automaton and each gene equal to an action of an automaton.

In this paper, we use LAGA performing subtasks of scheduling in high-level synthesis, and to trade-off conflicting design objectives the process of scheduling based on DFG with weights. And we set weights for DFG according to the constraints of resource.

The paper is organized as follows. Section II provides a brief review of the related work, with particular attention to the evolutionary approaches. Then, Section III describe in details methodology of LAGA, while Section IV presents the results of the experimental. Section V we summarize the paper and draw conclusions based on our experimental results.

II. RELATED WORK

A. High Level Synthesis Methods

A large number of scheduling and allocation techniques have been developed for HLS over the past two decades. It is well-known that there is a strong interdependence between the HLS subtasks, and there is no clear consensus on their order of execution [18]. Such decision often has a large impact on the quality of solutions found and most of the early HLS systems performed those two subtasks separately, obtaining poor results. In literature, the high level synthesis techniques can be classified into four categories: constructive approaches, iterative transformational approaches, exact approaches and non-deterministic approaches. The constructive approaches operate on one operation or resource at a time until all elements are considered. Important algorithms following this approach, for example for scheduling, include common as-soon-as-possible (ASAP) and as-late-as-possible (ALAP) scheduling, list based scheduling [15], force-directed scheduling [16] and path-based scheduling [17]. The iterative transformational approaches perform continuous refinements to the set of solutions while exact approaches [18] exploit a mathematical formulation of the problem to find the optimal solution, but the execution time of these algorithms grows exponentially with the number of variables and the number of inequalities. Therefore, these methods are impractical for large designs. Several high-level synthesis systems use nondeterministic approaches, and in particular GAs, to perform some or all of the synthesis subtasks. Most of them consider two phases and
problems separately like in [19] where GAs are used to schedule the operation while in [20] they are used to allocate and bind a scheduled graph. In the last years, the design of algorithms for DSE is becoming crucial to consider the effects of all the HLS subtasks.

B. Genetic Algorithm

Among the optimization methods inspired by the living nature, genetic algorithm, which is based on the principles of natural evolution, is considered one the best and most sophisticated [21]. Genetic Algorithm is a non-classic and random search optimization method that deals with the function itself, not its derivations, and is based on the theory of the survival of the fittest, inspire by Darwin's evolution theory, and natural genetics [22]. In this method, search begins from several points in solution space simultaneously and through point to point search. The variables of target function are evaluated and, finally, the point which the most or the least absolute is introduced as the optimal point [23]. Optimization is the most important and function of the Genetic Algorithm. In common optimization methods target function must necessarily be coherent and consistent [24]. In Genetic Algorithm, however, a consistent and devisable function is not needed. In accordance with Genetic Algorithm, a sample from among all decision variables, that affects the function, is regarded as a member, and a certain number of these samples, makes up a set of members[25]. In this method, a set of the population of variables is used in the process of search. As a result, as the chance of creating better variables is boosted, the possibility of finding the absolute or general optimal point is heighted. This quality is specifically suitable for functions sudden changes and possessing several situational optimal points. Complete information concerning Genetic Algorithm is brought in [26].

C. Learning Automata

A learning automaton is an abstract model that randomly selects an action from a set of the finite actions and applies it to the environment. The environment evaluates the selected action and informs the result of its evaluation, by a boosted signal, to the learning automata. By using the selected action and boosted signal, the learning automata results its internal situation and then selects its next action.

We can present the environment by $E=\{\alpha,\beta,\gamma\}$ in which $\alpha=\{\alpha_1,\alpha_2,\ldots,\alpha_r\}$ is the set of inputs, $\beta=\{\beta_1,\beta_2,\ldots,\beta_r\}$ is the set of outputs and $\gamma=\{\gamma_1,\gamma_2,\ldots,\gamma_r\}$ is the set of penalty possibilities. When $\beta$ is a two-member set, the environment is P type. In such environment, $\beta_1=1$ is considered penalty and $\beta_2=0$ reward. In a type Q environment, $\beta$ set of processes an infinite number of members. $ci$ is the possibility of a action's being penalized. Learning automata are directed into two groups: those with fixed structures, those with variable structures [27].

III. SCHEDULE USING LAGA

LAGA algorithm is constructed of two phase. In first phase with use of Genetic Algorithm, the result is optimized and in second phases the obtained results from Genetic algorithm improved using learning automata. In first phase the genetic Algorithm is endeavor to optimize the chromosomes and then the obtained chromosomes are putting into learning automata. Then Learning Automata is focus on Chromosome Genes and finding the most suitable place of Genes in Chromosomes. Figure 2 shows the flowchart of proposed algorithm.

A. Initial Population

At first, P (number of population) random generated chromosome and then all tasks are allocated to genes of chromosomes. Then a number is allocated randomly to all genes. The random allocated number to genes includes 2 concepts:

1. Task priority.
2. Processor’s number, which executes the task.

After allocating random numbers to chromosome genes, the values of the genes are interchanged with the number of 2/N load.

We are going to perform, the shown graph in Figure 3, referring to Figure 4, you can observe how tasks are allocated to interior status in order. Since there are two processors in the system, so odd numbers indicate P1 processor and even numbers indicate P2 processor. If the system includes more than two processors, processor’s number will consist of the result of the allocated number to the number of all processors.
B. Task Execution on Processors

When implementing a program on parallel processors, the data dependence between tasks should also be taken into consideration. In fact, a task cannot be implemented unless all of its parent tasks are implemented. In this section, how tasks are implemented on processors has been described [28]. Each task will be implemented in its relevant processor with regard to the automata of figure 4. From among all ready tasks, one, priority is higher than other tasks is implemented. In case, priorities of two tasks are the same as such other, one is randomly chosen for implementation. Ready task is one whose all parent tasks have been implemented.

For example, by considering the automata of figure 4 and the task graph of figure 3 in the first phase, it becomes evident that only T1 task is ready. In the first phase, then, T1 will be executing. After the execution of T1 task, T2, T3, T4, T5 tasks will be on the ready. In this phase, priority of T5 task is higher than other ready tasks, so it is executed. In the same fashion, all tasks are implemented on their own specific processors [29].

With regard to Figure 4 automata, tasks T1, T5, T2, T4, T6, T7, T9 will be run on P1 processor and T3, T8 on P2 processor as well. Tasks execution order will be as TABLE I.

TABLE I.

<table>
<thead>
<tr>
<th>J</th>
<th>S'</th>
<th>v(j)</th>
<th>J*</th>
<th>TS</th>
</tr>
</thead>
<tbody>
<tr>
<td>0</td>
<td>{1}</td>
<td>v(1)-15</td>
<td>1</td>
<td>S-{1}</td>
</tr>
<tr>
<td>2</td>
<td>{2,3,4,5}</td>
<td>v(2)-11,v(3)-14,v(9)-9,v(5)-17</td>
<td>2</td>
<td>S-{1,5}</td>
</tr>
<tr>
<td>1</td>
<td>{2,3,4}</td>
<td>v(2)-11,v(3)-14,v(9)-9</td>
<td>3</td>
<td>S-{1,5,3}</td>
</tr>
<tr>
<td>3</td>
<td>{2,4}</td>
<td>v(2)-11,v(9)-9</td>
<td>4</td>
<td>S-{1,5,3,2}</td>
</tr>
<tr>
<td>5</td>
<td>{4,6,7}</td>
<td>v(4)-9,v(6)-5,v(7)-3</td>
<td>5</td>
<td>S-{1,5,3,2,4}</td>
</tr>
<tr>
<td>4</td>
<td>{6,7,8}</td>
<td>v(6)-5,v(7)-3,v(8)-2</td>
<td>6</td>
<td>S-{1,5,3,2,4,6}</td>
</tr>
<tr>
<td>7</td>
<td>{7,8}</td>
<td>v(7)-3,v(8)-2</td>
<td>7</td>
<td>S-{1,5,3,2,4,6,7}</td>
</tr>
<tr>
<td>9</td>
<td>{8}</td>
<td>v(8)-2</td>
<td>8</td>
<td>S-{1,5,3,2,4,6,7,8}</td>
</tr>
<tr>
<td>6</td>
<td>{9}</td>
<td>v(9)-7</td>
<td>9</td>
<td>S-{1,5,3,2,4,6,7,8,9}</td>
</tr>
</tbody>
</table>

All tasks, now, will be executed according to table2. It shows task execution order on processors in details.

TABLE II.

<table>
<thead>
<tr>
<th>J*</th>
<th>S</th>
<th>pi</th>
<th>ti - vj + pj</th>
</tr>
</thead>
<tbody>
<tr>
<td>1</td>
<td>P1-[1],P2-[1]</td>
<td>1</td>
<td>T1 - 0+2-2</td>
</tr>
<tr>
<td>2</td>
<td>P1-[1,5],P2-[1]</td>
<td>1</td>
<td>T5 - 2+5-7</td>
</tr>
<tr>
<td>3</td>
<td>P1-[1,5],P2-[3]</td>
<td>2</td>
<td>T3 - 3+3-6</td>
</tr>
<tr>
<td>4</td>
<td>P1-[1,5,2],P2-[3]</td>
<td>1</td>
<td>T2 - 7+3-10</td>
</tr>
<tr>
<td>5</td>
<td>P1-[1,5,3,4],P2-[3]</td>
<td>1</td>
<td>T4 - 10+4-14</td>
</tr>
<tr>
<td>6</td>
<td>P1-[1,5,2,6],P2-[3]</td>
<td>1</td>
<td>T6 - 14+4-18</td>
</tr>
<tr>
<td>7</td>
<td>P1-[1,5,2,6,7],P2-[3]</td>
<td>1</td>
<td>T7 - 18+4-22</td>
</tr>
<tr>
<td>8</td>
<td>P1-[1,5,2,6,7,8],P2-[3,8]</td>
<td>2</td>
<td>T8 - 17+4-21</td>
</tr>
<tr>
<td>9</td>
<td>P1-[1,5,2,6,7,8,9],P2-[3,8]</td>
<td>1</td>
<td>T9 - 31+1-32</td>
</tr>
</tbody>
</table>

C. Fitness Function

In Genetic Algorithm, fitness function determines whether chromosomes are going to stay alive. In the problem of task scheduling, the object is to find a short makespan. Analysis function for scheduling problem is:

\[ \text{eval}(v_k) = 1/f_k, \quad k=1,2,..., \text{pop size} \]

\[ f_k : \text{the makespan resulting from } k^{th} \text{ chromosome.} \]

D. Crossover Operator

Crossover is a technique which produces off-springs when two parents mate together. The parents are selected by binary tournament selection method [30]. In this paper, a novel method for combining chromosomes has been put forward. The combination method used in this paper is a two-point one. First two points are randomly chosen as subclasses, and then their contents and orders are analyzed. For instance, the substring chosen from v1, has a weight order of 1-2-3-4. This weight order is used for changing the subclass chosen by v2. Thus, the 6-13-15-11 is changed to 15-13-11-6 and changes with the weight order of v1 subclass.

WMX algorithm is not one, which changes only the contents of two points selected from two chromosomes, but it also changes the contents of classes according to weight priorities. WMA is comprised of three steps.

Step1: random substring selection for two chromosomes. In figure 5, an example of the step1 of
combining chromosome by using WMA algorithm is displayed.

Step 2: defined genes mapping relation. Such as TABLE III.

| 4  2  1  3 |
| 2  5  17  3 |
| 1  2  3  4 |
| 15  13  11  6 |

Figure 6. Step 2 of WMA algorithm

Step 3: creating two new offspring generation, the result of above example indicated in Figure 6.

E. Mutation Operator

For operating mutation, two Genes are randomly selected from a chromosome and their amounts are changed with each other. The manner of swapping actions values of a chromosome is departed into three steps.

Step 1: automata action status after allocating values randomly.

Step 2: Selection of two random actions.

Step 3: The output for automata after swapping actions randomly.

F. Selection Operator

Selection operator in this paper is as follows: In each step of new population production, (1-p) % of chromosome, which has least amount of FT, are selected and enter the new population directly. The rest of the population is, than produced through combining chromosomes.

G. Reward and Penalize Operators

Since, each chromosome is presented as a learning automaton, in each automaton, after considering the fitness of a gene (either processor or action), which is selected on a random basis, that gene, is duly penalized or rewarded. As a result of rewarding or penalizing a gene, its position in the boundary position of an action, its punishment leads to a change in its action and, in consequence, creation of a new makespan. Deparing on the type of learning automata, reward and penalize operator will be different.

Reward action occurs when the fitness of a task is smaller than threshold.

Fitness of ti is: x/y

x: is the sum of connection cost of all parent and offspring nodes of tj node so that. \[ \sum c(t_i,t_j) \text{ if}(p_i \neq p_j) \]

pti: A processor that i t task is performed on it. pti: A processor that j t task is performed on it.
y: is the sum of costs of all parent and offspring nodes of ti node. \[ \sum c(t_i,t_j) \]
c(t_i,t_j): Communication cost between tj and ti tasks. Threshold rate is equal T/Ntasks

T: Consist of a number of related tasks to ti task that is executed on a processor which ti task is run in it.

Ntasks: The number of all graph tasks.

The more fitness level of tj task tends to zero of the connection cost between processors tends towards zero too. If, therefore, the fitness level of a ti task is equal to zero, It turns out that all related tasks of ti are performed on the same processors. T has a direct relation with x; as T increases x decreases and vice versa.

In case the fitness level of a task is lower or equal to the threshold amount, then the head of the task gets penalized. Two positions are possible when penalizing a head:

1. The head might be in a position other than frontier position. In the case, penalizing makes it less important. How the head of task T7 is penalizes, is shown in Figure 8.

2. The head might be in frontier position. In that case, we look for a head in the graph that has the greatest reduction in the amount of FT when processors (the numbers attributed to heads) are changed. Now if the
found head is in the frontier position, the positions of the two heads are changed with each other and if otherwise, i.e. if the found head is not in the frontier position, First the found head should be moved to its frontier position and then change occurs. Figure 9 shows how T8 task is penalized.

First by observing the task graph in figure 6, results obtained from various algorithms and the proposed algorithm is displayed in figure 10.

In this paper, the performance of the proposed algorithm is compared with well-known definite and indefinite algorithms. At first the proposed algorithm is simulated and evaluated on homogeneous platforms and then evaluated on heterogeneous platforms. Parameters that are used in the hybrid algorithm are shown in TABLE III.

### TABLE III. HYBRID ALGORITHM PARAMETERS

<table>
<thead>
<tr>
<th>Algorithm</th>
<th>Memory</th>
<th>Depth</th>
<th>Mutation Rate</th>
<th>Crossover Rate</th>
<th>Iteration</th>
<th>Population</th>
</tr>
</thead>
<tbody>
<tr>
<td>GA</td>
<td>-</td>
<td>0.2</td>
<td>0.7</td>
<td>40</td>
<td>50</td>
<td></td>
</tr>
<tr>
<td>LAGA</td>
<td>4</td>
<td>0.2</td>
<td>0.7</td>
<td>20</td>
<td>50</td>
<td></td>
</tr>
</tbody>
</table>

IV. EXPERIMENTAL RESULTS

The proposed LAGA-based high-level synthesis system has been implemented in the C language. The extended high level synthesis tool accepts ANSI C programs and generates RTL specifications in verilog.

To perform a qualitative assessment of algorithm of LAGA in the high-level synthesis, it was tested on a number of DSP benchmarks drawn from high-level synthesis literature.

For all the benchmarks tested, the synthesized designs were assumed to operate with a clock period of 20 ns. We used a 0.35- CMOS module library, where ALUs, multipliers, registers, and multiplexers are implemented as hard macro cells (cells having fixed aspect ratio and pin locations). The ALUs have a propagation delay of 6.5 ns, and multipliers have a propagation delay of 15 ns. We assume that the area cost and delay of a pipelined multiplier are the same as those of a nonpipelined multiplier, respectively.

In all the experiments, the size of the GA population was set to 100, the crossover probability was 0.90, and the mutation probability set to 0.20. Each of the GA runs was stopped after 10000 fitness evaluations. Since GA algorithms are stochastic algorithms, ten independent runs with different random number seeds were performed for each of the benchmark problem instances, and the best solution found by the GA in each of the ten runs was recorded.

We compared our results with those obtained from four different scheduling techniques commonly used in high-level synthesis, namely, the GA scheduling, ALAP scheduling, force-directed (FDS) scheduling, and simultaneous scheduling, allocation, and binding (SAM) technique. These scheduling algorithms were tested in a traditional high-level synthesis framework that performs the three synthesis subtasks of scheduling, allocation, and binding independently. The goal of this comparison was twofold: 1) to verify the performance gains from concurrently performing scheduling and allocation, over a traditional synthesis flow that carries out these subtasks independently and 2) to use the performance of these scheduling techniques as a baseline to compare our results. The same benchmarks are used for comparing the results.
For each of the benchmark problems, a design-space exploration was performed by setting different values to the weights (and) corresponding area constraints. Results are shown in TABLE IV and TABLE V.

In TABLE IV, the chip latency in bold indicate the Fastest designs for each areas value. From the table, it can be seen that LAGA algorithm finds better solutions than those of the other four scheduling techniques, for all the benchmarks tested.

The experimental results indicate that LAGA algorithm is very effective in high-level synthesis, especially under the area constraint.

TABLE IV. COMPARISON OF OUR LAGA-BASED METHOD WITH OTHER SCHEDULING ALGORITHMS

<table>
<thead>
<tr>
<th>Benchmark Example</th>
<th>Area Constraint</th>
<th>Chip Latency(ns)</th>
</tr>
</thead>
<tbody>
<tr>
<td>IIR</td>
<td>6.0mm²</td>
<td>62</td>
</tr>
<tr>
<td>FIR</td>
<td>5.0mm²</td>
<td>97</td>
</tr>
<tr>
<td>EWF</td>
<td>4.5mm²</td>
<td>272</td>
</tr>
<tr>
<td>ARF</td>
<td>10.0mm²</td>
<td>76</td>
</tr>
<tr>
<td>DCT</td>
<td>12mm²</td>
<td>56</td>
</tr>
<tr>
<td>FDCT</td>
<td>18mm²</td>
<td>65</td>
</tr>
</tbody>
</table>

TABLE V shows the improvement of the LAGA-based solutions over those of the other four scheduling techniques. The average improvements range from 0.74% (compared to the GA method) to 19.74% (compared to the FDS method).

The experimental results show that the LAGA algorithm to complete scheduling and allocation concurrently in high level synthesis is very effective, especially under area constraint. It can obtain a better delay performance than other algorithms.

TABLE V. INCREASE PROPORTION OF THE LAGA-BASED SOLUTIONS OVER OTHERS

The method is simulated on a number of DSP benchmarks. It can succeed in obtaining optimal solutions. The same problems have been also solved in a general way. The experimental results indicate that LAGA algorithm is very effective in high-level synthesis, especially under the area constraint.

V. CONCLUSION

In this paper, a new method LAGA is used in high level synthesis to deal with scheduling and allocation simultaneously. It can produce area and performance optimized designs. This algorithm utilizes Genetic Algorithm and Learning Automata methods sequentially to search for the mode space. It can find the Solutions quickly by using Genetic Algorithm and Learning Automata sequentially in search process.

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