# A Constructive Genetic Algorithm for Gate Matrix Layout Problems

Alexandre C. M. de OLIVEIRA, Luiz. A. N. LORENA

Abstract--This paper describes an application of a Constructive Genetic Algorithm (CGA) to the Gate Matrix Layout Problem (GMLP). The GMLP happens in very large scale integration (VLSI) design, and can be described as a problem of assigning a set of circuit nodes (gates) in an optimal sequence, such that the layout area is minimized, as a consequence of optimizing the number of tracks necessary to cover the gates interconnection. The CGA has a number of new features compared to a traditional genetic algorithm. These include a population of dynamic size composed of schemata and structures, and the possibility of using heuristics in structure representation and in the fitness function definitions. The application of CGA to CGMLP uses a 2-Opt like heuristic to define the fitness functions and the mutation operator. Computational tests are presented using available instances taken from the literature.

*Index Terms--* Constructive genetic algorithms, Gate Matrix Layout, VLSI layout design.

#### I. INTRODUCTION

Gate Matrix Layout problems (GMLP) are related to onedimensional logic arrays and programmable logic arrays folding [1],[2]. In very large scale integration design (VLSI design), the goal is to arrange a set of circuit nodes (gates) in an optimal sequence, such that the layout area is minimized, i.e., it minimizes the number of tracks necessary to cover the gates interconnection.

Fig.1 shows an example of gate matrix, where the gates are numbered from 1 to 9 and the dots are the connection requests (Fig.1a). A group of gates at the same connection is called net. There are 7 nets in circuit of Fig.1. To connect gates 1,3,4 and 7 of the net 1, it is necessary to cross gates that are not part of this net. Moreover, non-overlapping nets can be placed at the same connection track. To compute the number of tracks to cover all nets, it is enough to verify the maximum of overlapping nets. The number of tracks is an important cost factor of VLSI circuits.

Fig. 1b shows, at the bottom, the number of overlaps at each

gate. A typical instance of the GMLP is composed of an  $Ix\ J$  binary matrix, where I and J are the numbers of nets and gates, respectively. The original matrix is represented in the following by a matrix a where all dots are transformed in ones and the other positions receive zeros, and the derived interconnected matrix will be denoted by b, which is generated by transforming in ones all zeros between leftmost and rightmost ones in matrix a. Fig. 2 shows the matrices a and b related to the matrices in Fig. 1.

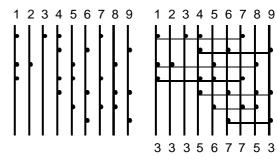


Fig.1. Gate matrix. a) Original matrix; b) Gate matrix derived by interconnection of gates at same net.

101100100	11111100
000101001	000111111
1 1 0 0 1 0 0 1 0	11111110
100110100	11111100
000101011	000111111
000010110	000011110
000001001	000001111
Matrix a	Matrix b

Fig. 2. Derived matrices a and b.

The number of tracks (TR) is easily calculated by:

$$TR = \max_{j \in \{1, \dots, J\}} \sum_{i=1}^{I} b_{ij} = \max\{3, 3, 3, 5, 6, 7, 7, 5, 3\} = 7.$$
 (1)

Besides, there is also the cost of connection of the gates that symbolizes the amount of necessary metal to cover the nets. In this example, the total wire length (*WL*) is:

$$WL = \sum_{i=1}^{J} \sum_{i=1}^{I} b_{ij} - I = (3+3+3+5+6+7+7+5+3)-7 = 35.$$
 (2)

*GMLP* is also a permutation problem, and considering that the initial sequence of gates is  $\{1,2,3,4,...,J\}$ , both TR and WL can be improved by permutated matrices that represent particular sequences of gates. Fig. 3 shows a permutated gate matrix corresponding to the sequence of gates  $\{6,9,4,8,7,5,1,2,3\}$ . The TR is reduced to 5 and the new WL is 22. Nets 3 and 7 can be placed at the same connection track (the same for nets 2 and 6).

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Alexandre C. M. de Oliveira is with the Departamento de Informática, Universidade Federal do Maranhão, São Luís, MA Brasil (e-mail: acmo@deinf.ufma.br).

Luiz A. N. Lorena is with the Laboratório Associado de Computação Aplicada, Instituto Nacional de Pesquisas Espaciais, São José dos Campos, SP Brasil (e-mail: lorena@lac.inpe.br).

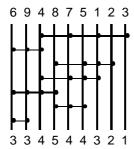


Fig. 3. Permutated gate matrix (optimal solution)

This paper describes the application of a *Genetic Algorithm* (*GA*) to problem *GMLP*. Traditional *GAs* work on a set of variables called structures. For applying them to optimization problems, the first step is the definition of a coding scheme that allows a one-to-one mapping between solutions and structures. A fitness function assigns a numeric value to each member of the current population (a collection of structures). The genetic operators used are selection (like tournament or biased roulette wheel) working together with a number of crossover and mutation operators. The best structure is kept after a predefined number of generations [3]-[5].

Holland put forward the "building block" hypothesis (schema formation and conservation) as a theoretical basis for the *GA* mechanism [4]. In his view, avoiding disruption of good schema is the basis for the good behavior of a *GA*. Schemata can be defined as parts of the structures, where some data are missing, and a major problem is that they are evaluated indirectly, via evaluation of their instances (structures). Goldberg and collaborators have addressed the problem of schemata evaluation and introduced the messy-GA [6]-[8]. In a different approach, Lorena and Furtado [9] proposed recently the *Constructive Genetic Algorithm (CGA)* as an alternative to a traditional *GA* approaches, particularly, for evaluating schemata directly [9].

The *CGA* evolves a population initially formed only by schemata, controlled by recombination, to a population of well-adapted structures (schemata instantiation) and schemata. It was applied to clustering problems [9], location problems and timetabling problems [10]. The *GMLP* is the first permutation like problem modeled and solved using the *CGA* approach.

The *CGA* application can be divided in two phases, the constructive and the optimal:

- The *constructive phase* is used to build a population of quality solutions, composed of well-adapted schemata and structures, through operators as selection, recombination and specific heuristics.
- The optimal phase is conducted simultaneously and transforms the optimization objectives of the original problem on an interval minimization problem that evaluates schemata and structures in a common way.

This paper is organized as follows. Section 2 presents the aspects of modeling for schema and structure representations and the consideration of the *GMLP* as a *bi-objective* optimization problem. Section 3 describes the *CGA* operators,

namely, selection, recombination and mutation, as well as a *CGA* pseudo-code. Section 4 shows computational results using instances taken from the literature.

#### II. CGA MODELING

In this section the modeling phase of the *CGA* is described. The *GMLP* is formulated as a *bi-objective optimization* problem. To attain the objective of evaluating schemata and structures in a common way, two fitness functions are defined on the space of all schemata and structures that can be obtained using a specific representation. An evolution process will be described in the next section. This process considers the two objectives on an adaptive rejection threshold, which gives ranks to individuals in population and yields a dynamic population.

Very simple structure and schema representations are implemented to the *GMLP*. A direct alphabet of symbols (natural numbers) represents the gate sequence and each gate is associated to a column of binary numbers, representing the connection request with each net. The 1 to J values are reserved to gate number and the symbol # is used to express indetermination (# - do not care) on schemata.

Fig.4 shows the representation for the GMLP instance of Fig.1 and a permutation over it, representing structures and an example of schema. The symbols "?" mean there is no information in this column, once the gate number is an indetermination "#".

101100100	100011010	1?0???01?
000101001	000101100	0 ? 0 ? ? ? 1 0 ?
110010010	011000011	0 ? 1 ? ? ? 0 1 ?
100110100	101001010	1?1???01?
000101011	000101101	0?0???10?
000010110	101000001	1?1???00?
$0\ 0\ 0\ 0\ 0\ 1\ 0\ 0\ 1$	000100100	0?0???10?
$s_i = (1\ 2\ 3\ 4\ 5\ 6\ 7\ 8\ 9)$	$s_i = (7\ 2\ 5\ 9\ 3\ 4\ 6\ 1\ 8)$	$s_k = (7 \# 5 \# \# \# 6 1 \#)$

Fig.4. a) Derived matrix (Fig.2a) and corresponding structure  $s_i$ ; b) Example of permutated matrix and corresponding structure  $s_j$ ; c) Example of schema  $s_i$ .

Let *X* be the space of all schemata and structures that can be created by this representation. The *GMLP* is modeled as the following Bi-objective Optimization *Problem (BOP)*:

Min 
$$\{g(s_k) - f(s_k)\}$$
  
Max  $g(s_k)$   
subject to  $g(s_k)$   $g(s_k)$  "  $g(s_k$ 

Function g is the fitness function that reflects the total cost of a given permutation of gates. Commonly, it is used in the formulation that considers track minimization as primary objective and wire length minimization as a secondary one. Therefore, it is defined as  $g(s_k) = I \times J \not R(s_k) + WL(s_k)$ , or

$$g(s_k) = I \cdot J \cdot \max_{j \in \{1, ..., J\}} \sum_{i=1}^{I} b_{ij}$$

$$+ \sum_{j=1}^{J} \sum_{i=1}^{I} b_{ij} - I$$
(4)

where the IJ product is a weight to reinforce the part of the objective considering the maximum number of tracks and to make it proportional to the second part of the objective concerning the wire length. If  $s_k$  is schema, the non-defined

columns (# label) are bypassed. It seems as these columns do not exist and the b matrix used to compute  $g(s_k)$  contains only columns with information. In the example of  $Fig.\ 4.c$ , the track number is max  $\{3,?,4,?,?,?,6,3,?\} = 6$  and the wire length is  $sum\{3+0+4+0+0+0+6+3+0\} - 7 = 9$ .

The other fitness function f is defined to drive the evolutionary process to a population trained by a heuristic. The chosen heuristic is the 2-Opt neighborhood. Thus, function f is defined by:

$$f(s_k) = g(s_v), s_v \in \{s_1, s_2, ..., s_V\} \subseteq \varphi^{2-Opt},$$

$$g(s_v) \le g(s_k)$$
(5)

where  $\mathbf{j}^{2-Opt}$  is a 2-Opt neighborhood of structure or schema  $s_k$ .

As the 2-Opt neighborhood can grow exponentially in function of problem size, in our implementation, only a fixed part, randomly chosen, of this space is used on each function call. Implementation details can be found in Section III.C.

Considering the definition of g, the maximization objective on BOP appears to be a contradiction, but it is needed at the constructive phase that gives distinct treatments to structures and schemata. By definition, f and g are applied to structures and schemata, just differing in the amount of information and consequently in the values associated to them. More information means larger values. In this way, the g maximization objective in BOP drives the constructive phase of the CGA aiming that schemata will be filled up to structures.

The optimal phase is conducted by the interval minimization g–f, which happens on structures near to a local 2-Opt minimum. It is interesting to note that a direct GMLP objective of tracks and wire length minimization is indirectly reproduced at the (g-f) minimization.

#### III. EVOLUTION PROCESS

The BOP defined above is not directly considered as the set X is not completely known. Alternatively is considered an evolution process to attain the objectives (*interval minimization* and g *maximization*) of the BOP. At the beginning of the process, two *expected values* are given to these objectives:

- g maximization: a non-negative real number  $g_{max} > max_{s\bar{1}x} \{g(s)\}$  that is an upper bound on the objective value;
- interval minimization: an interval length  $d_{\mathcal{R}_{max}}$ , obtained from  $g_{max}$  considering a real number  $0 < d \mathfrak{L} 1$ .

The evolution process is then conducted considering an adaptive rejection threshold, which contemplates both objectives in BOP. Given a parameter  $a^{3}0$ , the expression

$$g(s_k) - f(s_k) \stackrel{3}{\longrightarrow} d\chi_{max} - a\chi d\chi_{max} - g(s_k)$$
 (6)

presents a condition for rejection from the current population of a schema or structure  $s_k$ . The right hand side of (6) is the threshold, composed of the expected value to the interval minimization  $d_{\mathcal{R}_{max}}$ , and the measure  $g_{max}$  -  $g(s_k)$ , that shows the difference of  $g(s_k)$  and  $g_{max}$  evaluations.

Expression (6) can be examined varying the value of a. For a=0, both schemata and structures are evaluated by the difference g-f (first objective of BOP). When a increases, schemata are most penalized than structures by the difference  $g_{max} - g$  (second objective of BOP).

Parameter  $\boldsymbol{a}$  is related to time in the evolution process. Considering that the good schemata need to be preserved for recombination, the *evolution parameter*  $\boldsymbol{a}$  starts from 0, and then increases slowly, in small time intervals, from generation to generation. The population at the evolution time  $\boldsymbol{a}$ , denoted by  $P_{\boldsymbol{a}}$ , is dynamic in size accordingly the value of the adaptive parameter  $\boldsymbol{a}$ , and can be emptied during the process. The parameter  $\boldsymbol{a}$  is now isolated in expression (6), thus yielding the following expression and corresponding rank to  $s_k$ :

$$\alpha \ge \frac{d \cdot g_{\max} - [g(s_k) - f(s_k)]}{d[g_{\max} - g(s_k)]} = \delta(s_k).$$
 (7)

At the time they are created, structures and/or schemata receive their corresponding *rank value*  $\mathbf{d}(s_k)$ . These ranks are compared with the current evolution parameter  $\mathbf{a}$ . The higher the value of  $\mathbf{d}(s_k)$ , and better is the structure or schema to the BOP, and they also have more surviving and recombination time.

For the *GMLP*, the overall bound  $g_{max}$  is obtained at the beginning of the *CGA* application, by generating a random structure and making  $g_{max}$  receive the g evaluation for that structure. In order to ensure that  $g_{max}$  is always an upper bound, after recombination, each new structure generated  $s_{new}$  is rejected if  $g_{max} \pounds g(s_{new})$ .

## A. Initial Population

The initial population is composed exclusively of schemata, considering that for each schema, a proportion of random positions receive a unique gate number. The remaining positions receive labels #. Along the generations, the population increases by addition of new offspring generated out of the combination of two schemata.

#### B. Selection

There are two purposes on the evolution process: to obtain structures (good solutions to the *g maximization* objective on the *BOP*), and that these structures be good ones (best solutions to the *interval minimization* objective on the *BOP*). The selection of structures and/or schemata for recombination will be conducted to attain these two objectives, selecting for recombination schemata with small number of labels # and structures or schemata with small

$$d_k = \frac{g(s_k) - f(s_k)}{g(s_k)} \tag{8}$$

The structures and schemata in population  $P_a$  are maintained in ascending order, according to the key:

$$\Delta(s_k) = \frac{(1+d_k)}{\eta} \tag{9}$$

where  $\eta$  is the number of genes containing information (not #). Thus, good individuals with more genetic information

(structures or semi-complete schemata) appear in first order places on the population.

Two structures and/or schemata are selected for recombination. The first is called the *base*  $(s_{base})$  and is randomly selected out of the first positions in  $P_a$ , and in general it is a good structure or a good schema. The second structure or schema is called the *guide*  $(s_{guide})$  and is randomly selected out of the total population. The objective of the  $s_{guide}$  selection is the conduction of a guided modification on  $s_{base}$ . Fig. 5 represents the base-guide selection process.

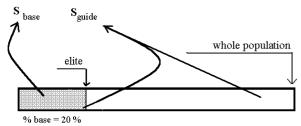


Fig.5. Base-guide selection

#### C. Recombination

In the recombination operation, the current labels in corresponding positions are compared. Let  $s_{new}$  be the new structure or schema (offspring) after recombination. Structure or schema  $s_{new}$  is obtained by applying the following operations:

```
{ Recombination }
For i from 1 to individual length
```

```
if \quad s_{base}(i) = \#
                                and
                                             s_{guide}(i) = \#
                                                                  then
                                                       s_{new}(i) = \#
                                             set
             s_{base}(i) <> \#
II
                                             s_{guide}(i)
                                                                  then
                                  and
                                                       = #
             s_{base}(i) is not
                                                                  then
                                  in
                                                       s_{new}(i) = s_{base}(i)
                                             set
                                                       s_{new}(i) = \#
                                             set
III
            s_{base}(i) =
                                             s_{guide}(i) <> \# then
       if
                            #
                                  and
            s_{guide}(i) is not
                                                                  then
                                                       s_{new}(i) = s_{guide}(i)
                                             set
                                                       s_{new}(i) = \#
                                             set
            s_{base}(i) <> \#
                                             s_{guide}(i) <> \#
                                  and
                                                                 then
            s_{base}(i) is not
                                                                  then
                                             S_{new}
                                                       s_{new}(i) = s_{base}(i)
                                             set
             else
             if s_{guide}(i) is not in
                                             S_{nev}
                                                       s_{new}(i) = s_{guide}
                                             set
                  else
                                                       S_{new}(i) = \#
                                             set
```

Observe that  $s_{base}$  is a privileged individual to compose  $s_{new}$ , but it is not totally predominant. There is a small probability of the  $s_{guide}$  gene information to be used instead of  $s_{base}$  one.

# D. The algorithm

The Constructive Genetic Algorithm can be summed up by the following pseudo-code. The  $\varepsilon$  increment is a linear step that increases the adaptive rejection threshold. Each distinct value of a corresponds to a generation. The stop conditions occur with an emptied population (assured by a sufficiently higher a) or at a predefined number of generations. The population increases, after the initial generations, reaching an upper limit (in general controlled by storage conditions), and decreases for higher values of the evolution parameter a.

```
CGA { Constructive Genetic Algorithm }
Given g_{max} and d; a := 0; e := 0.005;
Initialize P_a;
for all s_k \hat{\mathbf{I}} P_{\mathbf{a}} do compute g(s_k), f(s_k), \mathbf{d}(s_k);
while (not stop condition) do
    while (number of recombination) do
        Select Base and Guide from P_a;
        Recombine Base and Guide;
        Evaluate Offspring;
        Update Offspring in Pa
    end_while
    a := a + e;
    for all s_k \tilde{I} P_a satisfying a > d(s_k) do
        Eliminate s_k from P_a;
    end for
end_while
```

The CGA algorithm begins with the recombination procedures (over schemata only) and the constructive process builds structures (full individuals) progressively at each generation. The constructive process repeatedly uses genetic information contained in two individuals to generate another one. However, the constructive process can be complemented using especially designed *mutation* and *filling* heuristics, searching for a better overall performance. This filling process is called *unary-filling operator* and is applied before the recombination. It consists to fill the  $s_{base}$  substituting the # labels for gate numbers. For GMLP, this feature was implemented as follows:

This code fragment replaces the innermost while loop of the *CGA* pseudo-code. If the selected base is a schema, it is combined with the guide individual (schema or structure) giving a new individual. In any way, local search mutation is applied to a structure (full or filled base) and the offspring are updated depending on its rank. If a best structure is found, it is kept for the end.

For filling schemata it is used an algorithm based on neighborhood minimization. The principle is that a schema  $s_k$  must be filled with a good sequence of gate columns that do not belong to it yet.

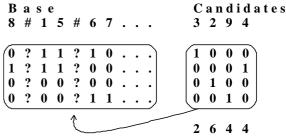


Fig.6. Rule for filling schemata

In the example on *Fig.6*, there are 4 candidate gate columns (on right) to include in a partially showed schema at the marked position (# position). For each # position will be tested all candidate gate columns. It will be chosen the gate column that minimizes the bit-to-bit *xor* relation, considering, in the matrix, the immediate neighbors gate columns to the left and to the right (if exists) of the # point. The bit-to-bit *xor* sum of first candidate column is 2, and is the smaller sum. This mechanism intends to decrease the possibility of zeros between ones at the included gate columns.

The local search mutation is always applied to structures, no matter how they are created (after recombination or after the filling process). The search at 2-Opt neighborhood of the structure (like the f definition on expression 2.3) avoids that the computational effort increases defining a constant number of neighbors to be inspected until the best is found. The neighbors are generated by all the 2-move changes in a constant length part of the structure. A position is chosen at random and an iterative process starts from it, inspecting all possible 2-move changes in the structure until a maximum length previously established. Each 2-move generates a neighbor structure that will be evaluated and the best one will be hold on.

The example of 2-move change is showed in *Fig.*7. The marks in positions of the structures mean reference points to be changed. Non-consecutive references cause the first change type, as showed in *Fig.*7a. Consecutive points cause the second change type in *Fig.*7b. Inspecting all or part of 2-Opt neighborhood needs several moves (catching 2-to-2 points). The number of 2-move changes (neighborhood width) on each local search mutation is a *CGA* parameter of execution to be tuned. This and all the others settings will be described in next section.

The offspring is included in  $P_{\alpha}$  if its rank value d is greater than the adaptive rejection threshold  $\alpha$  and also it is not equal to any other individual in  $P_{\alpha}$ .

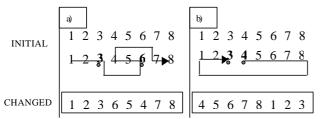


Fig.7 - Examples of one-move in 2-Opt neighborhood; a) non-consecutive reference points change: b) consecutive reference

### IV. COMPUTATIONAL RESULTS

The *CGA* for *GMLP* was coded in *ANSI C* and it was run on Intel Pentium II (266Mhz) hardware. For the computational tests, some *CGA* parameters were adjusted. The *d* parameter was set to 0.15 (usually between 0.10 and 0.20 values, for others applications [9]). This configures the interval  $d_{\mathcal{R}_{max}}$ , establishing the survival time of each individual, once the expected *d* values are proportional to this interval. The *e* was set to 0.005 and also contributes to the survival time of each individual in  $P_a$ . In this case, it is set to slow the increase on

**a**. These parameters avoid the premature termination with an empty population.

Each schema of the initial population received 50% of # genes (indetermination percentage), and 10% up to 20% of population were considered base individuals for base-guide selection, determining a small degree of diversification in selection process. In this case, 10% up to 20% of better individuals of the population are considered for the selection of individuals base.

Local search mutation rate was fixed in 100%, which means a constant improvement of individuals. The number of individuals initially generated was proportional to problem length (at least the number of gates). Other important parameter to be tuned is the neighborhood width (nw) to each local search mutation. After some simulations the better results set nw = 20. The ideal situation would be to use greater values for nw, but this would turn the mutation very slow.

The *CGA* was applied to several instances taken from the literature. Previous tests showed that *CGA* did not found any difficulties on small problems (smaller than v4090 problem), and the tests for some small problems are not reported in this work.

Table I presents the problem circuits considered, along with the corresponding number of nets and gates, a lower bound to the number of tracks, the previous best-known number of tracks (solution) and its reference. Chosen problems were the largest ones found in the literature and others for which authors have obtained recent improvements in results. The lower bound is computed by the maximum column sum of the original matrix. Most of the best previous results comes from Microcanonical Optimization (MCO) approach [11], but other ones with same best results were included as in references: GM\_Learn [12] and GM\_Plan [13]. The MCO overcame (or matched) all the other approaches and it was chosen for comparison with the results of CGA.

TABLE I CHOSEN PROBLEMS TO COMPARE

CHOSEN PROBLEMS TO COMPARE										
gate	net	LB	best	Refs						
			tracks							
10	11	4	4	[11],[13]						
25	17	4	8	[11],[13]						
17	10	5	5	[11]						
16	13	5	5	[11],[12]						
27	23	9	10	[11],[12]						
47	37	5	9	[11],[12]						
48	40	6	11	[11],[12]						
21	18	4	4	[11],[13]						
33	48	14	14	[11],[13]						
70	84	11	18	[11]						
141	202	18	27	[11]						
	gate  10 25 17 16 27 47 48 21 33 70	gate net  10 11 25 17 17 10 16 13 27 23 47 37 48 40 21 18 33 48 70 84	gate         net         LB           10         11         4           25         17         4           17         10         5           16         13         5           27         23         9           47         37         5           48         40         6           21         18         4           33         48         14           70         84         11	gate         net         LB tracks           10         11         4         4           25         17         4         8           17         10         5         5           16         13         5         5           27         23         9         10           47         37         5         9           48         40         6         11           21         18         4         4           33         48         14         14           70         84         11         18						

Table II presents the comparison among the best results of MCO. The MCO was run initially for 10 replications, but only after 1000 replications the authors reached their best results for problems wli, v4000, v4470, x0, w3 and w4. The CGA best results were obtained after 10 replications at same hardware environment. One replication means a complete execution

over the same problem with the same parameters settings. For the *CGA*, the stop condition of one replication is when the best-known solution is reached or at a pre-defined number of generations with no improvement.

TABLE II
CGA COMPARISON TO MCO

CGA COMPARISON TO MCO										
		MCG	)				CGA			
Prob	10 Repl: Tracks	1000 Repl: Tracks	one trial time	1000 trial time	10 Repl: Tracks	wire length	Gen. (avg)	One trial time (avg)	10 trial time	
wli	5.00	4.00	(avg) 0.01	10.0	4.00	24	5.00	0.50	5.00	
wsn	8.00	-	0.01	10.0	8.00	98	7.00	1.50	15.00	
v4050	5.00	-	0.01	10.0	5.00	41	5.00	0.50	5.00	
v4000	6.00	5.00	0.01	10.0	5.00	53	5.00	0.50	5.00	
v4470	10.00	9.00	0.70	700.0	9.00	246	33.00	66.50	665.00	
v4090	10.00	-	0.10	100.0	10.00	95	13.50	2.03	20.33	
x0	11.00	11.00	0.70	700.0	11.00	303	92.57	75.56	755.60	
w1	4.00	-	0.01	10.0	4.00	39	5.00	1.00	10.00	
w2	14.00	-	0.40	400.0	14.00	235	19.50	18.50	185.00	
w3	21.00	18.00	3.90	3900.0	18.00	677	186.00	306.25	3062.50	
w4	32.00	27.00	61.70	61700.0	27.00	1730	225.00	5224.67	52246.67	

To compare different number of replications with respect to execution time, the total time of all the replications was computed based on average of one trial execution time. Thus, the total time of CGA experiments is 10 times the average of one trial execution. By its turn, the total time of MCO experiments is 1000 times the average one trial execution. The "-" execution time, reported in [10] as close to zero, to our computations, was turned to 0.01 seconds, that represents the smallest fraction of one second with two decimal digits. The "-" symbol in the "tracks: 1000 replications" column means that MCO was not run 1000 times. According to the authors, for these problems, it was not necessary to improve the results obtained with 10 replications. The "gen" column is the number of generations (average) in 10 replications of CGA. The numbers of tracks columns refer to the best result found for each problem. An additional CGA column presents the wire length values, not reported in other references. Observing Table II, the CGA found the same results of MCO, but the total time of all replications to find the best results is smaller.

In [10], the authors remark the frequency of 36.3% to find the best result for *wli* problem (relatively small) in 1000 replications. No remarks are reported about larger problems. In our experiments, this percentage is 100% in 10 replications. In *CGA*, smaller percentages of best tracks were found only in larger problems (w3 and w4). There are no available data for a complete comparison of these two approaches, but, at least, *CGA* seems to be more robust than *MCO* in small problems, like the *wli* problem.

The *Table III* shows the *CGA* 5 best results (wire length and tracks) obtained in 10 replications. Problems *w3*, *w4*, *x0 v4470* have less than 100% of frequency in reaching their best solution. The w4 problem appears to be the most difficult one, but these results show that no more than 10 trials were necessary to *CGA* to find the best-known solution.

TABLE III

FIVE BEST RESULTS AND FREQUENCIES FOUND FOR BEST TRACKS											
CGA 5 best solutions											% best tracks
wsn	104	8	104	8	105	8	107	8	113	8	100.00
wli	18	4	18	4	18	4	18	4	18	4	100.00
v4050	41	5	41	5	41	5	42	5	42	5	100.00
v4000	53	5	53	5	53	5	54	5	55	5	100.00
v4470	246	9	253	9	265	9	266	9	282	9	100.00
v4090	95	10	96	10	96	10	98	10	99	10	90.00
x0	303	11	304	11	305	11	306	11	308	11	80.00
w1	39	4	39	4	39	4	39	4	43	4	100.00
w2	235	14	236	14	267	14	273	14	275	14	100.00
w3	677	18	687	18	834	18	843	18	717	18	50.00
w4	1730	27	1836	27	1849	27	1745	28	1745	28	30.00

#### V. Conclusion

This work describes an application of the *Constructive Genetic Algorithm* (*CGA*) to *Gate Matrix Layout problems* (*GMLP*). The *CGA* adapted to work with *GMLP* presents some new specific features, like the filling operator, and a 2-*Opt* heuristic used as mutation and on definition of the two fitness functions (*f* and *g*). Regarding the computational tests, the *CGA* reached all the best-known results (number of tracks) for instances taken from the literature, but it appears to be more robust than the *Microcanonical Optimization* approach. There is no published information about wire length in literature to completely compare the approaches. Applications of *CGA* to other classes of permutation problems are foreseen for future works.

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