

# Multi-objective Genetic Algorithms for Courses of Action Planning

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**Abstract - Planning military courses of action is a very complex and difficult activity. Planners should take into consideration environmental information, predictions, the end state targeted and resource constraints. Development of courses of action involves solving simultaneously planning and scheduling problems. In this work, a new approach based on genetic algorithms (GA) and multi-objective optimisation is proposed to support resource-constrained courses of action development where both cardinal and ordinal objectives are considered. A vector of fitness evaluations is proposed to control the proportion of the infeasible solutions. Crossover and mutation operators are designed to diversify the search space and improve solutions on all objectives from one generation to another. In the replacement strategy, a selection procedure, based on the dominance concept and a multi-criteria filtering method, is proposed. Such a strategy is applied when the population reaches a critical size. Different GA schemes are compared and their strengths and weaknesses are discussed. The multi-criteria filtering procedure used in the replacement strategy proved very efficient in the diversification of the Pareto front.**

## 1 Introduction

The Course of Action (COA) Development step of the Military Operational Planning Process (OPP) involves the entire staff. The Commander's guidance and intent helps the staff to focus on the development of comprehensive and flexible plans within the time available. These COAs "should answer the fundamental questions of when, who, what, where, why and how" [CFC Toronto, 2000, US Army, 1997]. Each COA should be suitable, feasible, acceptable, exclusive and complete. A good COA positions the force for the future operations and provides flexibility to meet unforeseen events during its execution. The "who" in a COA does not specify individual units, but rather uses generic assets and capabilities. During the development step, the staff analyses the relative combat power of friendly and enemy forces, and generates comprehensive COAs.

During the mission analysis, the staff should identify the assigned and implied tasks to perform the mission.

These tasks can be decomposed into sub-tasks. Tasks and sub-tasks can be represented by means of a hierarchical structure called work break-down structure. Leafs of this hierarchical structure are called elementary tasks. Synchronization analyses lead to identify temporal and spatial relationships between elementary tasks (e.g. End-Start, Start-Start, End-End, Time Laps, Same Spatial Zone...). Staff should then consider all available resources and capabilities and assign them to the tasks. Synchronizing COA requires scheduling starting and ending times of all tasks according to resource availability, deployment constraints and task relationships. Any resource or capability has an availability calendar, in-use costing, required preparations, required staffing, etc...

In summary, the challenge for the planning managers is to generate complex, spatially and temporally interdependent activities with precedence relationships, subject to resource constraints, and satisfying multiple incommensurable and often conflicting criteria.

Guitouni et al. (2000, 2002) proposed to model a COA planning as a multiple mode resource-constrained project-scheduling problem (MRCPS) since, from a methodological point of view, planning and scheduling are not much different. The model consists of representing generic activities (tasks with specific combinations of resources) into elementary (or primitive) actions interrelated to accomplish the mission objectives. This process implies the identification of the tasks (when and where) as well as their precedence relationships, the pool of available resources with their localization, and finally the objectives of the mission. A COA is then represented as an oriented time-space graph (see Fig. 1). Depending on the combination of resources allocated and the action's position in the schedule, different COA networks could be obtained. They constitute variants (or alternatives) of a mission with different evaluations on objectives.

Solving COA planning problems is NP-Hard. To obtain promising feasible alternatives with respect to multiple objectives, we thought to explore the potential of evolutionary algorithms (EA), a meta-heuristic that has proven reliable for solving combinatorial NP-hard problems.



state-of-the-art evolutionary methods such as NSGA-II (Deb, 2001), NPGA2 (Erickson et al., 2001) and SPEA2 (Zitzler et al., 2001), based on dominance, elitism and niching or crowding, aim to achieve this two-sided goal. The approach proposed in this work includes these two aspects (convergence towards Pareto frontier and diversification of efficient solutions) by using a first order multi-criteria filter (MFP) to select solutions making the next generation.

### 3 Construction of a GA-based multiple objectives RCPS

Multi-objective COAs could be characterised by a set of tasks, a set of resources, precedence relationships, resources availability constraints and global performance functions (criteria)  $F$ .

The problem formulation states as follows:

$$\begin{aligned} \text{Optimize } F_z, z = 1, \dots, Z & \quad (1) \\ \text{s.t. } t \in D & \quad (2) \\ \text{s.t. } R \in C & \quad (3) \end{aligned}$$

with the vector of tasks  $t = \{t_1, t_2, \dots, t_n\}$  having the following attributes for each task  $t_i$ :

- Starting and ending time  $[td(i), tf(i)]$  considered as integer variables. The earliest and latest starting and ending time, respectively.  $[\tau_s(i), \tau_e(i)]$  are used to generate different initial solutions.
- A localization spatial coordinate  $(x, y, z)$
- Type and quantity of resources required, represented by a set  $\mathbf{R}$  composed of renewable and non-renewable resources available in limited quantities.  $R_k(t_i) = \{r_{1i}, r_{2i}, \dots, r_{mi}\}$  is the  $k^{\text{th}}$  set (or combination) of resources required to accomplish the task  $t_i$ .
- Set of predecessors  $\{PR\}$  characterized by the tasks that temporally and/or spatially precede  $t_i$ .

Resources with the following attributes:

- Starting and ending time of availability  $[t_{rs}(k), t_{re}(k)]$  (resource's timetable)
- Quantity available during this interval of time
- Localization of the resources (depot)  $(x, y, z)$
- Type of resource
- Other specific characteristics such as in-use cost, mean speed (for mobile resources), reliability, etc.

Equation (2) concerns the constraints that ensure that each task is processed once in its time interval and precedence conditions are fulfilled (feasible tasks). Equation (3) express the resource constraints (e.g. availability).

The mGA is used to find different task-resource combination networks where all activities are completed, the resources and precedence constraints satisfied, and the best compromise between criteria reached. The optimisation is carried out using a randomly initialised population for a low size problem (COAs with less than 10 actions). For large size problems (COAs with more than 10 actions), a heuristic method based on the network approach and Cplex is used (Urli et al., 2003). Crossover and mutation are used for the exploration and exploitation of the search space. The principle of the variable

neighbourhood search (VNS) method is used during the optimisation. This is achieved by introducing three binary (crossover, mutation) and one unary (mutation) operators in order to explore a great number of neighbourhoods. Unlike classical GA procedures, the population size is not kept constant from generation to generation but increases by reproduction and mutation until it reaches a critical value determined empirically as explained in section 5.1. A replacement strategy is then applied to select the best candidates for the next generation (survivals). The selection is achieved based on dominance and/or a first order multi-criteria filtering procedure (MFP) detailed in section 3.4 and summarised in Figure 2.

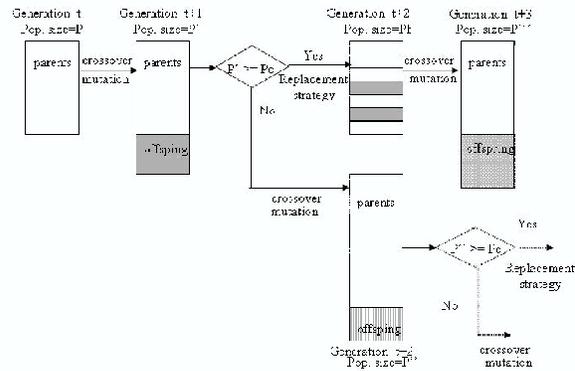


Figure 2. Evolution of the population from generation to generation

#### 3.1 Encoding variables

Each solution, representing a COA network, is encoded in the following form:  $x(i) = [(t_1, R_{1,1}), (t_2, R_{2,4}), \dots, (t_n, R_{n,1})]$ , for  $i = 1, \dots, \text{popsize}$ , where  $\text{popsize}$  is the population size.  $t_j$  is the  $j^{\text{th}}$  task to be scheduled and  $R_{j,k} = R_k(t_j)$  is one of the sets of resources available to accomplish this task.

#### 3.2 Evaluation of a solution

As it is difficult to apply genetic algorithms to a multi-objective constrained problem, one of the options proposed in the literature is to integrate the constraints into the objective functions using penalty functions. However, the aggregation of penalty functions to the objectives in a highly constrained problem may lead to a large proportion of unfeasible solutions and there is no control of this number. In this work, the constraints are considered as functions to be optimised as the objectives and the solutions generated will be retained regarding both their objectives and constraint violation. Thus, to each solution is assigned a fitness vector.

##### 3.2.1 Constraint functions

###### Resource availability

Each resource,  $r_{kj}$ , used by task  $t_j$  must be available, in the desired quantity, during the period  $[t_{rs}(j), t_{re}(j)]$ .

A penalty coefficient  $DP(j)$  associated to each task is introduced and defined by:

$$DP(j) = \begin{cases} 1, & \text{if all the resources to accomplish the task } t_j \text{ are available} \\ 0, & \text{otherwise} \end{cases}$$

For each solution, it follows:

$$fitn_1(i) = \sum_{j=1}^n PP(j) \quad (4)$$

This function is to be maximised (arg [max fitn<sub>1</sub>(i)] = n) to force the search into the feasible region.

*Precedence*

Some tasks have successors and predecessors. Other tasks are free. For each task j, if all its predecessors have been correctly scheduled, we set tPP(j) = 1 otherwise tPP(j) = 0. The precedence penalty is defined as:

$$fitn_2(i) = PP(i) = \sum_{j=1}^n tPP(j) \quad (4a)$$

As with fitn<sub>1</sub>(i), fitn<sub>2</sub>(i) is to be maximised.

### 3.2.2 Objective functions

The objectives considered here could be cardinal (quantitative) such as cost, reliability, make-span, or ordinal (qualitative) such as the impact of a COA. For a uniform optimisation (maximisation), the objectives to be minimised are reformulated and denoted fitn<sub>1</sub>(i), j = 3, ..., Z.

### 3.3 Genetic operators

*Crossover (or recombination)*

Selection of two parents for the crossover is done using the roulette wheel selection. Two candidates are selected using the following procedure repeated popsize times:

- Choose randomly two fitness components fitn<sub>k</sub> and fitn<sub>1</sub> (objectives or constraints)
- Evaluate selection probabilities of the population according to these two fitness components:

$$ps_k(i) = \frac{fitn_k(i)}{\sum_{i=1}^{popsize} fitn_k(i)}, \quad ps_1(i) = \frac{fitn_1(i)}{\sum_{i=1}^{popsize} fitn_1(i)} \quad (5)$$

- Select randomly one candidate per fitness component according to its selection probability.

This selection procedure, favouring one criterion per candidate, aims to produce offspring characterised by the best features present in their parents, and ultimately induces higher population diversity.

Two crossover procedures are proposed, to be used alternatively, in order to explore a greater number of search spaces. The first one is the uniform crossover operator (Syswerda, 1989) which has been shown to be superior to traditional crossover strategies for combinatorial problems. When two chromosomes are selected for crossover, a random mask is generated and their genes are exchanged according to the mask. This mask is simply a binary string with the same length as a COA vector (Sec. 3.1). The parity of each bit determines which genes will be exchanged. The gene in this procedure is represented by an action, (t<sub>j</sub>, R<sub>j,k</sub>).

The second operator is the partial mapped crossover (PMX) proposed by Goldberg and Lingle (1985) and is an

extension of two-point crossover to permutation representation.

A repairing procedure is used, in both operators, to resolve illegitimacy of the offspring if some activities are missing or duplicated. This procedure is achieved by simply transferring these activities from one child to the other one.

*Mutation*

Mutation is applied randomly on the population and the probability of mutation is inversely proportional to the population size as recommended by De Jong (1975). The two operators used for the mutation consist of:

- Exchanging, with a probability p<sub>α</sub>, a randomly selected combination of resources between two COAs. The offspring, which received the combination with the best criterion, is retained.
- Switching the quantity of two resources in a combination of resources associated to a randomly selected task j. This option is used every four generations.

The probability of mutation p<sub>α</sub> of a set of resources R<sub>i,j</sub> in a chromosome is based on a criterion related to the resource characteristics such as cost or reliability. In this way, only the best resource combination is exchanged, concerning the selected criteria. These probabilities are used alternatively from one generation to the other to produce COAs with improved combinations of resources.

### 3.4 Replacement strategy

The solutions generated from the crossover and the mutation operations are evaluated. When the population size attains or exceeds a critical value P<sub>c</sub>, P<sub>f</sub> individuals are selected, based on their fitness vector, among the parents and the offspring using a replacement procedure. Otherwise, new offspring are generated (see Fig. 2). Three strategies are tested for the replacement:

- The multi-criteria filtering procedure (MFP), proposed by Guitouni et al. (2001), returns (a user-defined) P<sub>f</sub> non-dominated diversified individuals. MFP is based on multi-criteria dynamic conjunctive and disjunctive procedures. The retained solutions are characterised by at least one best-scored objective or by all objectives achieving minimal threshold values,
- the non-dominated sorting approach (NDS), as proposed by Srinivas and Deb (1995), but used only to rank the best P<sub>f</sub> individuals,
- the mixed approach: NDS+MFP procedure.

#### 3.4.1 Strategy 1: MFP method

The filtering procedure is performed in two steps using a disjunctive procedure then a conjunctive procedure. Let A be the set of parents and offspring to be filtered and  $\hat{A}$  the set of individuals retained for the replacement. Let Card(A) = P<sub>c</sub> and Card( $\hat{A}$ ) = P<sub>f</sub>. Let  $\underline{e}_0$  be a threshold vector representing the lower limits (lower thresholds) of the objectives imposed to the selection:  $\underline{e}_0 = (e_{01}, e_{02}, \dots, e_{0Z})$ . Let  $\underline{e}_1$  be a threshold vector representing the upper limits (higher thresholds) of the objectives imposed to the selection:  $\underline{e}_1 = (e_{11}, e_{12}, \dots, e_{1Z})$ . S<sub>1</sub>: {fitn<sub>1min</sub>, fitn<sub>2min</sub>, ...}

$\text{fitn}_{Z_{\min}}$  the set of anti-ideal points of A,  $\text{fitn}_{z_{\min}} = \min_{i=1, P_c} \{\text{fitn}_z(i)\}$ ,  $z = 1 \dots Z$ ,  $S_2: \{\text{fitn}_{1_{\max}}, \text{fitn}_{2_{\max}}, \dots, \text{fitn}_{Z_{\max}}\}$  the set of ideal points of A,  $\text{fitn}_{z_{\max}} = \max_{i=1, P_c} \{\text{fitn}_z(i)\}$ ,  $z = 1 \dots Z$ .

#### Disjunctive procedure

This procedure selects individuals characterised by at least one objective having a maximal value. Let  $\hat{A}_d$  be the set of individuals selected by the disjunctive method. The following steps could describe it:

- $\hat{A}_d = \emptyset$
- Compute  $e_{1z}$  for each objective ( $e_{1z} = S_{2z} = \text{fitn}_{z_{\max}}$ ,  $z = 1, \dots, Z$ )
- Select individuals  $i$  that  $\exists z \in \{1, \dots, Z\}$ ,  $\text{fitn}_z(i) \geq e_{1z}$ ,  $i = 1 \dots P_c$
- Add these individuals in  $\hat{A}_d$ .

#### Conjunctive procedure

Candidates characterised by objective values less than the thresholds are discarded. The thresholds are automatically computed using a dichotomist method between  $S_1$  and  $S_2$ . The procedure has the following steps:

- Compute the threshold value  $e_{0z}$  for each objective
- Select individual  $i$  so that  $\forall z$ ,  $\text{fitn}_z(i) \geq e_{0z}$ ,  $i = 1 \dots P_c$
- Add this individual in the set  $\hat{A}_d$  until  $\text{Card}(\hat{A}_d) = \text{Card}(\hat{A})$ .

This multi-objective filtering procedure allows the selection of a diversified population of individuals characterised by the fittest objective or all objectives higher than a threshold value.

### 3.4.2 Strategy 2: NDS method

The solutions are ranked on the basis of non-domination. All non-dominated individuals in the current population are placed at the top of a list and assigned a rank of 1. These solutions are removed from the remaining population and the next set of non-dominated solutions is identified and assigned rank 2. The process is repeated until the entire population is ranked. The top  $P_f$  individuals in the list are then selected for the next generation.

### 3.4.3 Strategy 3: NDS+MFP method

This procedure combines the features of the two previous ones. First, the population is ranked using the NDS method. The set identified with rank 1, *i.e.*, the non-dominated solutions, is selected. Let  $A_1$  be this set.

- If  $\text{card}(A_1) = P_f$ , then these solutions are individuals of the new generation
- If  $\text{card}(A_1) > P_f$ , use the MFP method to return  $P_f$  solutions
- If  $\text{card}(A_1) < P_f$ , select the set of solutions identified with rank 2. Let  $A_2$  be this set
- If  $\text{card}(A_1 \cup A_2) > P_f$ , use the MFP method to return  $P_f$  solutions
- If  $\text{card}(A_1 \cup A_2) < P_f$ , select the set of solutions identified with rank 3, and repeat the process until  $P_f$  solutions are selected.

This strategy differs from NDS because it selects the solutions based on their evaluations besides being non-dominated. The efficiency of these three replacement strategies is examined in section 5.

## 4 Indicators for performance assessment

Different investigations to evaluate the performance of evolutionary algorithms have been proposed, taking into account the optimisers' stochastic characteristics (see for example: da Fonseca et al., 2001) or not (see for example Zitzler et al., 2003) and often assuming that the Pareto front is known. As pointed out by Bosman and Thierens (2003), comparing the performances of multi-objective evolutionary techniques is not an easy task. There are several criteria of finding a good approximation of the Pareto front and most current methods do not outperform each other but work better regarding different performance indicators. In this work, we examine the performance of the proposed mGA strategies using three indicators, assuming that no a priori information is available on the Pareto set, and taking into account the stochastic aspect of the methods. These indicators are the cardinality of the Pareto set  $S$  (Van Veldhuizen, 1999),  $N_s$ , the diversity of solutions in the decision space,  $D_s$ , and the spread of solutions on the Pareto front,  $COV$ .

#### Cardinality of the Pareto set, $N_s$

The non-dominated solutions are stored in an external list updated at each iteration. Thus, the approximation Pareto set  $S$  is represented by this population when the stopping criterion is reached and  $N_s = |S|$ .

#### Diversity of the Pareto set, $D_s$

When designing a GA algorithm, the diversification of the population (solutions) is often the major goal through the exploration and exploitation operators. Ultimately, the diversity of the solution in the Pareto set represents a desirable result as it gives to the decision maker a larger range of options. The diversity of a population used here is inspired from the Shannon theory to evaluate the value of information based on the entropy  $H$ . If the entropy of a message ( $0 \leq H \leq 1$ ) is high, then the redundancy of the information will be low and the value of this message will be high.

Here, the diversity of the solutions is based on the number of alternatives of resource combinations that could be allocated for each task  $i$  in the population. So the total number of COA variants is given by the number of different resource combinations available for all the tasks.

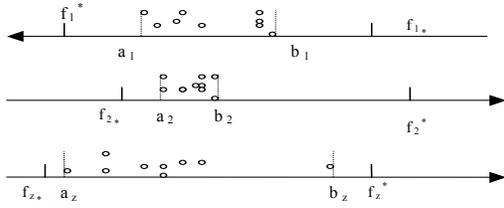
$$D_s = \frac{1}{n} \sum_{i=1}^n H_i \quad (8)$$

$$\text{and } H_i = \frac{1}{\log(q_i)} \sum_{j=1}^{q_i} \frac{n_{ij}}{\text{pop-size}} \log\left(\frac{n_{ij}}{\text{pop-size}}\right) \quad (9)$$

where  $D_s$  is the diversity of the population,  $H_i$  the entropy of a task  $i$ ,  $n$  the number of tasks in the COA,  $q_i$  the number of resource' combinations per task found in the population, and  $n_{ij}$  the number of solutions (COAs) with different combinations of resources for task  $i$ . The higher the number of COAs with different resource configurations per task, the higher the diversity of the approximation set.

#### Diversity of the Pareto front $COV$

We propose to express the diversification of the solutions spread all over the Pareto front by the extent of coverage,



**Figure 3. Spread of solutions over the Pareto front in multidimensional space**

in each objective's dimension separately as illustrated in Figure 3. This indicates the size of the objective space covered by the Pareto set. The extent of coverage relative to each objective  $f_i$  is defined as follows:

$$\text{Cov}_i = \frac{\max_{x,y \in S} |f_i(x) - f_i(y)|}{|f_i^* - f_i^*|} \quad (10)$$

where the numerator expresses the maximum distance among solutions of the approximated Pareto set  $S$  in the dimension  $i$ , and  $f_i^*$  and  $f_i^*$  are, respectively, the anti-ideal and ideal values of  $f_i$  (true extremes of the objectives or its worst and best values by comparing simultaneously all the approximation sets).

Zitzler and Thiele (1998) have proposed a similar indicator, but aggregating the  $\text{Cov}_i$  for  $i=1, \dots, Z$ , by considering the union of all bounding-boxes covered by the Pareto set is misleading. A single, global score for the coverage does not reflect how the diversification is with respect to a given criteria. For a better appreciation of the dispersion, it is more appropriate to consider the coverage array on all the objectives:

$$\text{COV} = (\text{Cov}_1, \text{Cov}_2, \dots, \text{Cov}_Z) \quad (11)$$

Comparing these coverage vectors could be achieved using dominance analysis, multiple criteria decision analysis or statistical techniques.

## 5 Planning a course of action: computational results

In order to evaluate the relative performance of the methods, we first check if an approximation Pareto set dominates another set. This set domination is defined as:

$S_1$  strictly dominates  $S_2$  if  $\forall x \in S_2, \forall y \in S_1,$

$\forall i, 1 \leq i \leq Z, f_i(y) \geq f_i(x)$  and  $\exists k, f_k(y) > f_k(x)$ .

We, then compare the difference between their performance indicators and test their significance by using the Wilcoxon signed-rank statistical test since the data distributions are not necessarily symmetrical.

If no conclusion can be drawn, we can examine which method puts more emphasis on diversity or on getting a high number of efficient solutions.

The efficiency of the method proposed here is investigated using the example of courses of action with four objectives: the cost and the make-span to be minimised (cardinal and linear objectives), the resource reliability and the impact on the enemy, to be maximised. The impact is measured on a qualitative scale (ordinal

objective). Resource availability and tasks precedence constraints are considered. Three examples are studied to examine the effect of the problem size. The performance indicators of an algorithm are compiled from 5 to 10 runs of each test application. Comparison between algorithms is done based on the means of these indicators. Since the objectives are the cost, the make-span, the reliability, and the impact, the diversity COV of the approximated Pareto set is defined as (Eq. 11):  $\text{COV} = (\text{cov}_{\text{cost}}, \text{cov}_{\text{reliability}}, \text{COV}_{\text{impact}}, \text{COV}_{\text{makespan}})$ .

In the first example, we consider 6 tasks and 5 generic (types) resources, combined in different sets, to generate randomly 10 initial COAs. This example is denoted by **6t-problem**. The resources are defined by their cost (fixed and in-use cost) and their reliability used to compute the cost and reliability of the COA (Guitouni et al., 2002). The impact of a COA is calculated as the median value of the actions' impact, which are ordinal values given by the decision makers (Guitouni et al., 2002). In this study, these data are randomly generated. The make-span considered here is the total delay between all the actions in the COA network. In the second example, we consider COAs defined by 50 tasks and 3 generic resources denoted by **50t-problem** and a third example with 100 tasks and 3 generic resources denoted by **100t-problem**. Initial populations of 15 and 22 COAs, for respectively the second and third examples, are created using a heuristic based on the network approach and CPLEX (Urli et al., 2003).

For the sake of comparison, we have coded the elitist non-dominated sorting genetic algorithm method (ENGA) proposed by Bagchi (1999) which has been found more efficient to discover the Pareto front compared to the well-known NSGA (Srinivas and Deb, 1995). We have chosen this method because it uses non-dominated sorting as in our approach and the diversification is based on niche formation and sharing fitness based on Pareto ranking. Moreover, in regard to the highly constrained problem treated here, ENGA was found to be easy to implement and less CPU time consuming compared to other efficient algorithms such as SPEA2 or NSGA-II.

The GA algorithm was implemented in C++. The parameters used in the procedure are:

- probability of crossover  $p_c = 0.6$ ,
- probability of mutation  $p_m = 1/\text{popsize}$  which gives a better result compared to a fixed value,
- critical population size to apply the replacement procedure  $P_c$ ,
- population size returned by the filtration procedure  $P_f$ ,
- stopping criteria  $\text{gen}_{\text{max}} = 500$  generations (6t-problem), 250 generations (50t-problem), 150 generations (100t-problem).

### 5.1 Effect of population size on the method's performance

In this section, we examine the effect, on the algorithm's performance, of the parameters  $P_c$  and  $P_f$  used in the replacement procedure (Sec. 3.4). The values of the

performance indicators  $N_s$ ,  $D_s$  and COV, summarised in Table 1, are calculated for the test application (five runs for each test) where the NDS procedure is used for the 6t-problem.

**Table 1. Population size effect on the quality of the final results**

$P_c \rightarrow P_f$	$N_s$	$D_s$	COV
30→10	1.8	0.73	(7.4; 2.6; 0; 9)
40→20	2.2	0.60	(4.6; 6.6; 5; 10)
60→30	4.7	0.70	(20; 10.5; 7; 22)
$P^* = 30$	3.5	0.73	(13; 11.5; 30; 11.3)

$P^*$ : constant size at each generation.

The results show that larger and diversified approximation sets (high  $N_s$ ,  $D_s$  and COV) are obtained when  $P_f = 30$ . This corresponds to the size of the problem with 6 tasks  $\times$  5 combinations of resources. Applying the replacement procedure when the population size reaches 60 ( $P_c = 60$ ) appears to be more adequate. Similar conclusions are derived with the two other replacement strategies (MFP and NDS+MFP).

## 5.2 Comparison between methods

First, performance of the three replacement strategies proposed for mGA (mGA-MFP, mGA-NDS, mGA-NDS+MFP) are compared. The results are for  $P_c = 60$  and  $P_f = 30$ . This choice is based on the empirical results presented in Table 1. The mean values of the performance indicators are tabulated on Table 2. Replacement strategy 3 (MFP+NDS) clearly outperforms the MFP and the NDS procedures regarding the number of non-dominated solutions  $N_s$  as well as the extent of coverage COV. The multi-criteria-filtering procedure does not generate all non-dominated solutions because some of these solutions are eliminated by the conjunctive procedure. In the NDS replacement method, among the  $P_f$  solutions filtered, those, which have the same rank, thus not dominated, are selected randomly with respect to their position in the list. In the MFP+NDS replacement strategy, these solutions are selected in accordance with their objectives by the conjunctive and disjunctive methods even if, globally, they are incomparable (not dominated). Such a strategy leads to a population of better quality, as their evaluations are greater than the threshold values.

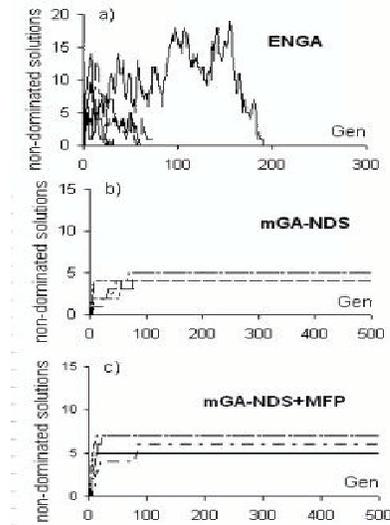
**Table 2. mGA performance using the three replacement strategies in the 6t-problem**

Method	$N_s$	$D_s$	COV
mGA- MFP	3.6	0.33	(18.7; 18.4; 0; 19)
mGA- NDS	4.7	0.70	(20; 10.5; 7; 21.7)
mGA- MFP+NDS	6.8	0.79	(26.3; 30; 25; 39)

We have also compared the features of mGA-NDS+MFP to the ENGA algorithm of Bagchi (1999) for the 6t-problem. For ENGA, the population size is fixed at

30 (such as the  $P_f$  value in our method) and the niche size,  $\sigma_{share}$ , is set to 5 which gives the best results after several trials between 1 and 5. As stated by the author, the sharing distance is computed in the phenotype space and the coefficient of the sharing function  $\alpha = 2$ . Five runs are compiled as well for each test application.

mGA- MFP+NDS is also better than ENGA as shown in Figures 4a-c. ENGA is not able to maintain a high number of feasible non-dominated solutions, from one generation to the other, for such a highly constrained problem. Only one run in five generated an interesting solutions set. The constraints in the fitness vector are not always fulfilled: they do not reach their maximum value (Eqs. 4, 4a) even if the solutions are non-dominated having the best-scored dummy fitness. In our method, feasible non-dominated solutions are archived and updated at each generation. This prevents the loss of interesting solutions during GA processing.



**Figure 4. Progress in finding non-dominated solutions on the 6t-problem. 5 runs are compiled for each method.**

## 5.3 Effect of problem size

The efficiency of these methods is also examined large size problems, which are more constrained. For the 50t-problem,  $P_c = 150$  and  $P_f = 50$  were found to represent the best compromise considering the quality of results and CPU time. For the 100t-problem,  $P_c = 150$  and  $P_f = 50$  yield the best compromise in terms of CPU time even if higher  $P_f$  values generate higher  $N_s$ . ENGA, as for the 6t-problem, failed to maintain the non-dominated solutions up to the end of the GA processing. To compare the sharing fitness approach used in ENGA to the MFP procedure, we have modified the former by archiving and updating the approximation Pareto set  $S$  discovered at each generation. Mean values for the performance indicators compiled from 10 runs are tabulated in Table 3.

The extent of the coverage on the third objective (the impact) is nil ( $cov_{impact} = 0$ ) for these large-size problems.

This result is the consequence that all non-dominated solutions always score the median value for this objective.

The mGA-NDS+MFP generates a larger number of efficient solutions for both problems and in the 50t-problem a better diversity in the decision and the objective spaces. For the 100t-problem, even if the COV is similar for all the methods, the simulations depicted in Figure 5 illustrate how mGA-NDS+MFP outperforms the modified ENGA in its ability to find better extremes. Data correspond to the 10 processed runs results. Its superiority comes from the multi-criteria filter based on the disjunctive and dynamic conjunctive methods.

**Table 3. Algorithm comparison on different problem size**

	Ns	Ds	COV
<b>Algorithm</b>	<b>50t-problem</b>		
mGA- NDS	13	0.75	(44; 77; 0; 50)
mGA- MFP+NDS	16	0.76	(68; 94; 0; 34)
modified ENGA	10	0.71	(43; 51; 0; 30)
<b>Algorithm</b>	<b>100t-problem</b>		
mGA- NDS	16.7	0.62	(20.7; 13; 0; 18)
mGA- MFP+NDS	22.6	0.61	(17; 11.2; 0; 13)
modified ENGA	12.1	0.64	(18; 9; 0; 12)

## 6 Conclusion

A new approach is proposed to optimise multi-objective, large-size, highly constrained problem such as COAs planning. This approach, based on genetic algorithms, uses non-dominance and a first order multi-criteria filtering method in the replacement procedure. This approach was compared to the classical one, based on niche formation and sharing function. This is done by analysing the approximation Pareto sets generated by the new algorithm, mGA-NDS+MFP, and by the elitist non-sorting dominated algorithm, ENGA (Bagchi, 1999). Regarding the cardinality and the diversity of these approximation sets, it was possible to conclude that mGA-NDS+MFP is more efficient than ENGA. Moreover, mGA-NDS+MFP does not present the drawbacks inherent to the classical approach such as fine-tuning hard parameters (e.g. niche size and distance sharing). By embedding the dynamic conjunctive and disjunctive methods in the replacement procedure, mGA-NDS+MFP was shown to be able to find solutions that are more consistently closer to the Pareto front.

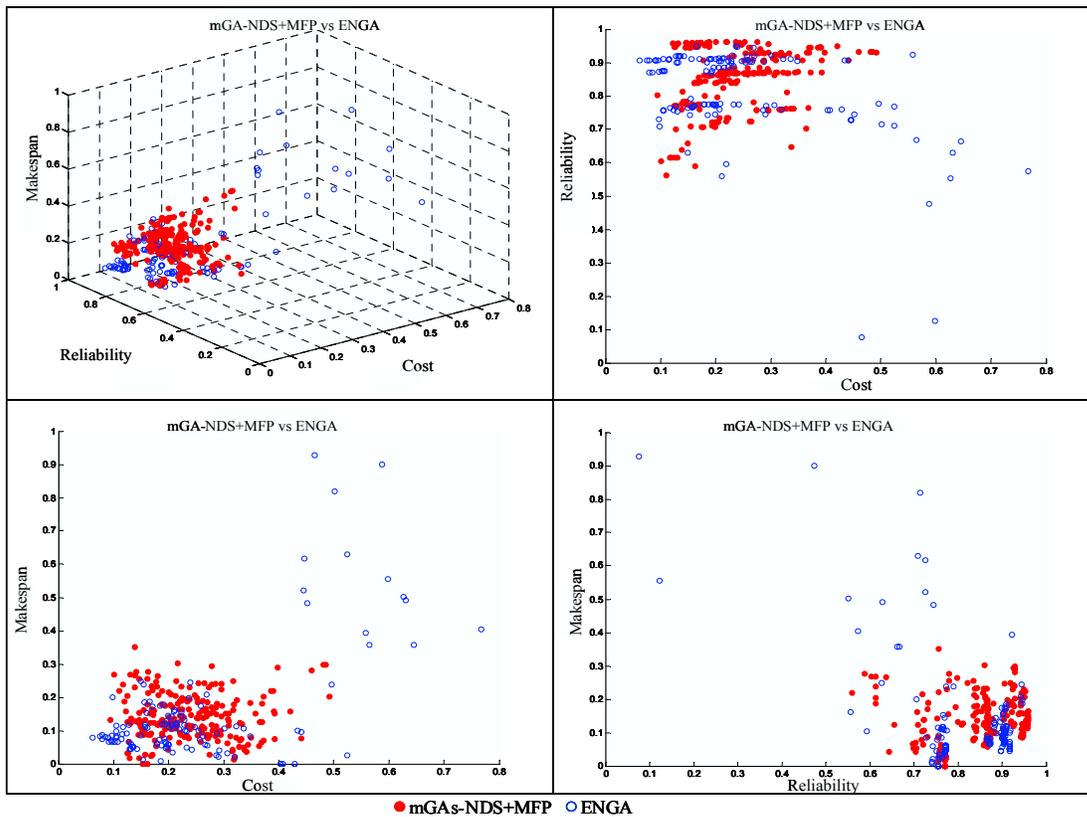
It is also important to underline the limitations of performance metrics (indicators) used in this paper. In fact, we think that future works should propose new metrics to characterise the Pareto frontier. This characterisation could include for example, the dispersion, the shape, and the lower and upper bounds of the frontier.

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**Figure 5: Empirical comparison of mGA-NDS+MFP (filled points) and modified ENGA (empty points) for the 100t-problem**

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