

# Approximating multi-objective time-dependent optimization problems

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## **Approximating Multi-Objective Time-Dependent Optimization Problems**

Said Dabia, El-Ghazali Talbi, Tom van Woensel, Ton de Kok

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# Approximating Multi-Objective Time-Dependent Optimization Problems

Said Dabia

Eindhoven University of Technology, School of Industrial Engineering, The Netherlands, s.dabia@tue.nl

El-Ghazali Talbi

University of Lille, INRIA, CNRS, France, El-ghazali.Talbi@lil.fr

Tom Van Woensel, Ton De Kok

Eindhoven University of Technology, School of Industrial Engineering, The Netherlands,  
{t.v.woensel@tue.nl, a.g.d.kok@tue.nl}

In many practical situations, decisions are multi-objective in nature. Furthermore, costs and profits are time-dependent, *i.e.* depending upon the time a decision is taken, different costs and profits are incurred. In this paper, we propose a generic approach to deal with multi-objective time-dependent optimization problems (MOTDP). The aim is to determine the set of Pareto solutions that capture the interactions between the different objectives. Due, to the complexity of MOTDP, an efficient approximation based on dynamic programming is developed. The approximation has a provable worst case performance guarantee. Even though the approximate Pareto set consists of less solutions, it represents a good coverage of the true set of Pareto solutions. Numerical results are presented showing the value of the approximation.

*Key words:* Multi-objective optimization; Time-dependent costs; Approximation; dynamic programming;  $\epsilon$ -dominance

*History:*

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## 1. Introduction

Many optimization problems encountered in practice are multi-objective in nature, *i.e.* different objectives are conflicting and equally important. Many times, it is not desirable to drop some of them or to optimize them in a composite single objective or hierarchical manner. For instance, while designing a product, many criteria are taken into account: *e.g.* the product's reliability should be maximized, while the cost and the environmental impact should be minimized. Obviously, the amount of examples that can be formulated is infinite.

Contrary to single-objective optimization problems where the optimal solution is a single candidate, the aim of Multi-Objective optimization Problems (in short, MOPs) is the determination of the set of points representing the compromise solutions between the different conflicting objectives. This set of points is defined as the set of Pareto solutions or the Pareto front. A solution is Pareto, if it is not possible to improve an objective without deteriorating at least another one. In this line of thought, decision makers are presented with the entire Pareto front (rather than a single solution) such that they can select their preferred solution depending on their specific situation. Although the roots of multi-objective optimization go back to the nineteenth century in the work of Edgeworth (1881) and Pareto (1896), in the literature, most optimization problems dealt with are mono-objective. In fact, multi-objective cost functions are usually reduced to a composite single objective cost function by using a (weighted) sum of the various objectives. It is argued that solutions obtained as such are only a small subset of the entire set of Pareto solutions, and therefore could lead to suboptimal or infeasible managerial decisions (Ehrgott (2005), Miettinen (1999) and Talbi (2009)).

When dealing with Multi-Objective optimization Problems, a number of complicating factors influence both the quality and easiness of obtaining the Pareto fronts. First, the number of Pareto solutions increases with the size of the problem, mainly with the number of objectives. Therefore, multi-objective decision making becomes very challenging. In fact, most multi-objective problems are NP-hard. Hence, it is computationally very expensive to compute the complete Pareto front. Furthermore, it is not straightforward to select a solution from the Pareto fronts. Consequently, many researchers direct their efforts on approximating the Pareto front in the hope of reducing the complexity of the applied algorithms and producing good approximations (*i.e.* approximate Pareto fronts) of the Pareto front. Approximate Pareto fronts contain less solutions, which facilitate the selection of a final solution. However, a good approximate Pareto front should have enough solutions to properly represent the real Pareto front. Second, in many practical settings, the cost parameters change over time which makes scheduling problems harder. For instance, in a traffic network, travel costs are a function of travel times that change depending on the time of the day a vehicle is traveling.

In this paper, an approximation algorithm based on dynamic programming is proposed for *Multi-objective Time-Dependent Optimization Problems* (MOTDPs). The multi-dimensional profit space is partitioned into intervals with exponentially increasing size. Each interval defines a cluster of solutions that are considered to be *very close* to each other. From each cluster only one solution is kept and the dynamic programming is adapted to the partitioned profit space. In this way, in each iteration of the dynamic programming only a polynomial number of solutions is processed. The approximation has a provable performance guarantee. Moreover, it is easy for a decision maker to select a solution as he is provided with a limited number of solutions. Even with less solutions, the resulting approximate Pareto front still properly covers the real Pareto front in the sense that each optimal Pareto solution is represented by at least one approximate Pareto solution. The evaluation of the proposed approach is demonstrated on a time-dependent capacitated traveling salesman problem with time windows, though the approximation presented is applicable for a generic MOTDP. In fact, the proposed approximation could be applied to a variety of well-known optimization problems for which a dynamic programming formulation is possible (*e.g.* knapsack problems, shortest path problems, traveling salesman problems, variants of vehicle routing problems, sequence alignment problems, ...).

The contributions of this paper are summarized as follows. A generic approximation is proposed which can be applied to structured optimization problem that can be solved by means of dynamic programming. The approximation generates an approximate Pareto front with less solutions (a polynomial number of solutions). However, the approximate Pareto front represents a very good coverage of the real Pareto front. Additionally, the approximation's worst case performance guarantee is provable. Hence, it is easier for the decision to select a solution for which he has a good feeling about its quality. Furthermore, the approximation is very flexible in the sense that the decision maker can choose different precision levels for the different objectives. In fact, the decision maker might be willing to tolerate more error for objectives that are less sensible. Finally, we are dealing with a realistic MOP for which costs are dynamic. In fact, in most real-life situations costs are time-dependent. To our knowledge, this is the first time a multi-objective time-dependent optimization problem is approximated.

The paper is organized as follows. Section 2 reviews the literature relevant to MOPs. Section 3 is devoted to the introduction of the main concepts related to MOTDPs. Section 4 describes a generic MOTDP and the assumptions made to guarantee the non-dominance principle. In Section 5, an approximation of the Pareto front is developed and the main results of the paper are derived. In section 6, the main results are validated based on a specific traveling salesman problem with time windows. Finally, Section 7 concludes with a summary of the main results.

## 2. Literature Review

As in mono-objective optimization, MOPs can be divided into two categories: those whose solutions are encoded with *real-valued* variables, also known as *continuous MOPs*, and those where the solutions are encoded using *discrete* variables, known as multi-objective combinatorial optimization problems (*MOCO*). In the class of continuous MOP, an infinite number of Pareto solutions composes the Pareto front whereas in combinatorial MOPs, both the feasible set  $S$  and the Pareto front are finite. Most heuristics for solving MOPs are designed to deal with continuous MOPs using, for instance, multi-objective simplex (Zeleny (1982) and Steuer (1986)). In the last decade, there is also a growing interest in solving combinatorial MOPs. However, in most of the cases, they are bi-objective optimization problems. Furthermore, there is a lack of test instances for real-life combinatorial MOPs, especially problems with many objectives (Ishibuchi et al. (2008)), uncertainty (Liefoghe et al. (2007)) and dynamicity (Farina et al. (2004)).

The study of computational complexity classes for MOCO started with the work of Serafini (1986), and Papadimitriou and Yannakakis (2002). They made a connection between the results obtained in mono-objective combinatorial optimization and the multi-objective field for several canonical problems. Serafini (1986) depicted nine possible definitions for MOCO problems and established reductions between them in order to facilitate obtaining complexity results. He showed that the following definition (denoted as V8 in his article) can be considered as a standard reference version to measure the computational complexity of MOCO problems. The definition can also be seen as the decision problem associated with a MOCO problem.

**DEFINITION 1 (GENERIC DEFINITION OF MOCO BY SERAFINI (1986)).** . Given  $z \in Z^n$ , does there exist  $x \in X$  such that  $f_i(x) \leq z$ ?

A NP-hard mono-objective problem implies a NP-hard character to its multi-objective extensions. In the multi-objective case the NP-hard class appears for the majority of problems. For some of them derived from mono-objective NP-hard problems, it is easy to prove also their NP-hardness. This is a reason why the study must focus on determining a specific stronger class as NP-hard in the strong sense or results about the NP-completeness. For example, NP-completeness is proved for shortest path problems, assignment problems and minimum maximal matching by Serafini (1986); for the minimum weight spanning tree by Camerini and Vercellis (1984); and for the max-linear spanning tree by Hamacher and Ruhe (1994).

Similarly to mono-objective optimization problems, MOPs can be solved by means of *exact* and *approximate* algorithms. In the literature, more attention has been devoted to bi-criteria optimization problems by using exact methods such as branch and bound algorithms (Sen et al. (1988), Ulungu and Teghem (1995), Visée et al. (1998), Sayin and Karabati (1999) and Lemesre et al. (2007a)), branch and cut (Jozefowicz et al. (2007)), *A\** algorithm (Stewart and White (1991), Mandow and Millan (1996)), and dynamic programming (White (1982) and Carraway et al. (1990)). Because of the complexity of MOPs, exact methods are only effective for problems with small instances and with no more than two criteria. However, there exists some new advances in this area, with several exact approaches proposed in the literature for bi-objective (Lemesre et al. (2006), Laumanns et al. (2004) and Lemesre et al. (2007b)) and multi-objective problems (Lemesre et al. (2006)). Approximate methods are mainly used to solve large-scale problems and when multiple criteria are involved. They can be divided into two classes: on the one hand algorithms that are only applicable to a specific problem. Such algorithms are developed based on some knowledge on the structure of the problem at hand. On the other hand, meta-heuristics which are of general purpose, in the sense that they can be applicable to a large variety of MOPs. A unifying view for analyzing, designing and implementing multi-objective meta-heuristics is provided in the book Talbi (2009). The main drawback of meta-heuristics is that they do not guarantee the performances related to the Pareto front. Moreover, the resulting approximate Pareto fronts might not properly cover the real Pareto front as they might contain very few solutions.

In the context of mono-objective optimization problems, an  $\epsilon$ -approximation scheme is an algorithm that, for every instance of the problem, finds an approximate solution that is guaranteed to be within a constant factor from optimal. Two classes of approximation schemes are mainly considered: Polynomial Time Approximation Scheme (PTAS) and Fully Polynomial Time Approximation Scheme (FPTAS). For any  $\epsilon > 0$ , a PTAS runs in time polynomial in the size of the instance, while an FPTAS runs in time polynomial in the size of the instance and  $1/\epsilon$ . From a computational complexity point of view, FPTAS are the strongest approximation schemes with performance guarantee that can be obtained for optimization problems. The notion of approximation schemes can be generalized to the case of multi-objective optimization problems by considering, for each solution on the approximate Pareto front, worst case performance guarantees with regard to all criteria.

### 3. Definitions, Variables and Background

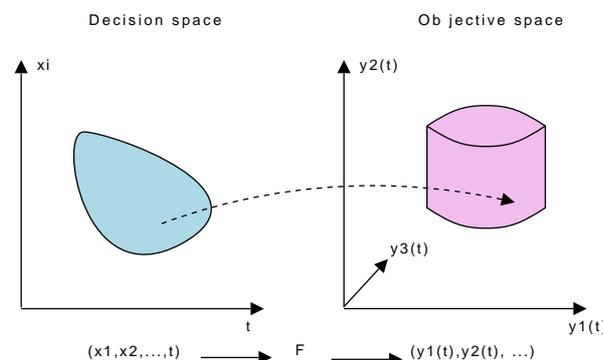
This section aims to give the most relevant definitions and notations used in the remainder of the paper.

**DEFINITION 2 (MULTI-OBJECTIVE TIME-DEPENDENT OPTIMIZATION PROBLEM).** A multi-objective time-dependent optimization problem (MOTDP) is defined as:

$$(MOTDP) = \begin{cases} \text{vmin } F(x, t) = (f_1(x, t), f_2(x, t), \dots, f_n(x, t)) \\ \text{s.t. } (x, t) \in S \end{cases} \quad (1)$$

where  $n$  ( $n \geq 2$ ) is the number of objectives,  $(x, t) = (x_1, \dots, x_k, t)$  is the vector representing the decision variables depending upon the starting time  $t$ , and  $S$  represents the set of feasible solutions associated with equality and inequality constraints, and explicit bounds.  $F(x, t) = (f_1(x, t), f_2(x, t), \dots, f_n(x, t))$  is the vector of objectives to be optimized. Note that the value of MOTDP also depends on the starting time  $t$ , which could also be considered as a decision variable.

The search space  $S$  represents the decision space or parameter space of the MOTDP. The space to which the objective vector belongs is called the objective space.  $F$  is a mapping from the decision space to the objective space which evaluates, for a starting time  $t$ , the quality of each solution  $(x_1, \dots, x_k)$  by assigning an objective vector  $(y_1(t), \dots, y_n(t))$ . The objective vector represents the quality (or *fitness*) of the solution (Fig. 1). The decision maker is usually interested in the value of a solution on each criterion. Therefore, the analysis of MOPs is done in the objective space. The set  $Y = F(S)$  represents the feasible points in the objective space, and  $y(t) = F(x, t) = (y_1(t), y_2(t), \dots, y_n(t))$ , where  $y_i(t) = f_i(x, t)$ , is a point in the objective space.



**Figure 1** Decision space and objective space in a time-dependent MOP.

DEFINITION 3 (PARETO DOMINANCE). An objective vector  $u(t) = (u_1(t), \dots, u_n(t))$  is said to dominate  $v(t') = (v_1(t'), \dots, v_n(t'))$  (we write  $u(t) \prec v(t')$ ) if and only if no component of  $v(t')$  is smaller than the corresponding component of  $u(t)$  and at least one component of  $u(t)$  is strictly smaller, *i.e.*

$$\forall i \in \{1, \dots, n\} : u_i(t) \leq v_i(t') \wedge \exists i \in \{1, \dots, n\} : u_i(t) < v_i(t').$$

DEFINITION 4 (PARETO OPTIMALITY). A solution  $(x^*, t^*) \in S$  is Pareto Optimal<sup>1</sup> if for every  $(x, t) \in S$ ,  $F(x, t)$  does not dominate  $F(x^*, t^*)$ , *i.e.*  $F(x, t) \not\prec F(x^*, t^*)$ .

A MOTDP involves the determination of a set of solutions known as the *Pareto optimal set*. The image of this set in the objective space is denoted as the *Pareto front*. We define the Pareto optimal set and Pareto front as follows:

DEFINITION 5 (PARETO OPTIMAL SET). For a given MOTDP  $(F, S)$ , the Pareto optimal set is defined as  $\mathcal{P}^* = \{(x, t) \in S / \nexists (x', t') \in S, F(x', t') \prec F(x, t)\}$ .

DEFINITION 6 (PARETO FRONT). For a given MOTDP  $(F, S)$  and its Pareto optimal set  $\mathcal{P}^*$ , the Pareto front is defined as  $\mathcal{PF}^* = \{F(x, t), (x, t) \in \mathcal{P}^*\}$ .

The generation of the Pareto optimal set often turns out to be practically impossible or computationally too expensive. Therefore, good approximations of  $\mathcal{PF}^*$  are desirable. We define the  $\epsilon$ -Pareto concept as follows:

DEFINITION 7 ( $\epsilon$ -DOMINANCE). An objective vector  $u(t) = (u_1(t), \dots, u_n(t))$  is said to  $\epsilon$ -dominate  $v(t') = (v_1(t'), \dots, v_n(t'))$  (we write  $u(t) \prec_\epsilon v(t')$ ) if and only if no component of  $v(t')$  is dominated by the corresponding component of  $u(t) - \epsilon$ , *i.e.*

$$\forall i \in \{1, \dots, n\} : u_i(t) - \epsilon \leq v_i(t') \wedge \exists i \in \{1, \dots, n\} : u_i(t) - \epsilon < v_i(t').$$

DEFINITION 8 ( $\epsilon$ -PARETO OPTIMALITY). A solution  $(x^*, t^*) \in S$  is  $\epsilon$ -Pareto Optimal if for every  $(x, t) \in S$ ,  $F(x, t)$  does not  $\epsilon$ -dominate  $F(x^*, t^*)$ , *i.e.*  $F(x, t) \not\prec_\epsilon F(x^*, t^*)$  (Fig.2).

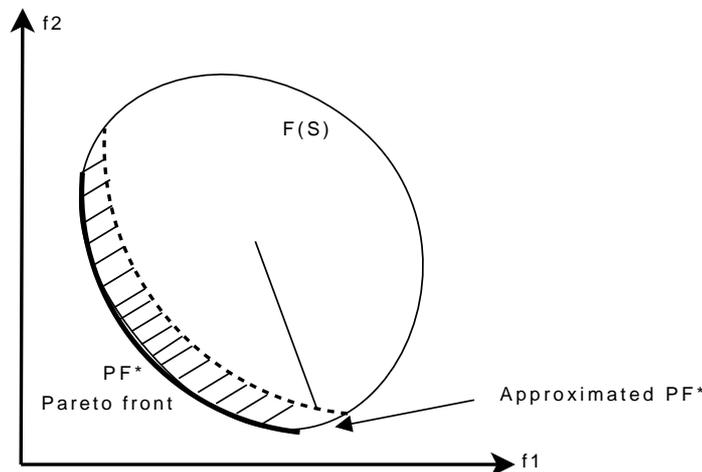


Figure 2  $\epsilon$ -Pareto concept. Sets of Pareto and  $\epsilon$ -Pareto solutions.

<sup>1</sup> The Pareto optimal solutions are also known under the name of *acceptable solutions*, *efficient*, *not-dominated*, *non-inferior*.

#### 4. A Dynamic Programming formulation for the MOTDP

We consider a generic Multi-Objective Time-Dependent Optimization problem (MOTDP) related to a set of items  $H$ . Practically,  $H$  represents jobs to be processed on a machine, or customers to be served by a truck or project tasks to be executed. The MOTDP involves the joint minimization of  $n$  objectives. We assume that each instance of MOTDP naturally decomposes into  $|H|$  arrays of vectors  $\{X^1, X^2, \dots, X^{|H|}\}$ . For every  $k \in \llbracket 1, |H| \rrbracket$ , the array  $X^k$  consists of the vectors  $\{I_1^k, I_2^k, \dots, I_{|X^k|}^k\}$  such that for all  $k \in \llbracket 1, |H| \rrbracket$  and  $p \in \llbracket 1, |X^k| \rrbracket$ ,  $I_p^k$  is an instance of MOTDP. By considering  $(x^\alpha)_{\alpha \in \mathbb{N}}$  to be a set of mappings from  $H$  to  $H$ ,  $I_p^k$  is represented by the vector  $[x^\alpha(1), x^\alpha(2), \dots, x^\alpha(k)]$  for some  $\alpha \in \mathbb{N}$ . Note that  $|X^k|$  depends on MOTDP and not on any instance of it.  $I_p^k$  is then a schedule consisting of at most  $k$  items, and  $X^k$  the set of schedules with at most  $k$  items. To each item  $i \in H$  a time-dependent vector costs  $c^i(t) = [c_1^i(t), c_2^i(t), \dots, c_n^i(t)]$  is associated. The cost  $c_r^i(t)$  is related to the  $r^{\text{th}}$  objective and is incurred when item  $i$  is processed at time  $t$ . We consider the first objective to be the total processing time of the items in  $H$ . Consequently,  $c_1^i(t)$  is the processing time of item  $i$ . Without loss of generality we assume that MOTDP is bounded in all objectives and let  $B$  denote the vector bounds such that  $B_r$ ,  $r \in \llbracket 1, n \rrbracket$ , is the upper bound corresponding to the  $r^{\text{th}}$  objective. Table1 summarizes the notations used in the paper.

**Table 1** Notation used in this paper.

Variable	Description
$H$	: A set of items
$t$	: Time. Time origin is always taken to be 0
$t_0$	: Starting time, $t_0 \geq 0$
$S$	: The set of feasible solutions
$t^i$	: starting time of processing item $i$
$t_X$	: Processing time of the subset of items $X \subseteq S$
$ X $	: Size of the set $X$
$[y_r]_{r=1}^{r=n}$	: The vector (or array of vectors) $y_1, y_2, \dots, y_n$
$\lfloor x \rfloor$	: Nearest integer smaller or equal to the real number $x$
$\lceil x \rceil$	: Nearest integer larger or equal to the real number $x$
$\llbracket a, b \rrbracket$	: The interval of integer numbers between $a$ and $b$ ( $a$ and $b$ are also integers).
$\mathbb{N}$	: The set of natural numbers
$\mathbb{R}$	: The set of real numbers
$\epsilon$	: The approximation worst case precision

The MOTDP is solved by a Dynamic Program ( $DP$ ).  $DP$  goes through  $|H|$  iterations such that, in the  $k^{\text{th}}$  iteration it processes the input  $X^k$  and generates  $X^{k+1}$ , the input to the next iteration. For simplicity and without loss of generality, we assume that the starting time  $t_0$  is given. Therefore, when unnecessary the time index is omitted from the notation. The value of  $DP$  is then denoted  $F(X^k)$  instead of  $F(X^k, t_0)$ , and is represented by the array of vectors  $\{f^k(I_1^k), f^k(I_2^k), \dots, f^k(I_{|X^k|}^k)\}$  where  $f^k$  is a mapping from  $X^k$  to  $\mathbb{R}^n$  such that:

$$f^k(I_p^k) = [f_r^k(I_p^k)]_{r=1}^{r=n} \quad \text{and} \quad f_r^k(I_p^k) = \sum_{i=1}^k c_r^{x^\alpha(i)}(t^{x^\alpha(i)}) \quad (2)$$

$t^{x^\alpha(i)}$  is calculated in an iterative way as follows:

$$t^{x^\alpha(1)} = t_0 \quad \text{and} \quad t^{x^\alpha(i+1)} = t^{x^\alpha(i)} + c_1^{x^\alpha(i)}(t^{x^\alpha(i)}) \quad (3)$$

The structure of  $DP$  is assuming the generation of Pareto solutions in all iterations. However, because of time-dependency, the generation of the Pareto front is not guaranteed. In fact, time-dependency makes it possible to generate non Pareto solutions from Pareto ones. This problem

can be solved by imposing some structure on the vector costs. This structure is illustrated in the following assumption:

ASSUMPTION 1 (FIFO). For every item  $i$ , and times  $t_1$  and  $t_2$ ,  $t_1 \leq t_2$  implies that:

$$t_1 + c_1^i(t_1) \leq t_2 + c_1^i(t_2) \quad (4)$$

$$c_r^i(t_1) \leq c_r^i(t_2), \text{ for all } r \in \llbracket 2, n \rrbracket \quad (5)$$

Remember that the first objective reflects the total processing time of the items in  $H$ . Assumption 1 thus means that for every item  $i \in H$  the processing time adheres to the well-known FIFO assumption (*i.e.* tasks cannot overtake each other). Furthermore, all other objectives costs are monotonically increasing in time.

The recursion for  $DP$  is formulated as follows:

$$F(X^k) = \underset{\substack{1 \leq p \leq |X^{k-1}| \\ i \in H \setminus I_p^{k-1}}}{vmin} \left\{ \left[ f_r(I_p^{k-1}) + c_r^i(t_0 + t_{I_p^{k-1}}) \right]_{r=1}^{r=n}, (I_p^{k-1}, t_0) \in S \right\} \quad (6)$$

In most cases  $|X^k|$  is exponential in the size of MOTDP. Consequently, the running time of  $DP$  is exponential too. Therefore, the determination of the optimal Pareto front for most MOTDP is very time consuming. In the next section, an approximate dynamic program ( $DP^\epsilon$ ) is presented.

## 5. Approximation based on Dynamic Programming, $DP^\epsilon$

In order to reduce the complexity of  $DP$ , we impose extra structure during its execution. In each iteration  $k$ , we trim the generated set  $F(X^k)$  of Pareto solutions by eliminating the solutions that are very close to each other. The trimmed set  $F(\tilde{X}^k)$  is then used in the dynamic program to compute the untrimmed set  $F(X^{k+1})$ . The idea of adding this type of structure to the execution of algorithms was first introduced by Ibarra and Kim (1975). Sahni (1976) and Woeginger (2005) applied it to a variety of single-objective and time-independent scheduling problems.

### 5.1. Setting up $DP^\epsilon$

The new approximate dynamic program recursion is formulated as follows:

$$F(X^k) = \underset{\substack{1 \leq p \leq |\tilde{X}^{k-1}| \\ i \in H \setminus \tilde{I}_p^{k-1}}}{vmin} \left\{ \left[ f_r(\tilde{I}_p^{k-1}) + c_r^i(t_0 + t_{\tilde{I}_p^{k-1}}) \right]_{r=1}^{r=n}, (\tilde{I}_p^{k-1}, t_0) \in S \right\} \quad (7)$$

Formally, the set  $F(X^k)$  can be represented by geometric points in the polyhedron  $[0, B_1] \times [0, B_2] \times \dots \times [0, B_n]$ . The polyhedron is cut into multiple boxes of exponentially increasing size. Points contained by the same box are considered to be very close to each other. In each box, only one point is retained. Figure 3 illustrates the trimming of the set of Pareto solutions in case  $n = 2$ . Obviously, the size of boxes depends on the precision vector  $\epsilon = [\epsilon_r]_{r=1}^{r=n}$ . Smaller precisions result in smaller boxes. Moreover, boxes with an exponentially increasing size result in a sort of logarithmic scale with a polynomial number of boxes (polynomial in the input size). Therefore, after trimming, only a polynomial number of solutions is kept. Note that if the boxes' size increases linearly, the number of solutions kept will still be exponential.

The cuts on the axis corresponding to the  $r^{th}$  objective are executed at the coordinates  $\Delta_r^{m_r}, m_r = 1, \dots, L_r$ , such that:

$$\Delta_r = 1 + \frac{\epsilon_r}{2|H|} \quad (8)$$

Where  $\epsilon = [\epsilon_r]_{r=1}^{r=n}$  is a vector of real numbers between 0 and 1 representing the approximation's precision vector.  $\epsilon_r$  is the precision related to the  $r^{th}$  objective.

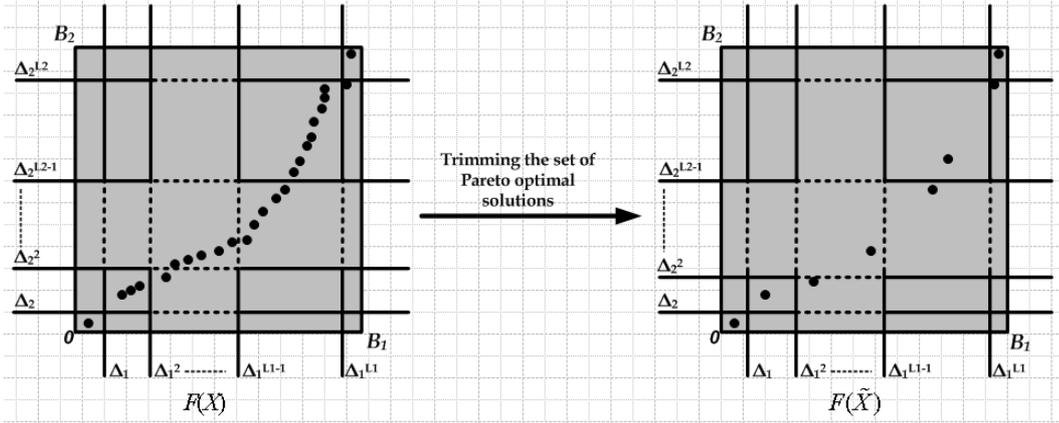


Figure 3 The reduction of the set of Pareto-optimal solutions in case of two objectives.

Note that different precisions can be assigned to the different objectives. Objectives that are less sensitive to errors (*e.g.* with a flat cost structure) could be assigned larger precision values. Furthermore, the form of the vector  $\Delta = [\Delta_r]_{r=1}^{r=n}$  given by equation (8) is justified by two reasons. First, the values of the vector  $\Delta = [\Delta_r]_{r=1}^{r=n}$  will have values very close to 1. Hence, two solutions in the same box are indeed very close to each other. Second, we know how the sequence  $(1 + \frac{x}{a})^a$  behaves when  $a$  goes to infinity.

The values of  $L_r$  are chosen such that  $\Delta_r^{L_r} \leq B_r$ . We choose:

$$L_r = \left\lceil \frac{\ln B_r}{\ln \Delta_r} \right\rceil \leq \left\lceil \left(1 + \frac{2|H|}{\epsilon_r}\right) \ln B_r \right\rceil \quad (9)$$

The trimmed set of Pareto solutions contains at most  $\prod_{r=1}^n L_r$  solutions. We can compute the complexity of  $DP^\epsilon$  as being proportional to:

$$\begin{aligned} \sum_{k=1}^{|H|} |F(\tilde{X}^k)| &= O\left(|H| \prod_{r=1}^n L_r\right) \\ &= O\left(|H|^{n+1} \prod_{r=1}^n \left(\frac{\ln B_r}{\epsilon_r}\right)\right) \end{aligned} \quad (10)$$

Given all elements of the arrays  $X^k$  are binary coded, Equation (9) and (10) shows that the running time of  $DP^\epsilon$  is polynomial in the input size and in the vector  $\frac{1}{\epsilon} = \left[\frac{1}{\epsilon_r}\right]_{r=1}^{r=n}$ . Moreover, the running time is still exponential in the number of objectives. However, in most cases the number of objectives is limited. Obviously there is a trade-off between the values of the elements of the vector  $\epsilon$  and the running time of  $DP^\epsilon$ . In fact, for small values of  $\epsilon$  more solutions are kept during the execution of  $DP^\epsilon$  as the boxes illustrated in Figure 3 are smaller. Therefore, more data is processed which results in an increase of the running time. Intuitively, the quality of the approximation depends on  $\epsilon$ . In fact, because of the trimming action,  $DP^\epsilon$  generates incorrect data. However, less data is trimmed for small values of  $\epsilon$  which limits the error caused by the trimming action.

## 5.2. Worst Case Performance of $DP^\epsilon$

We show that the worst case performance guarantee is such that every solution generated by  $DP$  is at most a constant factor  $\epsilon$  from that of a  $DP^\epsilon$  solution. For the sake of a clear presentation, the proofs needed in this section are available in the Appendix. Henceforth, we assume the following structure regarding the vector costs:

ASSUMPTION 2. For every item  $i \in S$  and real number  $1 \leq \alpha \leq 2$ , it holds that for every time  $t$  and  $r \in \llbracket 1, n \rrbracket$ :

$$c_r^i(\alpha t) \leq \alpha c_r^i(t) \quad (11)$$

Assumption 2 means that processing item  $i$  at a later time  $\alpha t$ , instead of time  $t$ , will not multiply costs by more than a coefficient  $\alpha$ . In other words, sudden increases in costs are not allowed (which seems reasonable from a real-life point of view). For instance, the costs functions  $c_r^i(t) = t$ ,  $c_r^i(t) = \sqrt{t}$  and  $c_r^i(t) = \ln(t)$  satisfy assumption (2). Note that the time origin is taken to be 0. We show that if assumption 2 is satisfied, the following lemma holds:

LEMMA 1. For all  $p \in \llbracket 1, |X^k| \rrbracket$ , there exists  $q \in \llbracket 1, |\tilde{X}^k| \rrbracket$  such that for all  $k \in \llbracket 1, |H| \rrbracket$  and  $r \in \llbracket 2, n \rrbracket$  (given assumptions 1 and 2):

$$t_{\tilde{I}_q^k} \leq \Delta_1^k t_{I_p^k} + (\Delta_1^k - 1)t_0 \quad \text{and} \quad f_r(\tilde{I}_q^k) \leq \max(\Delta_1^k, \Delta_r^k) f_r(I_p^k) \quad (12)$$

Lemma 1 presents an important result as upper bounds on the quality of the approximation are proved. Furthermore, it shows that the approximate Pareto front covers well the real Pareto front. In fact, every Pareto solution is closely approximated by at least one solution from the approximate Pareto front. In the following theorem, we further demonstrate that the upper bounds depends on the vector  $\epsilon$  and on the starting time  $t_0$ :

THEOREM 1. For all  $p \in \llbracket 1, |X^k| \rrbracket$ , there exists  $q \in \llbracket 1, |\tilde{X}^k| \rrbracket$  such that for all  $k \in \llbracket 1, |H| \rrbracket$  and  $r \in \llbracket 2, n \rrbracket$  (Given assumptions 1 and 2):

$$t_{\tilde{I}_q^k} \leq (1 + \epsilon_1) t_{I_p^k} + \epsilon_1 t_0 \quad \text{and} \quad f_r(\tilde{I}_q^k) \leq (1 + \max(\epsilon_1, \epsilon_r)) f_r(I_p^k) \quad (13)$$

Theorem 1 also shows that there is a trade-off between the quality of approximation and the starting time. In fact, a later starting time ( $t_0 > 0$ ) might result in reduced processing time (e.g. congestion might be avoided). However, a starting time  $t_0 \neq 0$  triggers an additional small error  $\epsilon_1 t_0$  in the first objective. Moreover, upper bounds of the other objectives are affected by  $\epsilon_1$ . This is not a surprise since costs are time-dependent and the starting time of processing item  $i$  is affected by  $\epsilon_1$ . Furthermore, in the special case where  $t_0 = 0$ ,  $\max(\epsilon_1, \epsilon_r) = \epsilon_r$  for all  $r \in \llbracket 2, n \rrbracket$ , and the number of objectives  $n$  is fixed,  $DP^\epsilon$  belongs to the family of FPTAS algorithms.

## 6. Computational results

To validate the material presented in the previous sections, a time-dependent capacitated traveling salesman problem with time windows and multiple tours is considered. The described problem is relevant in the situation where a vehicle is required to fill up ATMs located at different places from a central bank. For security reasons, it is not allowed to carry a large amount of money. Consequently, the vehicle is forced to make several short tours during its operating period (e.g. a working day) going back and forth to the central bank. Similarly, in the case of food home delivery, tours are relatively short as products are perishable (e.g. should remain warm) and thus need to be delivered as soon as possible to their destinations (Azi et al. 2007).

A single vehicle performs several tours to serve a set  $H = \{1, 2, \dots, N\}$  of geographically dispersed customers. The vehicle has a finite capacity  $Q$  and is only available for a limited amount of time  $T$ . Moreover, the tours' duration is restricted (e.g. due to quality or security issues), to last not more than  $t^{lim}$  time units. The same single vehicle could of course do multiple tours, if time allows for this. Because of road congestion, travel times are time-dependent: depending on the departure time  $t^i$  at customer  $i$ , a different travel time  $tt_{ij}(t^i)$  to customer  $j$  is incurred. Furthermore, all customers need to get delivered in their specific hard time windows. Service at customer  $i$ , with duration  $s_i$ , can only be started after its opening time  $t_i^l$  and no delivery is allowed after its closing

time  $t_j^u$ . We consider a situation with tight time windows, *i.e.* the size of time windows is small compared to travel time.

We consider a multi-objective cost function: simultaneously minimizing the total time traveled, including service time and waiting times at customers due to time windows, and maximizing the total demand fulfilled. Therefore, the vector costs is such that:

$$[c_{ij}(t^i)] = \begin{bmatrix} tt_{ij}^*(t^i) = tt_{ij}(t^i) + \max(0, t_j^l - t_i - tt_{ij}(t^i)) + s_j \\ d_j \end{bmatrix} \quad (14)$$

in which  $d_j$  is the stationary demand of customer  $j$ .

For our numerical study, the customers' location as well as the related demand are taken from the Solomon's instances (Solomon (1987)). Congestion is taken into account, by assuming that the speed on each link is time-dependent and derive the travel time profile by using the relation  $tt_{ij}(t) = \frac{d_{ij}}{v_{ij}(t)}$  where  $d_{ij}$ , distance between customers  $i$  and  $j$ , is computed based on Solomon's data sets.  $v_{ij}(t)$  is the time-dependent speed by which the vehicle traverse the link between customers  $i$  and  $j$ . The resulting travel times satisfy both assumptions 1 and 2. The congested speed is  $v_c = 30km/hr$ , and the free speed is  $v_f = 70km/hr$ . Furthermore, we take  $t^{lim} = \{500, 2000\}$ ,  $Q = \{100, 200\}$  and  $\epsilon = \{0.01, 0.05, 0.1, 0.3\}$ . We consider instances with 100 and 300 customers and a planning horizon  $T = 6000$  minutes.

To compare the different Pareto fronts generated, we introduce two measures defined by Zitzler et al. (2000): the two-set coverage metric and the average distance metric. The choice of the metrics is motivated by their intuitive explanation. The *two-set coverage* metric is defined as:

$$C(X \succ Y) = \frac{|\{y \in Y; \exists x \in X : x \text{ dominates } y\}|}{|Y|} \quad (15)$$

in which  $X$  and  $Y$  are two Pareto curves. The two-set coverage metric calculates the fraction of solutions in  $Y$  that are dominated by a solution in  $X$ .

The *average distance* metric is defined as:

$$M(X, Y) = \frac{1}{|X|} \sum_{x \in X} \min\{\|x - y\|; y \in Y\} \quad (16)$$

The metric  $M$  reflects how distant two Pareto curves are from each other.

The algorithms  $DP$  and  $DP^\epsilon$  are implemented on a Intel(R) Core(TM)2 CPU, 2.13 GHz, 3 GB of RAM computer, in a Matlab R2008b environment. All instances and software are available from the authors upon request.

### 6.1. Comparing $DP$ and $DP^\epsilon$

To illustrate the different Pareto fronts, Figure 4 depicts these corresponding to an instance with a relatively bad accuracy, namely the **R** instance with the input parameters  $t^{lim} = 500$  and  $Q = 100$ . As expected, smaller values of the worst case precision  $\epsilon$  result in a better  $\epsilon$ -Pareto front, *i.e.* closer to the real one. Furthermore, the deviation of a  $\epsilon$ -Pareto front from the optimal one increases in later iterations of  $DP$  (The size of boxes in Figure 3 increases). However, the deviation stays clearly within the worst case precision  $\epsilon$ .

Based on the results given in Table 2 and 3, we conclude that the accuracy of  $DP^\epsilon$  is excellent. For all instances, on average no more than 15% of the solutions generated by  $DP^{0.01}$  are dominated by any of  $DP$ 's solutions. Moreover, on average less than 32% of the  $DP^{0.05}$  solutions and 35% of the  $DP^{0.1}$  solutions are dominated by a  $DP$  solution. Furthermore, the distance of the 0.01-Pareto front from the optimal one is negligible (less than 2 in most cases). Moreover, even the distance of the 0.05-Pareto front and the 0.1-Pareto front from the optimal one can be consider as very small. We also observe that the fraction of  $DP^\epsilon$  solutions dominated by a  $DP$  solution increases slightly when randomness is added to the location of customers.

**Table 2** CPU and number of solutions generated.

Inst	$t^{lim}$	Q	DP		DP <sup>0.01</sup>		DP <sup>0.05</sup>		DP <sup>0.1</sup>		DP <sup>0.3</sup>	
			CPU	Nb sol	CPU	Nb sol	CPU	Nb sol	CPU	Nb sol	CPU	Nb sol
C100	500	100	1079	97	1091	97	947	68	775	46	640	23
		200	1182	97	1184	97	1080	62	814	44	670	23
	2000	100	1217	101	1232	102	1054	65	863	46	596	22
		200	1245	102	1270	102	1087	68	871	47	610	21
R100	500	100	1370	122	1376	119	1136	77	925	53	643	22
		200	1521	128	1453	122	1167	77	947	54	654	22
	2000	100	1762	143	1737	130	1273	74	999	53	596	22
		200	1555	136	1552	127	1210	76	944	53	638	22
RC100	500	100	1200	99	1195	97	935	55	748	42	541	19
		200	1300	103	1283	100	1023	56	831	40	587	19
	2000	100	1270	105	1261	98	969	56	757	40	510	20
		200	1335	107	1328	99	1036	60	809	40	575	19
C300	500	100	-	-	28770	88	25385	69	22877	46	15438	20
		200	-	-	24085	88	22208	67	20290	47	15545	19
	2000	100	-	-	-	-	-	-	21341	50	13810	24
		200	-	-	-	-	-	-	21895	51	15013	24
R300	500	100	-	-	32276	134	26670	87	21933	53	15814	23
		200	-	-	32862	129	29446	86	23005	57	16566	23
	2000	100	-	-	-	-	-	-	-	-	13628	21
		200	-	-	-	-	-	-	-	-	14152	24
RC300	500	100	-	-	-	-	27917	88	23646	48	16910	21
		200	-	-	-	-	28927	71	24002	48	17432	19
	2000	100	-	-	-	-	-	-	21600	48	15075	23
		200	-	-	-	-	-	-	23021	49	16954	23

**Table 3** The metrics C and M.

Inst	$t^{lim}$	Q	$C(DP^{0.01})$	$M(DP^{0.01})$	$C(DP^{0.05})$	$M(DP^{0.05})$	$C(DP^{0.1})$	$M(DP^{0.1})$	$C(DP^{0.3})$	$M(DP^{0.3})$
C100	500	100	0.01	0.04	0.34	29.62	0.37	76.17	0.30	358.36
		200	0.01	0.04	0.29	28.89	0.30	90.69	0.28	401.62
	2000	100	0.05	0.10	0.32	21.7	0.39	69.91	0.32	321.84
		200	0.04	0.59	0.34	21.88	0.38	60.85	0.35	315.69
R100	500	100	0.43	4.61	0.40	21.81	0.38	45.23	0.42	276.37
		200	0.45	1.43	0.39	16.29	0.42	45.77	0.39	266.63
	2000	100	0.45	0.67	0.39	27.47	0.43	54.63	0.41	281.85
		200	0.22	0.67	0.45	27.47	0.28	54.63	0.42	285.70
RC100	500	100	0.06	0.65	0.30	27.47	0.33	54.63	0.30	286.38
		200	0.17	0.88	0.30	24.81	0.25	59.38	0.21	290.64
	2000	100	0.08	1.40	0.34	24.27	0.30	62.48	0.24	321.86
		200	0.11	0.73	0.20	18.85	0.30	62.48	0.33	315.71

### 6.2. Impact of the Worst Case Precision $\epsilon$ for $DP^\epsilon$

The complexity of the  $DP^\epsilon$  algorithm increases with  $\frac{1}{\epsilon}$ . Hence, choosing smaller worst case precisions results in higher computation times. Table 2 shows the impact of  $\epsilon$  on the computation times of  $DP^\epsilon$ . We observe that with an  $\epsilon = 0.01$ , there is a small increase of about 0.1% in computation times with regard to  $DP$ . In fact, for small values of  $\epsilon$  not many solutions are deleted during the execution of  $DP^\epsilon$  and therefore it is not compensated for the trimming time. Furthermore, for  $\epsilon = 0.05$  and  $\epsilon = 0.1$ , CPU times are remarkably low, respectively 18% and 35% on average with regard to  $DP$ .

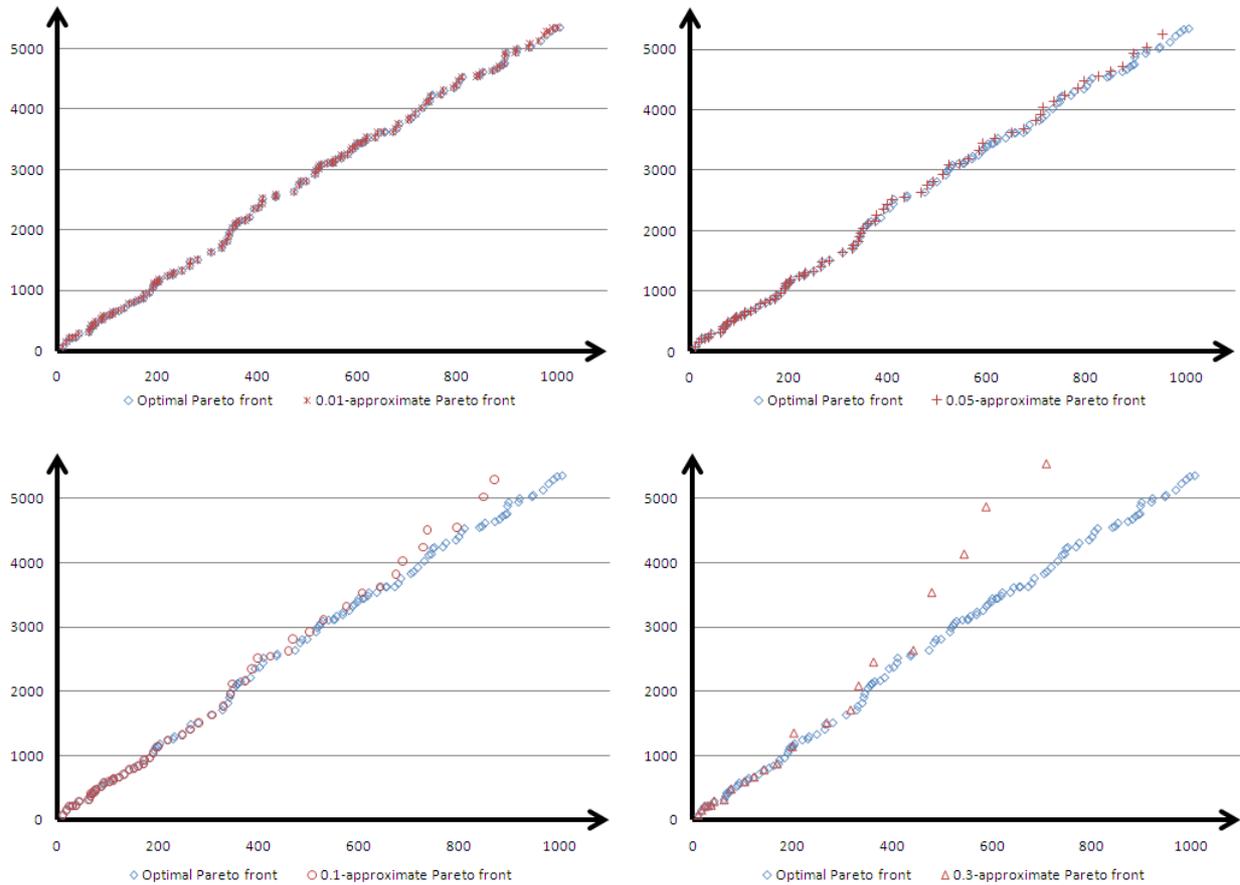


Figure 4  $\epsilon$ -Pareto fronts.

## 7. Conclusions

Multi-objective optimization problems are very challenging. They are at least as complex as their mono-objective version. Furthermore, cost and profits are hardly time-independent making the analysis and solution of multi-objective optimization problems even harder. Most existing algorithms fail to perform well in terms of both computation times and solutions quality. While exact algorithms can only deal with small problems, heuristics produce weak Pareto fronts that badly cover real Pareto fronts. In this paper, we propose a generic and flexible framework to deal with multi-objective time-dependent optimization problems. An efficient approximation based on dynamic programming is developed that generates good quality approximate Pareto fronts. Although they contain less solutions, the approximate Pareto fronts cover well the real Pareto fronts. The quality of the solutions can be decided on as the precision on each objective is an input to the algorithm and can be tuned by the decision maker. However, small precisions require more computation time. Moreover, per objective a different precision can be set. Therefore, larger errors might be allowed for less sensible objectives (*e.g.* with a flat cost structure). The approximation is tested on a time-dependent capacitated traveling salesman problem with time windows and multiple tours. Reasonably large instance with 300 customers are solved. The approximation reduces computation times considerably. Furthermore, it generates approximate Pareto fronts with few solutions, which makes the selection of a solution easier.

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## Appendix. Proofs

**Proof of lemma 1:** Note that if  $\alpha$  and  $\tilde{\alpha}$  are two vectors in the same box, then:

$$\frac{\alpha_r}{\Delta_r} \leq \tilde{\alpha}_r \leq \Delta_r \alpha_r \quad \text{for all } r \in \llbracket 1, n \rrbracket \quad (17)$$

Since  $\Delta_r$  is very close to 1,  $\alpha$  and  $\tilde{\alpha}$  are indeed very close to each other.

To prove lemma 1, we use induction on  $k \in \llbracket 1, |H| \rrbracket$ .

From (17), we conclude that lemma 1 is true for  $k = 1$ . Let us assume lemma 1 is true for  $k - 1$  and prove it for  $k$ .

Let  $I_p^k \in \{X^k\}$  for some  $p \in \llbracket 1, |X^k| \rrbracket$ . Per definition of the set  $X^k$ ,  $I_p^k$  is feasible. Hence, there exists a feasible solution  $I_u^{k-1} \in X^{k-1}$  for some  $u \in \llbracket 1, |X^{k-1}| \rrbracket$ , and some item  $i \in H$  such that  $I_p^k = I_u^{k-1} \cup \{i\}$  and for all  $r \in \llbracket 2, n \rrbracket$ :

$$t_{I_p^k} = t_{I_u^{k-1}} + c_1^i(t_0 + t_{I_u^{k-1}}) \quad \text{and} \quad f_r(I_p^k) = f_r(I_u^{k-1}) + c_r^i(t_0 + t_{I_u^{k-1}}) \quad (18)$$

On the other hand, because of the induction assumption, there exists  $\tilde{I}_v^{k-1} \in \tilde{X}^{k-1}$  for some  $v \in \llbracket 1, |\tilde{X}^{k-1}| \rrbracket$  such that for all  $r \in \llbracket 2, n \rrbracket$ :

$$t_{\tilde{I}_v^{k-1}} \leq \Delta_1^{k-1} t_{I_u^{k-1}} + (\Delta_1^{k-1} - 1)t_0 \quad \text{and} \quad f_r(\tilde{I}_v^{k-1}) \leq \max(\Delta_1^{k-1}, \Delta_r^{k-1}) f_r(I_u^{k-1}) \quad (19)$$

Furthermore,  $DP^\epsilon$  generates the vector  $[f_r(\tilde{I}_v^{k-1}) + c_r^i(t_0 + t_{\tilde{I}_v^{k-1}})]_{r=1}^{r=n}$  in the  $k$ th step. This vector point might be removed after trimming. However some vector  $[f_r(\tilde{I}_q^k)]_{r=1}^{r=n}$ , located in the same box, should be left. From (17), we obtain for all  $r \in \llbracket 2, n \rrbracket$ :

$$t_{\tilde{I}_q^k} \leq \Delta_1(t_{\tilde{I}_v^{k-1}} + c_1^i(t_0 + t_{\tilde{I}_v^{k-1}})) \quad \text{and} \quad f_r(\tilde{I}_q^k) \leq \Delta_r(f_r(\tilde{I}_v^{k-1}) + c_r^i(t_0 + t_{\tilde{I}_v^{k-1}})) \quad (20)$$

Because of the FIFO and the induction assumptions, we have:

$$t_0 + t_{\tilde{I}_q^k} \leq \Delta_1(\Delta_1^{k-1}(t_0 + t_{I_u^{k-1}}) + c_1^i(\Delta_1^{k-1}(t_0 + t_{I_u^{k-1}}))) \quad (21)$$

and for all  $r \in \llbracket 2, n \rrbracket$ :

$$f_r(\tilde{I}_q^k) \leq \Delta_r(\Delta_r^{k-1} f_r(I_u^{k-1}) + c_r^i(\Delta_1^{k-1}(t_0 + t_{I_u^{k-1}}))) \quad (22)$$

Using assumption 2, we obtain:

$$t_0 + t_{\tilde{I}_q^k} \leq \Delta_1^k(t_{I_u^{k-1}} + c_1^i(t_0 + t_{I_u^{k-1}})) + \Delta_1^k t_0 \quad (23)$$

and for all  $r \in \llbracket 2, n \rrbracket$ :

$$f_r(\tilde{I}_q^k) \leq \Delta_r(\Delta_r^{k-1} f_r(I_u^{k-1}) + \Delta_1^{k-1} c_r^i(t_0 + t_{I_u^{k-1}})) \quad (24)$$

Hence, for all  $r \in \llbracket 2, n \rrbracket$ :

$$t_{\tilde{I}_q^k} \leq \Delta_1^k t_{I_p^k} + (\Delta_1^k - 1)t_0 \quad \text{and} \quad f_r(\tilde{I}_q^k) \leq \max(\Delta_1^k, \Delta_r^k) f_r(I_p^k) \quad (25)$$

**Proof of theorem 1:**  $DP^\epsilon$  and  $DP$  have at most  $|H|$  iterations. Let us assume they converge after  $k$  iterations ( $k \leq |H|$ ). According to lemma 1, For all  $p \in \llbracket 1, |X^k| \rrbracket$ , there exists  $q \in \llbracket 1, |\tilde{X}^k| \rrbracket$  such that for all  $r \in \llbracket 2, n \rrbracket$ :

$$t_{\tilde{I}_q^k} \leq \Delta_1^k t_{I_p^k} + (\Delta_1^k - 1)t_0 \quad \text{and} \quad f_r(\tilde{I}_q^k) \leq \max(\Delta_1^k, \Delta_r^k) f_r(I_p^k) \quad (26)$$

The sequences  $\left(1 + \frac{\epsilon_r}{2|H|}\right)^{|H|}$  are increasing in  $|H|$  and converges to  $e^{\frac{\epsilon_r}{2}}$ . Hence, for every  $|H| \geq 1$ :

$$\left(1 + \frac{\epsilon_r}{2|H|}\right)^{|H|} \leq e^{\frac{\epsilon_r}{2}} \quad (27)$$

Furthermore, for  $0 < \epsilon_r < 1$  we have:

$$e^{\frac{\epsilon_r}{2}} \leq 1 + \epsilon_r \quad (28)$$

Therefore, we have:

$$t_{\tilde{I}_q^k} \leq (1 + \epsilon_1) t_{I_p^k} + \epsilon_1 t_0 \quad \text{and} \quad f_r(\tilde{I}_q^k) \leq (1 + \max(\epsilon_1, \epsilon_r)) f_r(I_p^k) \quad (29)$$

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