

Ant Colony Optimization on scheduling problem with job values given as a power function of their completion times

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Abstract— This paper deals with the problem of scheduling jobs on the identical parallel machines, where job values are given as a power function of the job completion times. Minimization of the total loss of job values is considered as criterion. This paper establishes the computational complexity of the problem by the application of Ant Colony Optimization on the problem. Strong NP hardness of its general versions and NP hardness of its single machine case. Moreover, some special cases of the problem in the polynomial time. Finally the project construct and experimentally test branch and bound algorithm along with some elimination properties improving its efficiency. In order to solve the significant real life problems a lot of new scheduling problems have been formulated. The problems with resource allocation and the problem with deteriorating jobs are two kinds of examples of such scheduling.

Index Terms— Scheduling, Scheduling Algorithm, Parallel Processor, Ant Colony, ACO, MMAS, job values, loss function

1 INTRODUCTION

The aim of this scheduling is to describe and solve a scheduling problem where job values (or losses of jobs values) changes during their execution and are described as a power function of the job completion times. In this paper we introduce a new non decreasing power model of loss of job values and formulate a problem of minimization of the total job values loss which can be used to solve the problem of establishing an order of object renovation. Here we introduce a new model of job value, defined as a difference between initial job value and non decreasing power function loss. [1]

Another new kind of the scheduling problems that we had studied in this paper. Namely the in which processing times of all the jobs are some values which are fixed in advance and constant during optimization process but their values deteriorate over time. The process description of the problem can be illustrated by an application example, which characterizes the utilization process of the components from some used up computers. Therefore assumption that there is given a set of some used up computer which cannot be used any more because their further utilization is connected with a high risk of a breakdown.

there are some application which require faster processors or simply of their components are already broken. However some of their components (e.g monitors, floppy disks drives, network drives, power suppliers) can be utilized as a spare parts in some other computers. Thus the problem of disassembling computers into their components appears nearest

can be shown that the values of the computer component decreases over time as a power function of the speed of this different for the particular component. In disassembling process we are interested in the component values which are determined at the moments at they are available for utilization. The component is ready to be used after it is completely removed from one computer and its proper functionality is confirmed. Thus the order in which the computer will be disassembled has a significant influence on the total profit, i.e. the sum of the component values. Therefore, maximization of the total component value is considered as an optimization criterion [2]

2 SCHEDULING

A schedule is a tangible plan or document, such as a bus or a class schedule. A schedule usually tells us when things are supposed to happen; it shows us a plan for the timing of certain activities and answers the question, "If all goes well, when will a particular event take place?"

The aim of this scheduling is to describe and solve a scheduling problem where jobs values (or losses of jobs values) changes during their execution and are described as a power function of the job completion times. [3]

2.1 Notations

Regardless of its nature, every scheduling problem S can be formulated as a quadruple, $S = (J, M, P, L)$, where J is a set of non preemptive jobs immediately available for processing time 0. M is a set of identical parallel machines entities that will perform the available jobs. P is the processing time of the job, where $p_j > 0$ and L is the loss function as $l_j(t)$ characterizing the loss of its value at time t . [4]

Processing time (p_{ij}) - The p_{ij} represents the processing time of job j on machine i . The subscript i is omitted if the processing time of job j does not depend on the machine or if job j

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is only to be processed on one given machine.

Exponential loss rate (aj)-This is the loss rate denoted by the a_j and $a_j > 0$ and it is associated with the job j_i performing on machine M .

Proportional loss rate (wj)-This is another loss rate denoted by the w_j and $w_j > 0$ and associated with the job j_i performing on machine M .

Loss function(lj(t))-This is loss rate function characterizing the loss of its value at time t . The loss rate function is a non decreasing power function of time. Moreover the loss rate is calculated only at the completion time C_j .

Three basic pieces of information that help to describe jobs in the single-machine case are:

- Processing time (p_j) the amount of processing required by job j
- Exponential loss rate (a_j) the exponential loss rate of job j
- Proportional loss rate (w_j) The proportional loss rate of job j
- Completion time (C_j) The time at which the processing of job j is finished.

2.2 Scheduling Algorithm and its Complexity

A useful perspective on the relation of scheduling problems and their solution techniques comes from developments in a branch of computer science known as complexity theory. The notion of complexity refers to the computing effort required by a solution algorithm. Computing effort is described by order-of-magnitude notation. For example, suppose we use a particular algorithm to solve a problem of size n . (Technically, n denotes the amount of information needed to specify the problem.) The number of computations required by the algorithm is typically bounded from above by a function of n . If the order of magnitude of this function is polynomial, as n gets large, then we say the algorithm is polynomial. For instance, if the function has order of magnitude n^2 , denoted $O(n^2)$, then the algorithm is polynomial. On the other hand, if the function is $O(2^n)$, then the algorithm is non polynomial (in this case, exponential). Other things being equal, we prefer to use a polynomial algorithm because as n grows large, polynomial algorithms are ultimately faster. [5]

3 RELATED WORK

This paper deals with the problem of scheduling jobs on the identical parallel machines, where job values are given as a power function of the job completion times. Minimization of the total loss of job values is considered as criterion. This paper establishes the computational complexity of the problem of strong NP hardness of its general versions and NP hardness of its single machine case. Moreover, some special cases of the problem in the polynomial time. Finally the paper construct and experimentally test branch and bound algorithm along with some elimination properties improving its efficiency. In order to solve the significant real life problems a lot of new scheduling problems have been formulated. The problems

with resource allocation and the problem with deteriorating jobs are two kinds of examples of such scheduling.

3.1 Mathematical Formulation

There are given a set of m identical parallel machines There are given a set of m identical parallel machines $M = \{M_1, \dots, M_m\}$ and a set of n independent and non preemptive jobs $J = \{J_1, \dots, J_n\}$ immediately available for processing at time 0. Each job $J_i \in J$ is characterized by its processing time $p_i > 0$ and a loss function $l_j(t)$ characterizing the loss of its value at time t . The loss function is given by non decreasing power function of time.

$$L_j(t) = w_j t^{a_j} \tag{1}$$

$$TLV(\Pi) = \sum_{i=1}^m \sum_{j=1}^{n_i} l_{\pi_i(j)}(C_{\pi_i(j)}) = \sum_{i=1}^m \sum_{j=1}^{n_i} w_{\pi_i(j)} C_{\pi_i(j)}^{a_{\pi_i(j)}} \rightarrow \min$$

The decrease of the computer components values can be modeled by the following time function $v(t) = \omega t^{-\alpha}$ where the values of α and ω were established experimentally for different kinds of the processors, RAM modules, hard disks, monitors etc. Therefore the values of the computer components decrease fast at the beginning when the components are quite new, and the values decrease slowly when the component are quite old. Thus such situation can be also modeled by the scheduling problem considered in the present paper as follows. The job is to recover a given component from some used up computer. The value $v_j(t)$ of job J_j decreases over time and can be decreased by the following expression.

$$V_j(t) = v_j^0 - l_j(t) = v_j^0 - w_j t^{a_j} \tag{2}$$

Where, $v_j^0 > 0$ is its initial value (since $l_j(0)$ and $a_j \in (0, 1)$ (the function $v_j(t)$ is convex for the concave function $l_j(t)$ given by Expression. The specific values of v_j^0, w_j , and a_j for particular components can be established experimentally, taking into account the appropriate values of ω and α . In such a case, we can also assume that the job values $v_j(t)$ is positive at any moment t of the optimization process. Thus, the following condition should be satisfied $v_j(t) = v_j^0 - w_j t^{a_j} > 0$ for t belongs $[0, \sum_{j=1}^n p_j]$. The problem is to find such a solution Π that the sum of job values calculated at their completion times is maximal, i.e.,

$$\left(\sum_{i=1}^m \sum_{j=1}^{n_i} v_{\pi_i(j)}(C_{\pi_i(j)}) \right) = \sum_{i=1}^m \sum_{j=1}^{n_i} v_{\pi_i(j)}^0 - \sum_{i=1}^m \sum_{j=1}^{n_i} l_{\pi_i(j)}(C_{\pi_i(j)}) \rightarrow \max, \tag{3}$$

Since, $\sum_{i=1}^m \sum_{j=1}^{n_i} v_{\pi_i(j)}^0 = \sum_{j=1}^n v_j^0$ is constant, thus, maximization of the sum of job values is equivalent to the minimization of the total loss of jobs values.

4 ALGORITHM FOR PARALLEL PROCESSOR

In this problem we have given a set of m identical parallel machines and a set of n independent and non-preemptive jobs immediately available for processing at time 0. Each job J_j is belongs to J characterised by its processing time and a loss function.

Firstly we assign the jobs to the machines according to their processing time. The job j having the minimum processing time is assigned to the first machine and as goes on. The number of jobs and number of machines are randomly generated

by the user. And we generate the sequence number with processing time.

And secondly we generate another sequence number by assigning the jobs to the machines on the basis of their loss function. The loss function is denoted by $I_j(t)$ characterizing the loss of its value at time 0. The loss function is denoted by non decreasing power function. [6]

For this section we describe the most efficient algorithms and the result of the performed experimental analysis. All the algorithms A1LST-A3LST are based on the list strategy but they use three different priority dispatching rules to obtain an order of jobs in an input list. [7]

4.1 Algorithm AiLST

Step 1 Construct the list of jobs L , according to the algorithm A1, $i \in \{1, 2, 3\}$.

Algorithm A1 and A2 construct solutions by sequencing all the jobs in the non increasing order of a_j and $(a_j \cdot w_j / p_j)$ respectively).

Algorithm A3

Step1. Set $\pi = 0$, $U = \{1 \dots n\}$ and $C = \sum_j p_j$.

Step2. Find a job $j \in U$ with the minimal value of $w_j C$ and move it from U at the beginning of π . Set $C = C - p_j$.

Step3. If $U \neq \emptyset$ then go to step 2 the job order is given by π .

The computational complexity of algorithm A1 and A2 is equal to $O(n \log n)$, while the complexity of algorithm A3 is equal to $O(n^2)$. Finally in order to improve the job sequence on the machines the following procedure is launched when the jobs are scheduled. [8] [9]

4.1.1 Algorithm DoubleSwap

Step1: For $j=1, \dots, n-1$ swap jobs in positions j and $j+1$ if

$$I_{\pi(j)} (C_{\pi(j)}) + I_{\pi(j+1)} (C_{\pi(j+1)}) > I_{\pi(j+1)} (C_{\pi(j-1)} + P_{\pi(j+1)}) + I_{\pi(j)} (C_{\pi(j+1)}) \quad (4)$$

Step2: For $j=n, \dots, 2$ swap jobs in positions j and $j-1$ if

$$I_{\pi(j)} (C_{\pi(j)}) + I_{\pi(j-1)} (C_{\pi(j-1)}) > I_{\pi(j-1)} (C_{\pi(j)} + P_{\pi(j)}) + I_{\pi(j)} (C_{\pi(j-2)} + P_{\pi(j)}) \quad (5)$$

5 ANT ALGORITHM

5.1 Algorithm used in ACO

There is a pool of jobs having constant processing time (prt_j) and initial job value (C_j) but varying deterioration (d_j) depending on completion time. The algorithm for parallel processor scheduling using Ant Colony Optimization may be represented as follows:

1. Initialize the pheromone $ph_{j,i}$ with initial pheromone val-

ue (τ_0 / τ_0) . Pheromone $ph_{j,i}$ is accumulated between jobs and deterioration interval. [10]

2. All ants start making its tour.

a) Ants select a job first. Selection of job depends on two parameter- (1) processing time and (2) next deterioration, which is drive by a probability distribution function. Job j is selected for deterioration interval i according to formula [11]

$$j = \max (ph_{j,i}, i * d_i^{1/2} / prt_j) \quad \text{if } q < q_0$$

$$j = p(j) = (ph_{j,i} * d_i^{1/2} / prt_j) / \sum (ph_{j,i} * d_i^{1/2} / prt_j) \quad \text{otherwise}$$

Where, d_i = deterioration for interval i
 prt_j = processing time for job j (6)

b) Ants select a processor (p) for selected job (j) according to formula

$$p = \max [(t_{pc} - t_{pi-1}) / (t_{pi} - t_{pi-1}) / i_p]$$

Where, t_{pc} = completion time on processor p

t_{pi} = end point of interval i for processor p

i_p = interval for processor p

c) Update the pheromone (local updation)

$$ph_{j,pi} = (1 - \rho) * ph_{j,pi} + \rho * \tau_0$$

d) Steps i to iii is repeated for all jobs.

3. Select the best tour among tours of all ants.

a) Best tour with maximum job value is selected.

b) Update the pheromone (global updation)

$$ph_{j,i} = (1 - \alpha) * ph_{j,i} + \alpha * (i/D)$$

where, D = total deterioration in best tour

Repeat steps 1 to 3 for each iteration

5.2 Steps for Solve the problem by ACO

The basic steps, which are to be followed for solving the problem by ACO, are as follows [12]

- Represent the problem in the form of sets of components and transitions, or by a set of weighted graphs, on which ants can build solutions
- Define the meaning of the pheromone trails
- Define the heuristic preference for the ant while constructing a solution
- If possible implement a efficient local search algorithm for the problem to be solved.
- Choose a specific ACO algorithm and apply to problem being solved
- Tune the parameter of the ACO algorithm

5.3 Ant System

Ant System is the first ACO algorithm proposed in the literature. Its main characteristic is that, at each iteration, the pheromone values are updated by all the m ants that have built a solution in the iteration itself. The pheromone τ_{ij} , associated

with the edge joining cities i and j , is updated as follows:[13]

$$\tau_{ij} \leftarrow (1-\rho) \cdot \tau_{ij} + \sum_{k=1}^m \Delta_{ij}^k$$

where ρ is the evaporation rate, m is the number of ants, and $\Delta\tau_{ij}^k$ is the quantity of pheromone laid on edge (i, j) by ant k :

$$\Delta\tau_{ij}^k = \begin{cases} Q/L_k & \text{if any } k \text{ used edge } (i, j) \text{ in its tour,} \\ 0 & \text{otherwise,} \end{cases}$$

where Q is a constant, and L_k is the length of the tour constructed by ant k . In the construction of a solution, ants select the following city to be visited through a stochastic mechanism. When ant k

is in city i and has so far constructed the partial solution s^p , the probability of going to city j is given by

$$p_{ij}^k = \begin{cases} \frac{\tau_{ij}^\alpha \eta_{ij}^\beta}{\sum_{i,l \in N(i^p)} \tau_{ij}^\alpha \eta_{ij}^\beta} & \text{if } c_{ij} \in N(s^p) \\ 0 & \text{otherwise,} \end{cases} \quad (7)$$

where $N(s^p)$ is the set of feasible components; that is, edges (i,l) where l is a city not yet visited by the ant k . The parameters α and β control the relative importance of the pheromone versus the heuristic information η_{ij} , which is given by:

$$\eta_{ij} = 1/d_{ij}, \text{ where } d_{ij} \text{ is the distance between cities } i \text{ and } j. [14]$$

5.3.1 MIN-MAX Ant System (MMAS)

This algorithm is an improvement over the original Ant System. Its characterizing elements are that only the best ant updates the pheromone trails and that the value of the pheromone is bound. The pheromone update is implemented as follows: [15]

$$\tau_{ij} \leftarrow [(1-\rho) \cdot \tau_{ij} + \Delta\tau_{ij}^{\text{best}}]^{\tau_{\min}^{\max}} \quad (8)$$

where τ_{\max} and τ_{\min} are respectively the upper and lower bounds imposed on the pheromone; the operator $[x]_b^a$ is defined as:

$$[x]_b^a = \begin{cases} a & \text{if } x > a, \\ b & \text{if } x < b, \\ x & \text{otherwise;} \end{cases}$$

And $\Delta\tau_{ij}^{\text{best}}$ is:

$$\Delta\tau_{ij}^{\text{best}} = \begin{cases} 1/L_{\text{best}} & \text{if } (i, j) \text{ belongs to best tour,} \\ 0 & \text{otherwise,} \end{cases} \quad (9)$$

where L_{best} is the length of the tour of the best ant. This may be (subject to the algorithm designer decision) either the best tour found in the current iteration—iteration-best, L_{ib} —or the best solution found since the start of the algorithm—best-so-far, L_{bs} —or a combination of both. Concerning the lower and upper bounds on the pheromone values, τ_{\min} and τ_{\max} , they are typically obtained empirically and tuned on the specific problem considered. Nonetheless, some guidelines have been provided for defining τ_{\min} and τ_{\max} on the basis of analytical considerations. [16]

5.4 Ant Colony System

The most interesting contribution of ACS is the introduction of a local pheromone update in addition to the pheromone update performed at the end of the construction process (called offline pheromone update). The local pheromone update is performed by all the ants after each construction step. Each ant applies it only to the last edge traversed: [17]

$$\tau_{ij} = (1 - \phi) \cdot \tau_{ij} + \phi \cdot \tau_0, \quad (10)$$

where $\phi \in (0, 1]$ is the pheromone decay coefficient, and τ_0 is the initial value of the pheromone.

The main goal of the local update is to diversify the search performed by subsequent ants during iteration: by decreasing the pheromone concentration on the traversed edges, ants encourage subsequent ants to choose other edges and, hence, to produce different solutions. This makes it less likely that several ants produce identical solutions during one iteration. The offline pheromone update, similarly to MMAS, is applied at the end of each iteration by only one ant, which can be either the iteration-best or the best-so-far. However, the update formula is slightly different:

$$\tau_{ij} \leftarrow \begin{cases} (1 - \rho) \cdot \tau_{ij} + \rho \cdot \Delta\tau_{ij} & \text{if } (i, j) \text{ belongs to best tour,} \\ \tau_{ij} & \text{otherwise} \end{cases} \quad (11)$$

As in MMAS, $\Delta\tau_{ij} = 1/L_{\text{best}}$, where L_{best} can be either L_{ib} or L_{bs} . Another important difference between ACS and AS is in the decision rule used by the ants during the construction process. [18][19]

In ACS, the so-called pseudorandom proportional rule is used: the probability for an ant to move from city i to city j depends on a random variable q uniformly distributed over $[0,1]$, and a parameter q_0 ; if $q \leq q_0$, then

$$p_{ij}^k = \begin{cases} \frac{\tau_{ij}^\alpha \eta_{ij}^\beta}{\sum_{i,l \in N(i^p)} \tau_{ij}^\alpha \eta_{ij}^\beta} & \text{if } c_{ij} \in N(s^p) \\ 0 & \text{otherwise,} \end{cases} \quad (12)$$

6 RESULTS

This paper solves the problem of parallel processor scheduling. It is analyzed that the final results of parallel processor scheduling using ant colony optimization technique are better than the results of heuristics that have been already developed earlier.

6.1 Results of Heuristics

Table 6.1 Result Of Scheduling For 100 jobs

	VALUE 1	VALUE 2	VALUE 3	VALUE 4	VALUE 5
A1	7.8965	8.9764	7.1234	9.8765	6.7453
A2	8.8989	7.8976	8.9907	5.4532	6
A3	3.2412	2.1342	4.3321	2.1311	4.6754
DA1	9.8765	7.8564	7.3521	8.3409	6.4409
DA2	7.8976	6.5643	5.3421	7.3452	8.9856
DA3	3.2143	2.1045	3.1232	1.5643	4.6867
ACS1	2.1321	2.1001	1.9987	1.3342	3.0934

Table 6.2 Result Of Scheduling For 200 jobs

	VALUE 1	VALUE 2	VALUE 3	VALUE 4	VALUE 5
A1	7.8675	8.4532	7.8889	5.3421	9.5632
A2	8.8956	6.7843	8.9786	6.2311	6.3444
A3	4.1234	3.1232	3	1.9906	3.0568
DA1	9.4532	6.5543	8.4251	9.1765	6.3245
DA2	8.9234	8.1232	7.5643	4.1242	8.6421
DA3	3.1232	3.1096	2.7865	1.9874	2.3421
ACS1	2.1342	1.2321	2.3339	1.0078	2.2222

Table 6.3 Result Of Scheduling For 300 jobs

	VALUE 1	VALUE 2	VALUE 3	VALUE 4	VALUE 5
A1	4.5632	6.6543	9.4532	6.6665	7.3421
A2	6.7654	6.7888	7.5643	7.5643	7.3421
A3	5.6754	3.2213	5.6543	4.5632	4.5621
DA1	8.8945	8.6554	8.7865	9.7854	8.4654
DA2	9.1123	6.3452	5.6754	7.7644	5.6754
DA3	4.4453	3.2218	4.3211	2.1068	3.5632
ACS1	4.4005	2.4321	3.3305	2.1005	2.5612

* **RED COLOR:** MAXIMUM LOSS VALUE
SKY BLUE COLOR: MINIMUM LOSS VALUE

7 GRAPHS

In these graphs

Y Axis represents the loss values of jobs

X Axis represents the algorithm used

Value 1 shows the result of first execution of the entire algorithm

Value 2 shows the result of second execution of the entire algorithm

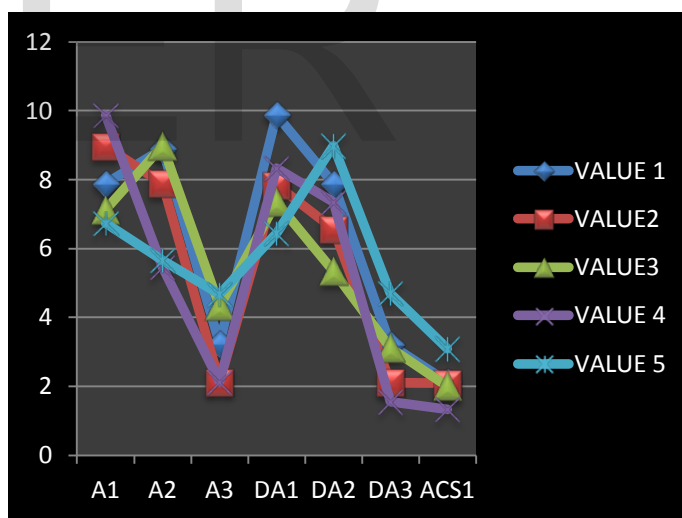
Value 3 shows the result of third execution of the entire algorithm

Value 4 shows the result of fourth execution of the entire algorithm

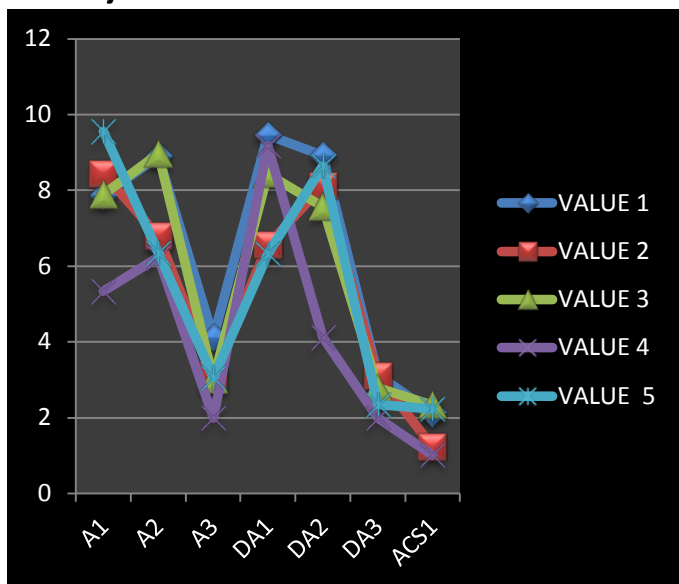
Value 5 shows the result of fifth execution of the entire algorithm

Here the number of processor are taken as 5. And value1, value2, value3, value4 and value5 are the different values they comes out at each execution. And with the help of graph its clear that the ant colony optimization gives the minimum loss values of the jobs.

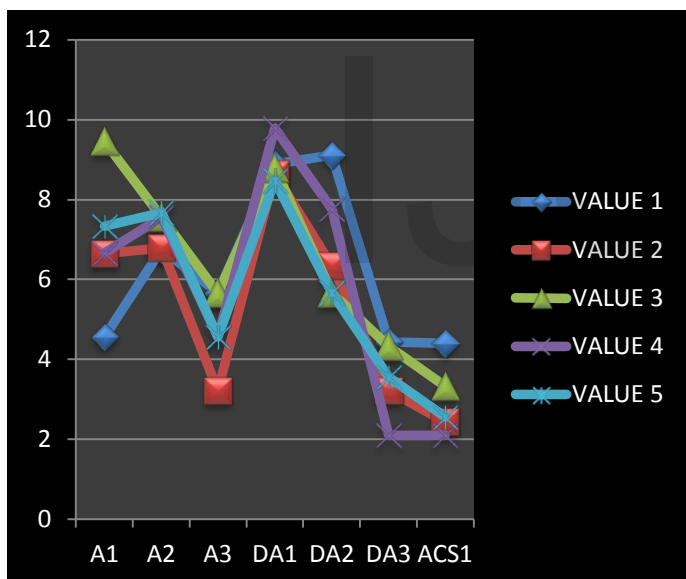
7.1 Resultant graph of multi processor scheduling for 100 jobs



7.2 Resultant graph of multi processor scheduling for 200 jobs



7.3 Resultant graph of multi processor scheduling for 300 jobs



7 CONCLUSION

In this paper the parallel machines scheduling problem with time dependent losses of job values has been investigated. It has been shown that the general version of the problem is strongly NP-hard and its single machine case is NP-hard. Therefore, in order to solve the problem the use of branch and bound algorithm, supported by some elimination properties which improve its efficiency. It delivers optimal solution of the instances of 25 jobs in a few minutes on the average, using some heuristics algorithm with low computational complexities – $O(n \log n)$ and $O(n^2)$. The experimental analysis has revealed, that their efficiencies strongly depend on the instance size and the values of the problem parameters. However, the analysis has delivered also a manner of choosing an appropri-

ate algorithm in order to obtain near solutions.

At last comparison is made with ant colony algorithms and we conclude that ant colony gives more optimized result as compare to heuristics.

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