

# Enhancing the Schedulability of Real-Time Heterogeneous Networks of Workstations (NOWs)

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**Abstract**—This paper proposes a Real-Time Duplication-Based Algorithm (RT-DBA) for scheduling precedence-related periodic tasks with hard deadlines on networks of workstations (NOWs). We have utilized selective subtask duplication that enables some tasks to have earlier start times, which enables additional tasks (and, hence, task sets) to finish before their deadlines, thereby increasing the schedulability of a real-time application. We strongly believe that duplication can be used as a tool for obtaining a better quality of service (QoS) from the real-time heterogeneous system, and this is our major contribution. We have taken both the computation and the communication heterogeneities into account while modeling such a system. Both data and control dependencies between the tasks have also been considered. Our algorithm exhibits scalability, fully exploits the underlying parallelism, and is capable of scheduling an application, even if the available number of processors is less than the required number of processors. Based on extensive simulation studies, we observe that RT-DBA offers an enhanced success ratio as compared to other scheduling schemes when communication is a dominant factor.

**Index Terms**—Favorite predecessor, heterogeneous network of workstations, periodic tasks, precedence constraints, real-time scheduling, subtask compaction, subtask duplication.



## 1 INTRODUCTION

IN addition to their correctness [1], a real-time system is defined as one which makes results available in a timely manner, due to decreasing costs of processors, real-time applications are being implemented and run on networks of workstations (NOWs). Real-time requirements may necessitate intensive communication [2], [3] in these NOWs. In order to provide a cost-effective solution while satisfying desired performance requirements, such multicomputer systems may very often be heterogeneous. This means that they usually consist of both programmable processors and dedicated hardware systems for enhanced performance, and heterogeneity can be present within these two domains. For example, programmable processors can be general-purpose microprocessors, microcontrollers, or specialized digital signal processors (DSPs). Therefore, each task could have potentially a different execution time if it were to be executed on a microprocessor, a microcontroller, or a DSP system, and it is imperative to devise efficient scheduling algorithms for extracting a better performance on a heterogeneous multiprocessor platform [4], [5]. Real-time scheduling algorithms can be broadly categorized into two types—static algorithms [6], [7] and dynamic algorithms [8], [9]. Real-time scheduling on multiprocessors has been proven to be NP-complete, even for a very simplified task model such as two processors and unit task execution time

[10]. This has encouraged various researchers to propose heuristics for solving the scheduling problem [2], [3], [4], and we also adopt a similar approach in this paper.

## 2 MOTIVATION

As a major contribution, we are proposing a way to select communicating subtasks for duplication so that there is no need for interprocessor communication, thereby eliminating such precedence delays. Hence, a larger number of tasks and, consequently, a larger task set are able to meet the deadlines. We strongly believe that our proposed algorithm can increase the schedulability of real-life applications such as audio/video on demand wherein the server relays audio or video information to the user's computer by transmitting a certain number of frames per second. By employing our proposed subtask duplication approach, many tasks can be processed relatively faster and therefore enable a larger number of frames to be transmitted within the same time duration, and the server could possibly provide a better QoS. By employing task duplication, a larger number of subtasks can be scheduled within the deadline window, and hence, the utilization factor of the hard real-time system can be improved. This means that a larger number of task sets can meet their deadlines. Therefore, our proposed scheme can be very useful for traditional hard real-time military and command and control applications.

The aim of this research is twofold: 1) to propose a scheduling algorithm (Real-Time Duplication-Based Algorithm (RT-DBA)) that employs duplication in order to improve the schedulability of a hard real-time system on NOWs and 2) to ensure that a real-time schedule is generated, even if an adequate number of processors are not available, and, in doing so, demonstrate downward scalability. The proposed scheme works in the following

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stages: 1) The tasks are scheduled based on the rate-monotonic algorithm [11], i.e., the tasks with lower periods are assigned higher priority and are scheduled first. The proposed scheme computes a set of mathematical parameters for each subtask and then generates clusters of subtasks based on these quantities. 2) It assigns each cluster to an appropriate heterogeneous processor and schedules the messages on appropriate links. Based on the initial number of clusters, the critical number of processors (or CNP) is obtained that leads to reduced computation time. We assume that  $CNP = \text{initial number of clusters}$ . Another quantity of interest is the number of processors that are available to schedule the real-time application. This quantity is called the available processors (or AP). 3) AP is compared with CNP and one of the following three actions is taken:

1. If  $AP < CNP$ , the proposed scheme demonstrates downward scalability and invokes the cluster compaction procedure to schedule all generated clusters on to the fewer number of processors.
2. If  $AP = CNP$ , the proposed scheme schedules the clusters to their appropriate processors.
3. If  $AP > CNP$ , the proposed scheme exploits the underlying parallelism and employs subtask duplication to reduce the final finish time of the application, thereby increasing its success ratio (SR).

The rest of this paper is organized as follows: In Section 3, we discuss the related work. Section 4 describes the task model along with some terminology. The proposed algorithm with pseudocode is elaborated in Section 5. The simulation results are described in Section 6. Finally, Section 7 concludes the paper. A running trace of the proposed algorithm, as well as the theoretical analysis, is shown in the Appendix.

### 3 RELATED WORK

Real-time scheduling on multiple processors has been addressed using various approaches, basically extending the single-processor model, e.g., [4]. In another approach, real-time tasks are grouped into clusters based on communication delays, each cluster is assigned to a processor, e.g., [12], and several heuristics have been proposed to schedule complex periodic tasks on homogeneous multiprocessors. Hence, the algorithms in [12] cannot be directly adopted for heterogeneous systems. Moreover, unlike the schemes discussed in [12], our scheme is scalable. A replication-based scheme that improves the schedulability of a real-time application has been discussed in [13]. However, the scheme makes the naïve assumption that there is only one communication channel in the network. Essentially, the scheme is based on dynamic programming and can potentially have a very large search space. In contrast, our scheme considers multiple channels and is a heuristic-based approach.

Scheduling of real-time tasks on a uniprocessor has been a well-researched problem [11], [14]. Reward-based scheduling has been discussed in [15]. Extensive research has also been carried out for the homogeneous multiprocessor model [8], [9], [16], [17], [18]. Real-Time Scheduling on Heterogeneous Multiprocessors has also been a topic of extensive research [4], [5], [6], [11], [20], [21], [22], [23], [24], [25]. Recent papers [4], [5] extend the popular EDF algorithm for heterogeneous multiprocessors. One of the

limitations is that they assume independence between various subtasks.

A reliability-driven scheduling scheme for tasks with precedence constraints on heterogeneous multiprocessor systems has been discussed in [21]. In our previous work [22], [23], we have proposed several scheduling algorithms for heterogeneous NOWs. Omari et al. [26] present an adaptive scheme that schedules fault-tolerant soft tasks on multiprocessors. A replication-based technique for multiprocessor scheduling has been discussed in [27]. A feedback control algorithm called EUCON has been proposed in [28]. In [29], the authors propose a solution to the problem of scheduling a set of precedence-related tasks on variable-speed processors. The roots of our scheme can be found in [30], [31], and [32], which do not consider real-time tasks. In [33] and [34], the authors presented a hardware-software cosynthesis system that partitions and schedules an embedded system consisting of multiple periodic tasks. In [35] and [36], the authors proposed algorithms for scheduling periodic tasks on identical multiprocessors. In this research, we assume processors of varying speeds and configurations. Task duplication has been suggested for I/O-intensive tasks in a cluster-based system [37] so that frequently copying of huge amounts of data can be avoided. Task duplication has also been used for improving the schedule length of the parallel application [38], [39], albeit for non-real-time systems.

### 4 PROBLEM SPECIFICATION

Various mathematical notations used for the real-time task parameters are shown in Table 1. The input to the algorithm is the real-time application ( $T$ ), which consists of a set of independent periodic tasks. Each periodic task ( $t_i \in T$ ) can be further subdivided into a set of precedence-related subtasks [24]. The real-time system also consists of the set of available heterogeneous processors,  $AP = \{p_1, p_2, p_3, \dots, p_n\}$ . A periodic task can be represented by a Directed Acyclic Graph (DAG) [16], [24], with a subtask  $s(i, j, k)$  of a task  $t_i$  being represented as a node in the DAG, where  $i$  denotes the periodic task ID,  $j$  denotes the subtask ID within a periodic task, and  $k$  stands for the invocation ID of a subtask. The worst-case execution time of a subtask is given by  $\mu(s(i, j, k), p_z)$ , where  $p_z$  is an element of an array  $P$  ( $p_z \in P$ ). The communication cost between two subtasks  $s(i, j, k)$  and  $s(i, l, k)$  of a task  $t_i \in T$  is given by  $c((s(i, j, k), s(i, l, k)))$ . The length of the longest path from the subtask to the exit node is given by  $p\text{-}e(s(i, j, k))$ . The set of clusters is given by  $CL = \{cl_1, cl_2, cl_3, \dots\}$ . The edges between the subtasks represent the data dependency between them in that a subtask cannot start till it has received the results from all of its predecessor tasks. The control dependency between tasks can be modeled based on the system functionality if certain conditions exist in the DAG. These conditions are represented on the edges of the DAG such that the transmission on such an edge will take place only if the associated condition is satisfied.

An example in Fig. 1 has two independent real-time tasks  $t_1$  and  $t_2$ . The edges between the tasks represent the data flow (solid edges) and the control flow (dotted edges). The number along each edge represents the volume of data that needs to be transmitted if the two tasks are allocated to

TABLE 1  
Mathematical Notations Used for Task Parameters

Notation	Task Parameter
T	The real-time application (set of independent periodic tasks)
AP	The set of available processors
C	Set of conditions for modeling control dependencies between the tasks
Q	Data structure queue used to select subtask clusters
PL	List of periodic tasks in increasing order of periods
i	Task id
j	Subtask id
k	Invocation id
$t_i$	Periodic task (such that $t_i \in T$ ) corresponding to a DAG
$pd_i$	Period of task $t_i$
$d_i$	Deadline of task $t_i$
$S(t_i)$	Set of subtasks belonging to task $t_i$
$s(i,j,k)$	$k^{\text{th}}$ invocation of $j^{\text{th}}$ subtask of task $t_i$
$p_z$	Processor to which current subtask is assigned
$\mu(s(i,j,k), p_z)$	Execution time of subtask $s(i,j,k)$ provided it will execute on processor $p_z$
$c(s(i,j,k), s(i,l,k))$	Cost of communication between subtasks $s(i,j,k)$ and $s(i,l,k)$
$p\_e(s(i,j,k))$	Length of longest path from subtask $s(i,j,k)$ to exit node
$pred(s(i,j,k))$	Set of immediate predecessors of $s(i,j,k)$
$succ(s(i,j,k))$	Set of immediate successors of $s(i,j,k)$
$est(s(i,j,k))$	Earliest start time of subtask $s(i,j,k)$ provided it will execute on processor $p_z$
$ect(s(i,j,k))$	Earliest completion time of subtask $s(i,j,k)$ provided it will execute on $p_z$
$fproc_m(s(i,j,k))$	$m^{\text{th}}$ favorite processor for subtask $s(i,j,k)$
$fpred(s(i,j,k))$	Favorite predecessor of subtask $s(i,j,k)$
$bst(s(i,j,k))$	Best start time of subtask $s(i,j,k)$
$bct(s(i,j,k))$	Best completion time of subtask $s(i,j,k)$
CL	Set of real-time subtask clusters

different processors. Two conditions—namely,  $C_1$  and  $C_2$  are shown. A condition  $C_x \in C$  has two possible outcomes—true ( $C_x$ ) or false ( $\neg C_x$ ). Based on the outcomes of various conditions for a task set, a subset of the tasks is executed. We assume that the same conditions exist for all invocations of the real-time tasks. It has been assumed that in each invocation ( $k$ ) during  $pd_i$ , the period of a task ( $t_i$ ), a task needs to be executed for all subtasks ( $s(i,j,k)$ ). The deadline of a real-time task is given by  $d_i$ , and it can be  $\leq$  the task

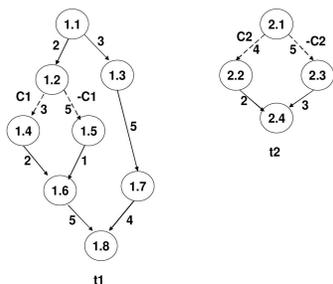


Fig. 1. Representing the data and control flow between two real-time tasks  $t_1$  and  $t_2$ .

period. The schedule is generated for a length equal to the *hyperperiod* (or *hp*), which is defined as the least common multiple (LCM) of all the task periods. Such a scheme may prove expensive if the periods for the real-time tasks are prime with respect to each other (as it will result in a very large hp). Some real-time literature states that “*in many realistic cases, this worst case may not be encountered*” [40] and “*in industry practice, task periods are typically assigned to be integer multiples of each other*” [41]. However, if the tasks in the application have coprime periods, then one can modify the periods of some selective tasks (increase or decrease) by a small amount [33]. Since the periods are modified by a small amount (1-5 percent), the feasibility of the real-time schedule is not affected drastically.

We assume that a set of processors are connected by a point-to-point network. A deadline-driven protocol called the earliest-due-date deadline (EDD-D) is used to schedule messages between the real-time subtasks [42]. When a subtask called the sender  $s$  wishes to send a message to a subtask called the destination  $d$ , EDD-D sets up a channel  $s \rightarrow d$ . The set of messages in the system is represented by  $M = \{m_1, m_2, m_3, \dots, m_n\}$ , and each message is characterized by its deadline, given by  $d(m_i)$ , given  $m_i \in M$ . Messages are scheduled to be transmitted on links between



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Input:  
 task set  $T = (t_1, t_2, \dots, t_n)$ , subtask set  $S(t_i)$  for all tasks in  $T$  such that  $s(i,j,k) \in S(t_i)$   
 processor set  $AP$   
 $pred(s(i,j,k))$ : set of predecessors/parents of subtask  $s(i,j,k)$   
 $succ(s(i,j,k))$ : set of successors/children of subtask  $s(i,j,k)$

Output:  
 real-time schedule

Begin:

1. obtain the  $pd_i$  values of all  $t_i \in T$  and calculate the  $hp$ .
2. **if** ( $hp$  is large)
3.     modify selected  $pd_i$  values by a small amount so that  $hp$  is reduced.
4.     calculate  $est, fproc_m, ect, fpred$  and  $p_e$  for all subtasks  $\in S(t_i)$  of all tasks  $\in T$ ;
5.     calculate  $bst$  and  $bct$  for all subtasks  $\in S(t_i)$  of all tasks  $\in T$ ;
6.     perform cluster generation;
7.     schedule the tasks on to the appropriate processors;
8.     perform task duplication or cluster compaction;
9.     perform message scheduling;

End

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Fig. 2. Steps for RT-DBA.

$fpred$  has already been assigned previously, then the remaining set of predecessors are examined; the predecessor that has the minimum execution time on the current processor is chosen. This set will also exclude the predecessors that cannot be executed because of control dependencies.

- e. Next, the best start time ( $bst$ ) and the best completion time ( $bct$ ) are calculated for each subtask in a bottom-up fashion. These quantities are used to find out if two subtasks are *critical*. Two subtasks  $s(i,y,z)$  and  $s(i,j,z)$  are critical with respect to each other if  $bst(s(i,y,z)) - bct(s(i,j,z)) < c(s(i,j,z), s(i,y,z))$ . The significance of subtask criticality implies that if two subtasks are not critical to each other, they can be assigned to different processors. For calculating the  $bct$  value of a subtask node, the  $bst$  values of all of its successors need to be calculated. If a subtask node has more than one successor, the minimum  $bst$  value is chosen as the  $bct$  value. In addition, if the subtask node is the  $fpred$  of its successor, the communication cost between them is zero. Otherwise, the communication cost is subtracted from the  $bst$  value (see (7) in Table 2).
  - f. Next, the execution cost of a subtask on a fictitious processor is calculated (see (9) in Table 2). This quantity is the average execution cost of a subtask on all available processors. This quantity is used for the execution cost of a subtask in case  $AP < CNP$  and the cluster compaction procedure needs to be carried out (please see stage 4).
3. **Stage 3 (perform cluster generation).** In stage 3, the cluster generation procedure is carried out (see Fig. 3). This stage involves the generation of a *queue* that contains all the subtasks in the graph arranged in increasing order of their  $p_e$  values. When calculating the  $p_e$  values, only the computation costs are taken into consideration (the communication costs are ignored). The clusters are formed by following the  $fpred$  chain from the bottom to the top

of the DAG. The generation of clusters is accomplished by performing a depth-first search, starting from the exit node. The search progresses by tracing the path from the initial subtask (the first *queue* element) to the entry subtask by following the  $fpred$  value. The first cluster completes when the entry subtask is reached for the first time. The next cluster then starts from the next unassigned subtask in the *queue*. If the  $fpred$  of a subtask has already been assigned and if it is not critical, then an unassigned subtask (among the current subtask's immediate predecessors), which has the lowest execution time on the current processor, is selected. The cluster generation process continues recursively, till all the subtasks have been assigned. Each cluster is then allocated to the first available  $fproc_m$  of the header subtask of that cluster.

The process outlined above assumes that the deadline of a task is equal to its period. If the deadlines of the real-time tasks are smaller than their periods, then the quantity  $(d_i - pd_i)$  is calculated for all the tasks. Then, the clusters are formed for the tasks based on a decreasing value of  $(d_i - pd_i)$ . The main idea behind this approach is that the tasks with higher  $(d_i - pd_i)$  values need to finish execution earlier. Hence, they deserve to be assigned to faster processors.

It may also be possible that the deadline of a task is greater than its period. Hence, at any time instant, there may be multiple invocations of a real-time task. The proposed scheme can be easily modified to handle such a case. The first invocation of the real-time task is assigned as before. The subsequent invocations will be assigned to the extra processors. After this assignment, the duplication phase can be carried out to enhance the SR.

4. **Stage 4 (compare AP and CNP and perform duplication or compaction).**
  - i. If  $AP > CNP$ , subtask duplication is carried out to increase the SR. A check is done to see if each subtask in a cluster is preceded by its  $fpred$

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Input: DAG, set of predecessor tasks, set of successor tasks, and queue.  
Output: Set of task clusters.  
Begin:

1.  $s(i,y,k)$  = first element of queue;
2. current processor =  $fproc_l$  of  $s(i,y,k)$ ;
3. assign  $s(i,y,k)$  to current processor;
4. **while** (NOT all tasks assigned) {
5.    $s(i,z,k) = fpred(s(i,y,k))$ ;
6.   **if**( $s(i,y,k)$  has more than one immediate predecessors) {
7.     **if**(( $bst(s(i,y,k)) - bct(s(i,z,k)) < c(s(i,z,k), s(i,y,k))$ )) /\* if they are critical to each other\*/
8.     assign  $s(i,y,k)$  and  $s(i,z,k)$  to the same processor;
9.     **else if**( $s(i,z,k)$  has already been assigned OR cannot be executed because of control dependency)
10.      $s(i,z,k)$  = predecessor of  $s(i,y,k)$  that has the lowest run time on this processor;
11.   }
12.   assign  $s(i,z,k)$  to current processor;
13.    $s(i,y,k) = s(i,z,k)$ ;
14.   **if**( $s(i,y,k)$  is the entry node) {
15.     assign  $s(i,y,k)$  to current processor;
16.      $s(i,y,k)$  = next unassigned element in *queue*;
17.     current processor = next available  $fproc_m$  of  $s(i,y,k)$ ;
18.   }
19. }  
End

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Fig. 3. Cluster generation phase of RT-DBA.

subtask. If this is not the case, then the tail of that particular cluster is replaced with its  $fpred$  chain, till the entry node. The cutoff tail is assigned to the next available  $fproc_m$  of the last node. Duplication is carried out till all subtasks are preceded by their favorite predecessors or till there are no more remaining processors.

- ii If  $AP < CNP$ , then the excess clusters are accommodated in fictitious processors. The procedure for cluster compaction is given in Fig. 4. The clusters are then arranged in decreasing order of  $c_c(cl_x, p_z)$ , which is defined as the sum of execution times of all the subtasks assigned to  $p_z \in P$ . The value of  $c_c(cl_x, p_z)$  for each cluster  $cl_x \in CL$  is given by  $c_c(cl_x, p_z) = \sum \mu(s(i, j, k), p_z)$ , where  $s(i, j, k) \in t_i$ , and  $t_i \in T$ . The *least populated* cluster of a system processor is merged with the *most populated* cluster of a fictitious processor. This process is carried out

till the number of clusters becomes equal to the number of processors. The  $p_e$  value of a subtask is used to find out the place where the subtask is to be inserted. The subtasks are inserted so that the increasing  $p_e$  value is maintained as we go from a larger numbered subtask to a smaller numbered subtask.

Our algorithm can also handle the case when AP is variable. Based on the initial AP, RT-DBA performs one of the following actions: 1) compaction if  $AP < CNP$  (mode<sub>1</sub>), 2) mapping clusters on to the processors if  $AP = CNP$  (mode<sub>2</sub>), and 3) duplication if  $AP > CNP$  (mode<sub>3</sub>). The initial value of AP can either increase or decrease. If AP increases, the action taken by RT-DBA will depend on the mode that it is in. If it is in mode<sub>1</sub>, then the compaction procedure is called with this new AP value. If the value of AP increases further, it could go to mode<sub>2</sub> or mode<sub>3</sub>. If it is in mode<sub>2</sub>, then the duplication procedure can be called. If it is in mode<sub>3</sub>, then further duplication can be carried out. If AP

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1. **while**( $AP < CNP$ ) {
2.   calculate  $c_c$  for each cluster  $cl_x \in CL$ ;
3.   sort available processors in ascending order of  $c_c$ ;
4.   temp =  $CNP/2$ ;
5.   **for**( $j=0; j < temp; j++$ )
6.     merge task lists of processors  $j$  and  $(temp*2-j-1)$ ; /\*use  $p_e$  values to determine relative positions \*/
7.     decrement  $CNP$ ;
8.   } //end while

End

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Fig. 4. Pseudocode for cluster compaction.

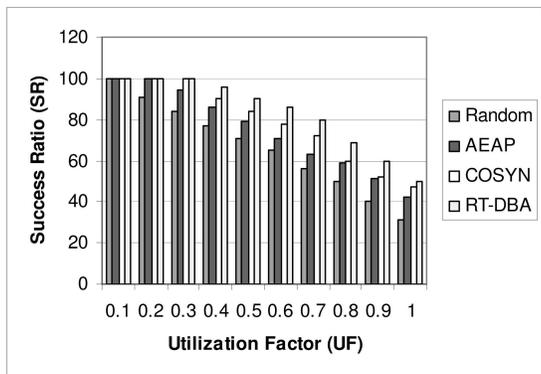


Fig. 5. SR versus UF for CCR = 0.1.

decreases, the action taken by RT-DBA depends on the mode that it is in. If it is in  $mode_1$ , then more compaction can be carried. If it is in  $mode_2$ , compaction can be carried out by calling the compaction procedure. If it is in  $mode_3$ , the amount of duplication can be reduced by calling the duplication procedure with the new AP value. If the AP value decreases further, then it could go to  $mode_2$  or  $mode_1$ .

## 6 SIMULATION RESULTS

In this section, we compare the performance of RT-DBA with relevant algorithms in the literature, namely, AEAP [21] and COSYN [33] (the performance of MOGAC is similar to COSYN [34]). Both AEAP and COSYN are heuristic-based algorithms that schedule precedence-related periodic tasks on heterogeneous multiprocessors. We have used both randomly generated task graphs and graphs that represent real-life applications [43] such as sparse matrix solver and SPEC fppp. Randomly generated graphs have been used extensively in the past by researchers [12], [21], [22], [23]. Input DAGs have also been generated using the popular Task Graphs for Free (TGFF) suite [44], as well as the benchmark Standard Task Graph Set [43]. The processor architectures have been generated in a random fashion. We have coded a software scheduler module that accepts a set of precedence-related periodic tasks and generates a real-time schedule. Four performance metrics have been used. The first metric is the SR and is defined as the ratio of the number of task sets found to be schedulable (all tasks meet deadlines) by the algorithm to the total number of task sets considered. The second metric is called the Compaction Factor (CF), and it is defined as  $AP/CNP$ . The third metric is called the Heterogeneity Factor (HF), and it is defined as the standard deviation of the subtask execution costs. To calculate the standard deviation of a set of execution costs, we

1. calculate the mean value of all execution costs,
2. calculate the deviation for each execution cost  $(cost - mean)$ ,
3. calculate the squares of all the deviations obtained in step 2,
4. calculate the mean of all squared deviations, and
5. calculate the square root of the mean obtained from step 4.

The fourth metric is called the total Communication-to-Computation Ratio (CCR), and it is defined as *average edge*

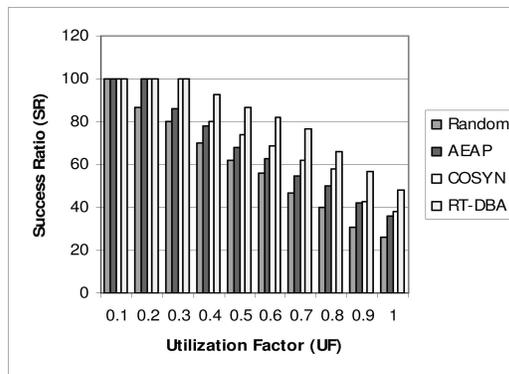


Fig. 6. SR versus UF for CCR = 1.0.

*cost/average node cost*. The execution cost for a node is the cost on its chosen processor. A low value of CCR would imply low levels of communication (versus computation) and vice versa.

The task and system parameters are generated in a way similar to [12]. The CCR values between subtasks in a periodic task are given by a uniform distribution and have been selected from the range (0, 10). The execution cost of each subtask is given by a uniform distribution and has been selected from the range (1, 100). The period of the first task in the task set has been set as follows:  $pd_i = C_{av} * task\_size$  [12]. Here,  $C_{av}$  is the average execution time of all the subtasks in the periodic task, and  $task\_size$  is the number of subtasks. The utilization factor is given by a uniform distribution, and its value is selected from the range (0, 1). The results shown here are for task sets with five periodic tasks  $\{T = t_1, t_2, t_3, t_4, t_5\}$ . Tasks  $t_1$  to  $t_5$  have  $\{2, 4, 6, 8, 10\}$  subtasks, respectively. The total number of edges has been selected from the range (20, 60). The number of predecessors and successors has been varied from 1 to 10. The periods of the tasks are  $pd_1, pd_2, pd_3, pd_4$ , and  $pd_5$ . If the period of the first task is  $pd$ , then the periods of the subsequent tasks will be  $2pd, 3pd, 4pd$ , and  $5pd$ . The length of the schedule generated is the LCM of the periods, i.e., it is  $60 * pd$ . Enough simulation runs are performed to provide 95 percent confidence interval. Each point in the following curves is the result of 1,000 runs of the algorithms. For each point in the following graphs, we generated 1,000 task sets. All the simulations were run on a Pentium 4 2.4-GHz PC running Red Hat Linux.

### 6.1 Effect of Communication to Computation Ratio

Figs. 5 and 6 indicate the effect of CCR on the proposed schemes. The number of processors ranged from 5 to 25. The utilization factor of the heterogeneous system is given by the equation  $UF_{hetero} = \sum \mu(s(i, j, k)) / (AP * hp)$  for all  $t_i \in T$ . The subtask execution times have been generated randomly from the range  $\{2, 200\}$ . The experiments have been run for CCR values of 0.1, 1.0, and 10.0, representing varying levels of communication. All of the four algorithms have been assigned an equal number of processors. In Figs. 5 and 6, we observe that as the system utilization increases, the SR decreases. This can be explained as follows: as the utilization increases, so do the subtask execution times (for a fixed period). Hence, fewer subtasks (and, consequently, task sets) succeed in meeting their deadlines, and the SR value reduces. We also observe that RT-DBA offers a higher

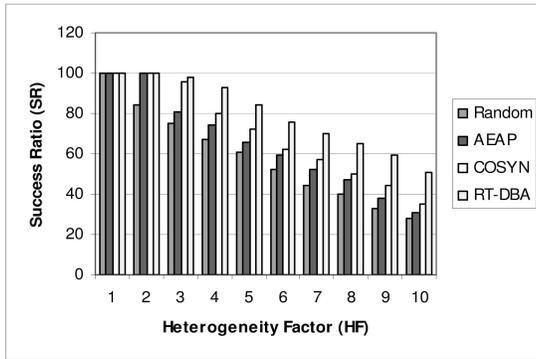


Fig. 7. Effect of heterogeneity.

SR as compared to the other schemes, as selective subtask duplication results in some subtasks finding earlier start (and, hence, finish) times. This results in an increase in the number of task sets that can meet their deadlines. Hence, the SR value is higher as compared to AEAP and COSYN that do not employ duplication. It can also be seen that the improvements in the SR are more pronounced for the case where the CCR value is 1.0 than that when  $CCR = 0.1$ . This is due to the fact that the aim of subtask duplication is to reduce the interprocessor communication between subtasks assigned to different processors. Hence, the gains (in terms of SR values) are greater when there is more communication in the system.

## 6.2 Effect of Heterogeneity

We have captured the effect of subtask heterogeneity (in terms of the execution time) on the performance of the real-time heterogeneous system in this simulation. The number of processors varied from 5 to 27. The  $min_e$  value has been kept fixed at 10, and the  $max_e$  value has been incremented by a constant factor of 4 time units. The results of this study are shown in Fig. 7. We observe that as the HF value of the system increases, its SR experiences a decrement. This is because with an increase in the HF value, there is more variation in the subtask execution times.

Also, a task set with larger variation in execution times will have a larger average subtask execution time. Hence, it will be assigned a larger period as compared to the same task set with lesser variation in execution times. But there is still a larger disparity between the average subtask execution time and the actual subtask execution times for the task set with a higher HF value. Hence, as the HF value is increased, the actual completion time of the subtasks increases. Therefore, a lesser number of subtasks succeed in meeting their deadlines. This leads to a decrease in the SR of the application. A CCR value of 1.0 has been assumed in this study.

## 6.3 Effect of Cluster Compaction

This study demonstrates the scalability aspect of RT-DBA. The results of this study are illustrated in Fig. 8. A low value of CF indicates that  $AP \ll CNP$ . On the other hand, a high value of CF indicates that  $AP \gg CNP$ . The number of processors ranged from 3 to 30. We observe that as the CF increases, the SR value also increases. This is due to the fact that as the CF value increases, processors are added to the heterogeneous system, which enables the system to exploit the parallelism offered. Hence, a larger number of subtasks (and, hence, task sets) meet their deadlines. If  $CF < 1$ ,

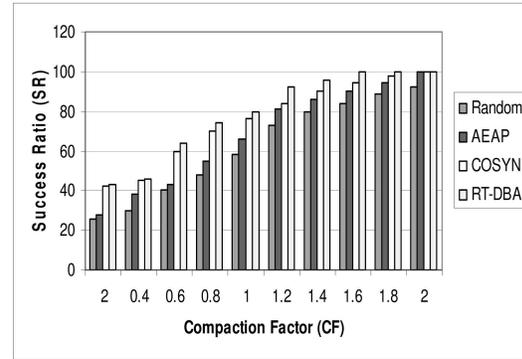


Fig. 8. Effect of compaction.

compaction is carried out; else, duplication is performed. Hence, more task sets are able to meet their deadlines, and the SR increases versus the compared schemes. Although AEAP and COSYN do not support compaction, we included this feature for the sake of fairness in comparison. A CCR value of 1.0 has been assumed in this experiment.

## 6.4 Effect of Message Load on Success Ratio

Fig. 9 shows the effect of the message load on the SR of the application. A CCR of 1.0 and  $N = 30$  have been assumed for this simulation. The number of messages has been varied from 6 to 60 in increments of six. The message load = number of messages/number of tasks in the application. It is observed that as the message load increases, the SR value decreases. That is because with increased message load, there are more messages in the system that need to be scheduled (for the same number of tasks and processors). Hence, there is a greater load on the processors, and consequently, more tasks (and, hence, task sets) miss their deadlines.

## 6.5 Results for Real-Life Application Graphs

We have tested and compared various schemes for the task graphs that represent real-life applications such as sparse matrix solver and SPEC fppp [43]. The results of these simulations are given in Figs. 10 and 11. We observe that our proposed scheme offers a higher utilization value (the utilization value at which task sets begin to miss deadlines) as compared to COSYN, AEAP, and the Random scheme. If there is no communication in the application (i.e.,  $CCR = 0$ ), the performance of COSYN and RT-DBA is similar. However, RT-DBA outperforms COSYN (and the other schemes) when there is significant communication in the system (i.e.,  $CCR = 1.0$ ). This is because, the aim of duplication is to

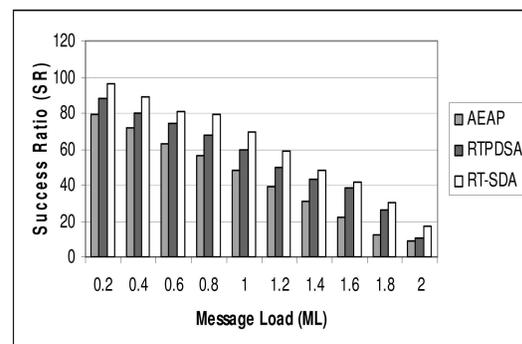
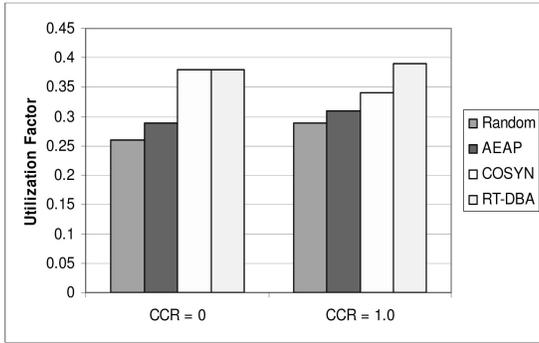


Fig. 9. Effect of message load on SR.

Fig. 10.  $U_b$  for the sparse matrix task graph.

reduce/minimize the IPC time between communicating subtasks assigned to different processors. This ensures that selected subtasks start (and, hence, finish) earlier. Hence, more subtasks (and task sets) are able to meet their deadline requirements. Therefore, it follows logically that the gain in SR will be more when there is more communication in the system. The sparse matrix solver graph has 96 tasks and 67 edges. The number of predecessors has been randomly chosen from  $\{0, 6\}$ . The execution costs have been randomly chosen from  $\{0, 35\}$ . The SPEC fppp graph has 334 tasks and 1,145 edges. The number of predecessors has been randomly chosen from  $\{0, 80\}$ . The execution costs have been randomly chosen from  $\{0, 404\}$ .

In addition, we have also tested all the algorithms using the TGFF suite [44]. The results are shown in Fig. 12. A CCR value of 1.0 has been used. We observe that as the utilization factor increases, the SR decreases. That is because the execution costs of the subtasks increases, and there is more tightness in the schedule. We also see that RT-DBA offers a better SR as compared to other existing schemes, owing to subtask duplication. For this experiment, we considered a task set  $T = \{t_1, t_2, t_3\}$ . Tasks  $t_1$ ,  $t_2$ , and  $t_3$  have 4, 8, and 12 subtasks, respectively. The task periods are  $pd_1$ ,  $pd_2$ , and  $pd_3$ . If the period of the first task is  $pd$ , then the periods of the second and the third tasks will be  $2pd$  and  $3pd$ , respectively. Hence,  $hp = 6pd$ . The execution times have been chosen uniformly from the range  $\{2, 200\}$ . For each subtask, the number of predecessor tasks and successor tasks were chosen uniformly from the range  $\{1, 10\}$ .

### 6.6 Effect of Data and Control Dependency

Fig. 13 shows the effect of control dependency on the system performance. This simulation has been conducted for a single periodic task consisting of 30 subtasks. The

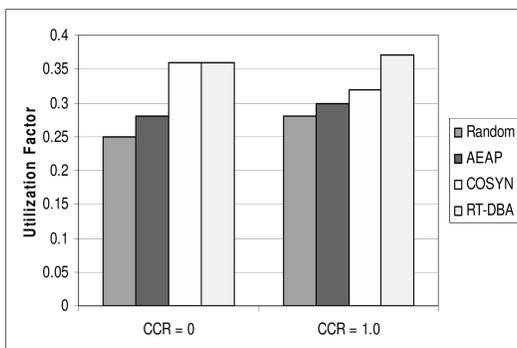
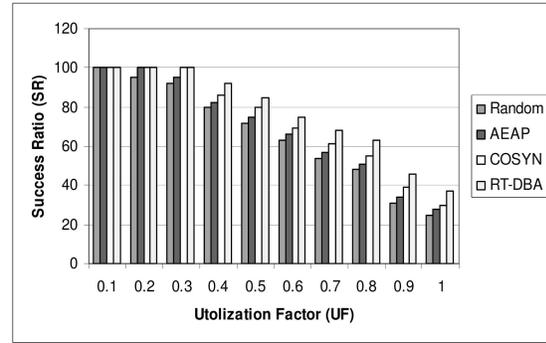
Fig. 11.  $U_b$  for SPEC fppp.

Fig. 12. SR versus UF for TGFF (CCR = 1.0).

number of conditions have been chosen as four for this task set. This resulted in the formation of eight subgraphs. In the figure, we observe that the addition of control dependency results in a higher SR for the real-time application. This can be explained as follows: with the addition of conditions between tasks, only a subset of them executes (as compared to the scenario in which there is no condition dependency between tasks and, hence, all tasks are executed). This leads to a reduction of load on the NOWs as multiprocessors, and this translates to a higher SR value.

### 6.7 The Case When Deadline Is Less than Period

In this section, we discuss the case when the deadline is smaller than the period. The results of our simulation are shown in Fig. 14. The same model that is described at the beginning of Section 5 has been used in this study. Looking at the figure, it can be seen that for the largest  $(d - pd)$ , the first scheme offers a better SR value than the plain RT-SDA. This can be explained as follows: The tasks with larger  $(d - pd)$  values have less time to complete execution than the tasks with smaller  $(d - pd)$  values and therefore should be executed first. This ensures that such tasks get their "favorite processors," and hence, the SR of the application could improve as compared to the plain RT-SDA.

## 7 CONCLUSION

The major contribution of this research is in proposing a selected duplication-based algorithm (RT-DBA) for enhancing the schedulability of precedence-related periodic tasks on real-time heterogeneous NOWs. When sufficient processors are available, subtask duplication cuts down

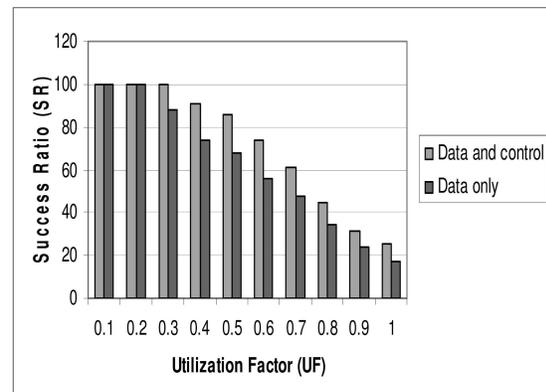
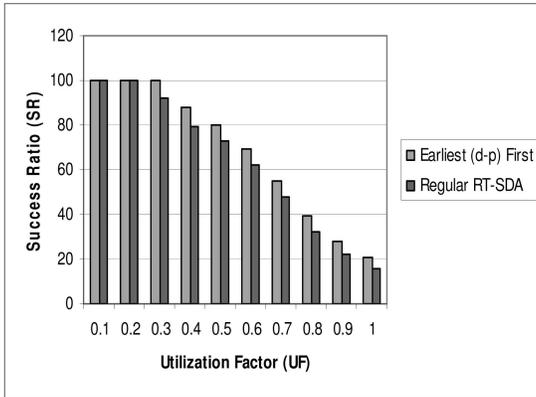


Fig. 13. Effect of control dependency.

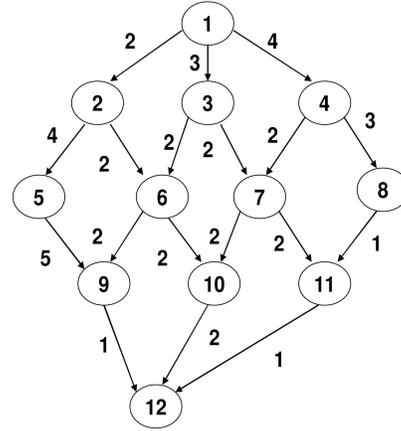
Fig. 14. Case when  $d_i < pd_i$ .

the interprocessor communication between communicating subtasks assigned to different processors. This enables some subtasks to find earlier start times. Hence, more subtasks (and task sets) are able to meet their deadline requirements. This phenomenon enables RT-DBA to offer a higher SR as compared to the other schemes. We strongly believe that the duplication can be used as a valuable tool to improve the schedulability of real-time applications, especially when communication is a dominant factor in the system. In addition, our scheduling scheme supports scalability for the heterogeneous NOWs in that the application is schedulable even if the required number of processors is not available, although this would result in a reduction in the schedulability of the application. Therefore, we conclude that a trade-off exists between scalability and schedulability—increased scalability results in a decrease in schedulability and vice versa.

## APPENDIX A

Now, we will illustrate the working of RT-DBA with the help of the sample DAG ( $t_1$ ) in Fig. 15. The period of  $t_1 = 50$ . The execution times of all the subtasks of  $t_1$  are shown in Table 3. The algorithm works as follows:

1. First, we obtain the  $est$  for all subtasks. Let us consider subtask 10, which has two immediate predecessors: 6 and 7. The  $fproc_1$  of 6, 7, and 10 is  $p_1$ . According to its equation,  $est(10) = \max. \{ect(7) \parallel ect(6) + c(6, 10)\} = \max. \{11 \parallel (10 + 2)\} = \max. \{11 \parallel 12\} = 12$ . The  $est$  values of all the subtasks in the DAG are shown in Table 4.
2. Next, we obtain the  $fproc_m$  for all subtasks. Let us consider subtask 1.10. The sum of its  $est$  plus its execution time on all the processors ( $p_1$  to  $p_8$ ) is calculated as  $\{(12 + 1), (12 + 5), (12 + 6), (12 + 7), (12 + 4), (12 + 8), (12 + 5), (12 + 6)\} = \{13, 17, 18, 19, 16, 20, 17, 18\}$ . On arranging these values in increasing order, we get  $\{13, 16, 17, 17, 18, 18, 19, \text{and } 20\}$ . These values correspond to processors  $p_1, p_5, p_2, p_7, p_3, p_8, p_4, \text{and } p_6$ , respectively, and they are the favorite processors for subtask 10 (in increasing order).
3. Now, we obtain  $ect$  for all nodes in the application. For a subtask 10,  $ect(10) = est(10) + \mu(10, fproc_1(10)) = 12 + 1 = 13$ . Hence, the  $ect$  value for 1.10 is 13.  $ect$  values for all subtasks are shown in Table 4.

Fig. 15. Sample application ( $t_1$ ).

4. Now, we obtain the  $fpred$  for each node in the application. Let us consider subtask 12.  $fpred(12) = \{\max. (ect(9) + c(9, 12)) \parallel (ect(10) + c(10, 12)) \parallel (ect(11) + c(11, 12))\} = \max. \{(19 + 1), (13 + 2), (15 + 1)\} = \max. \{20, 15, 16\} = 20$ . Hence,  $fpred(12) = 9$ . The  $fpred$  values of all subtasks are shown in Table 4.
5. Next, we obtain  $bst$  and  $bct$  for all subtasks. Let us consider subtask 3. It has two successor subtasks—6 and 7. Now,  $bst(6) = 11$ , and  $bst(7) = 12$ . According to the equation in Table 4,  $bct(3) = \min. \{bst(6) - c(3, 6), bst(7)\}$ . Hence,  $bct(3) = \min. \{(11 - 2), 12\} = \min. \{9, 12\} = 9$ . Therefore,  $bst(3) = bct(3) - \mu(3, fproc_1(3)) = bct(3) - \mu(3, p_1) = 9 - 2 = 7$ .
6. Now, we arrange all nodes in the *queue* and start cluster formation. Initially, the *queue* formation takes place. For task  $t_1$ , the *queue* is  $\{12, 11, 10, 9, 6, 7, 8, 5, 3, 4, 2, \text{and } 1\}$ . The generation of the first cluster starts from 12. It is preceded by  $fpred(12) = 9$ . Subtask 9 is preceded by its  $fpred$ , which is 5. Continuing in a similar fashion, the next subtasks will be 2 and 1. Hence, the first cluster is 12-9-5-2-1 and is assigned to its  $fproc_1 = p_5$ . Similarly, the second, third, and fourth clusters are 11-7-4-1, 10-6-3-1, and 8-4-1, and they are assigned to  $p_1, p_2, \text{and } p_7$ , respectively. The complete schedule is shown in Fig. 16. It has a schedule length of 34 time units. Since three free processors are available, duplication can be carried out. Subtask 7 in the second cluster is not preceded by its  $fpred$ . Hence, the subcluster 4-1 is cut off and is replaced with the  $fpred$  chain 3-1. The second cluster is modified to 11-7-3-1, and the cutoff subcluster 4-1 is allocated to  $p_6$ , which is the best available processor remaining. The schedule length now has reduced to 30 (see Fig. 17). Hence, duplication has resulted in a gain of 4 time units.

### A.1 RT-DBA Time and Space Complexity Analysis

Let  $m_j$  be the total number of subtasks in DAG  $d_j \in D$ . Let  $p$  be the number of heterogeneous processors and  $e$  be the total number of communication edges in the DAG. The first step involves calculating the  $est, ect, fpred, fproc_m, p-e,$  and *queue*. While calculating the  $est$ , for each node, all the predecessors need to be examined, and all edges need to be traversed. Hence, the number of predecessors =  $m_i$ . Therefore, for  $m_i$  nodes, the complexity for calculating  $est$  is

TABLE 3  
Task Execution Times

Tasks	$\mu(p_1)$	$\mu(p_2)$	$\mu(p_3)$	$\mu(p_4)$	$\mu(p_5)$	$\mu(p_6)$	$\mu(p_7)$	$\mu(p_8)$
1	4	6	7	8	4	6	5	6
2	2	5	9	5	5	6	5	6
3	2	9	4	3	2	6	5	6
4	1	6	9	8	5	4	5	6
5	4	9	5	4	3	3	5	6
6	2	3	4	6	2	5	5	6
7	4	5	6	7	4	6	5	6
8	6	7	8	9	5	7	5	6
9	8	5	6	4	6	7	5	6
10	1	5	6	7	4	8	5	6
11	3	7	6	5	5	5	5	6
12	5	7	8	9	4	6	5	6

TABLE 4  
Mathematical Quantities for the Tasks Shown in Fig. 15

Subtask	$fproc_1$	$fproc_2$	$fproc_3$	$fproc_4$	$fproc_5$	$fproc_6$	$fproc_7$	$fproc_8$	$est$	$ect$	$fpred$	$bst$	$bct$	$p\_e$
1	P1	P5	P7	P2	P6	P8	P3	P4	0	4	-	3	7	43
2	P1	P2	P4	P5	P7	P6	P8	P3	4	6	1	9	11	35
3	P1	P4	P5	P3	P7	P6	P8	P2	4	6	1	7	9	33
4	P1	P6	P5	P7	P2	P8	P4	P2	4	5	1	8	9	34
5	P1	P5	P6	P4	P3	P7	P8	P2	6	9	2	12	16	26
6	P2	P1	P5	P3	P6	P7	P4	P8	7	9	2	11	14	23
7	P1	P5	P2	P7	P3	P6	P8	P4	7	11	4	12	16	24
8	P1	P5	P7	P8	P2	P6	P3	P4	5	11	4	9	15	25
9	P4	P1	P2	P7	P3	P5	P8	P6	15	19	5	16	20	17
10	P1	P5	P2	P7	P3	P8	P4	P6	12	13	7	17	18	17
11	P1	P4	P5	P6	P7	P3	P8	P2	12	15	7	16	19	16
12	P5	P1	P7	P6	P8	P2	P3	P4	20	24	9	20	24	9

$O(m_i^2 + e)$ . While calculating the  $fproc$ , for each node, all  $p$  processors need to be examined. Therefore, for  $m_i$  nodes, the complexity is  $O(m_i * p)$ . While calculating the  $ect$ , all nodes need to be examined. Therefore, the complexity for this stage is  $O(m_i)$ . While calculating the value of  $p\_e$ , for each node,  $m_i$  nodes need to be examined, and so, the complexity is  $O(m_i^2)$ . The formation of the *queue* involves sorting nodes based on  $p\_e$  values (using Quick Sort). Therefore, the complexity is  $O(m_i * \log_2 * m_i)$ . While calculating the  $fpred$ , for each node, roughly  $m_i$  predecessors

need to be examined, so the complexity is  $O(m_i^2)$ . On adding all the above, the complexity becomes  $O(m_i^2 + e + m_i + m_i * p + m_i^2 + e + m_i * \log_2 m_i)$ . For most practical DAGs and dense graphs,  $m_i \gg p$ , and  $e$  is of the order of  $m_i^2$ . Hence, the complexity =  $O(m_i^2)$ . The second step deals with the calculation of the quantities  $bst$  and  $bct$ . This involves a bottom-up traversal of the DAG. The complexity of this step is  $O(e)$ . The next step involves the formation of the clusters. This is similar to a depth-first search of the DAG, and the complexity of this step is  $O(m_i + e)$ . For a

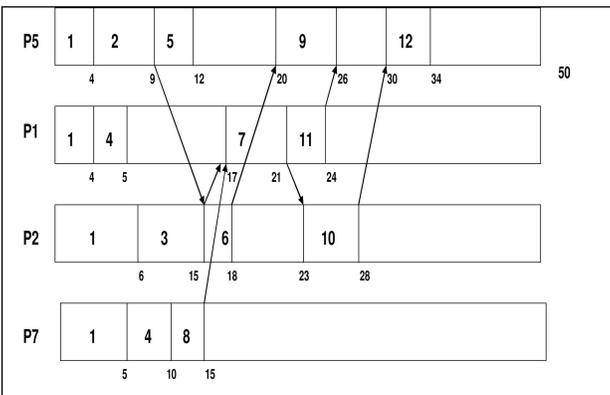


Fig. 16. Initial schedule.

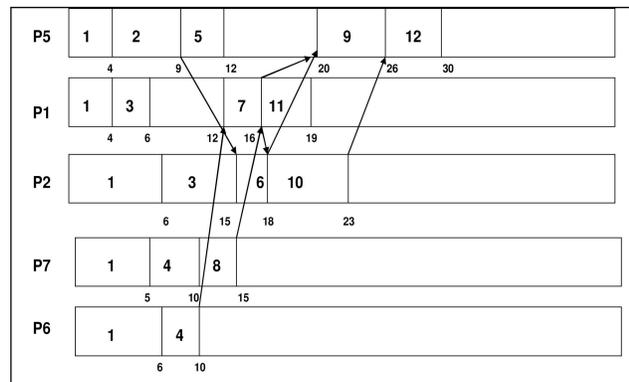


Fig. 17. Schedule after first duplication.

connected graph, this is equivalent to  $O(e)$ . The complexity of the duplication stage is given by  $O(t * m_i)$ , where  $t$  is the number of subtasks not preceded by their  $fpred$  subtasks. In the worst case, all the subtasks are duplicated. In other words,  $t = m_i$ . Hence, the complexity becomes  $O(m_i^2)$ . The next stage is the cluster compaction stage, and this step involves sorting of the processors based on the empty slots. Therefore, each pass through the compaction code is of the order of  $O(m_i + p * \log_2 p)$ . In the worst case,  $p = m_i$ . The number of required passes depends on the difference between AP and CNP and is given by the expression  $Q = \log_2(m_i)$ . Therefore, the complexity of compaction is  $O(m_i^2)$ . The calculation of the utilization factor values involves the traversal of all the invocations of all tasks. Therefore, the complexity of this step is  $O(m_i)$ . Hence, the overall complexity is  $O(m_i^2 + e + e + m_i^2 + m_i^2 + m_i)$ . For dense graphs,  $e = m_i^2$ . Hence, the time complexity for the specific DAG with  $m_i$  is  $O(m_i^2)$ . Now, DAG<sub>1</sub> will run  $hp/pd_1$  times. Similarly, DAG<sub>2</sub> will run  $hp/pd_2$  times, and so on. Hence, the time complexity for all  $n$  DAGs is given by the following equation:

$$\begin{aligned} & m_1^2 * (hp/pd_1) + m_2^2 * (hp/pd_2) + \dots + m_n^2 * (hp/pd_n) \\ & = hp(m_1^2/pd_1 + m_2^2/pd_2 + \dots + m_n^2/pd_n). \end{aligned}$$

Each real-time subtask can be represented as a floating point and can therefore be stored in one memory cell. The total number of real-time tasks is  $hp/pd_1 + hp/pd_2 + hp/pd_3 + \dots + hp/pd_n = hp * (1/pd_1 + 1/pd_2 + 1/pd_3 + \dots + 1/pd_n)$ . For the sake of brevity, let us assume that this is equal to  $x$ . Since each subtask needs one memory cell,  $x$  subtasks need  $x$  memory cells. For all DAGs, the total number of edges can be given by  $E$ . For very dense graphs,  $E = x^2$ . Hence, all edges need  $x^2$  memory locations. The execution cost array for a subtask depends on the number of processors. Each subtask needs AP elements. Therefore, all  $x$  subtasks will need  $x * AP$  elements. In the worst case, there will be one subtask and one processor, so  $AP = x$ . Hence, all  $x$  subtasks will need  $x^2$  memory locations. The number of subtasks that need to be duplicated  $= x^2$ . On adding all the above, we get  $x + x^2 + x^2 + x^2$ . Hence, the space complexity of RT-DBA is of the order of  $x^2$ .

**Schedulability test condition S-1.** A task  $t_i \in T$  is schedulable with RT-DBA if the following condition is satisfied:  $ect(s(i, j, k)) \leq hp$  such that  $|succ(s(i, j, k))| = 0$ .

The exit node will have zero successors. Because of the precedence constraints, the exit node will always be the last node to execute. A task set  $T = \{t_1, t_2, t_3, \dots, t_{ap}\}$  will be schedulable with RT-DBA if the above test condition holds for all  $t_i \in T$ .

**Schedulability test condition S-2.** If  $UF_{hetero}$  corresponds to the task load of the heterogeneous system,  $\Omega_{hetero}$  corresponds to the communication load, and  $\Phi_{hetero}$  corresponds to the idle load, then a task set is schedulable with RT-DBA if  $UF_{hetero} \leq AP - (\Omega_{hetero} + \Phi_{hetero})$ .

$CL_{hetero}$  refers to the communication load of the real-time application. The communication cost is the delay associated with data transmission and is given by  $c(s(i, j, k), s(i, l, k))$ , where  $s(i, j, k)$  and  $s(i, l, k)$  are the two communicating subtasks. Note that based on our model, subtasks that are part of the same task can communicate with each other (hence, the  $i$  term in both subtasks is the same). Also, data transmission is done

between the same invocations of two communicating subtasks (hence, the  $k$  term in both subtasks is the same). Let us define  $\Delta$  to be the set of all precedence-related subtask pairs that are not assigned to the same processor:

$$\begin{aligned} \Delta & = \{s(i, j, k), s(i, l, k)\} \text{ for all } t_i \in T \\ & \text{where } i, j, k \text{ and } l \text{ are all integers.} \end{aligned} \quad (1)$$

The sum total of communication edge costs for all such pairs would give us the total communication cost. Let us say that this cost  $= \Omega_{hetero}$ . Hence, we get

$$\Omega_{hetero} = \sum c(s(i, j, k), s(i, l, k)) \text{ for all } t_i \in T \text{ and all } ST(t_i). \quad (2)$$

Since RT-DBA is nonpreemptive, there is a certain amount of time that a processor can be idle. For a specific cluster  $cl_\tau \in CL$ , this time can be given as  $hp - (lact(head_{cl_\tau}))$ , where  $head_{cl_\tau}$  corresponds to the head subtask of cluster  $cl_\tau$ . For all clusters, this time can be represented by (if the total number of clusters in an application is  $\alpha$ )

$$\begin{aligned} \Phi_{hetero} & = (hp - lact(head_{cl1})) + (hp - lact(head_{cl2})) \\ & + \dots + (hp - lact(head_{cl\alpha})). \end{aligned} \quad (3)$$

A real-time application will be schedulable if the sum of the task load, communication delay, and idle time is less than the total processing capacity of the workstations. The total processing capacity  $= AP$ . Hence, we get

$$\begin{aligned} UF_{hetero} + \Omega_{hetero} + \Phi_{hetero} & \leq AP \\ \Rightarrow UF_{hetero} & \leq AP - (\Omega_{hetero} + \Phi_{hetero}). \end{aligned} \quad (4)$$

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