

Online Algorithms for a Generalized Parallel Machine Scheduling Problem

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Abstract: We consider different *online* algorithms for a generalized scheduling problem for *parallel machines*, described in details in the first section. This problem is the generalization of the classical parallel machine scheduling problem, when the make-span is minimized; in that case each job contains only one task. On the other hand, the problem in consideration is still a special version of the workflow scheduling problem.

We present several heuristic algorithms and compare them by computer tests.

Keywords: scheduling, parallel machines, online algorithm.

1 The generalized Parallel Machine Scheduling problem

We are given a list of (types of) **tasks** T_1, \dots, T_t , and another list of **machines** M_1, \dots, M_m , and the table $I[i, j]$ which shows the required **time** what machine M_i needs to solve the task T_j ($1 \leq i \leq m$, $1 \leq j \leq t$). $I[i, j] \neq 0$ is assumed, but $I[i, j] < 0$ indicates that M_i is unable to solve T_j .

We receive online the list of **jobs** J_1, \dots, J_L , where any job J_ℓ ($1 \leq \ell \leq L$) contains several tasks from the list T_1, \dots, T_t , in blocks (see /0a/ below). The number of the jobs, L , and the number of *tasks* from which J_ℓ are build up, are unknown in advance. We have to schedule all the tasks of the jobs J_ℓ ($1 \leq \ell \leq L$) to the machines, fulfilling the following requirements.

/0a/ J_ℓ contains of f_ℓ many **blocks**

$$J_\ell = (B_{\ell,1}, \dots, B_{\ell,f_\ell}) \quad (1)$$

/0b/ where any block $B_{\ell,\phi}$ ($\phi \leq f_\ell$) contains some tasks

$$B_{\ell,\phi} = (T_{\ell,\phi,1}, \dots, T_{\ell,\phi,K_\phi}) \quad (2)$$

K_ϕ is called the **length** of $B_{\ell,\phi}$, assuming $1 \leq K_\phi$. The case $f_\ell=1$ is allowed, too. The **size** of the job J_ℓ is clearly

$$\mathbf{size}(J_\ell) = \sum_{\phi=1}^{f_\ell} K_\phi \quad (3)$$

/0c/ Since $T_{\ell,\phi,k}$ are members of the list T_1, \dots, T_t we can define the function φ as

$$\varphi(\ell, \phi, k) = \tau \quad \text{if and only if} \quad T_{\ell,\phi,k} = T_\tau \quad (4)$$

/0d/ any machine M_i can solve any task, starting at any time, assuming $I[i, \tau] > 0$.

/1a/ For any $\ell \leq L$ and $\phi \leq f_\ell$ the tasks of the block $B_{\ell,\phi}$ (see /0b/) can be solved (started) at any time, independently from each other (using any machine),

however the blocks $B_{\ell,2}, \dots, B_{\ell,\phi}, \dots, B_{\ell,\phi+1}$ must wait for finishing the previous ones:

/1b/ any task of $B_{\ell,\phi+1}$ may start only when each task of $B_{\ell,\phi}$ has already been finished.

In other words: for any $\ell \leq L$, $\phi_1 < \phi_2 \leq f_\ell$ and $k_1 \leq K_{\phi_1}$, $k_2 \leq K_{\phi_2}$ the task T_{ℓ,ϕ_2,k_2} can be started *only after* the task T_{ℓ,ϕ_1,k_1} has been finished.

We underline that /1b/ refers to blocks of the *same job* J_ℓ :

/1c/ we have *no* restriction at all for the starting times of the tasks of the job J_{ℓ_1} when comparing to J_{ℓ_2} , for $\ell_1 \neq \ell_2 \leq L$.

This means, especially, that

/1d/ all *tasks* (in the first block) of any job may be started even at time 0,

/2/ each machine M_i in every time may work on *at most one* task $T_{\ell,\phi,k}$, and M_i can not stop until it finishes the current task,

/3/ scheduling the task $T_{\ell,\phi,k}$ to the machine M_i means, that we choose a (positive) number d such that M_i is able to solve the task T_τ in the interval $(d, d + Ido[i, \tau])$, assuming $\tau = \varphi(\ell, \phi, k)$ and fulfilling /2/.

The goal is to finish all jobs as early as possible:

/4/ We have to finish all tasks of the jobs J_1, \dots, J_L such that each machine finishes all its tasks until the time I , satisfying /0a/ through /3/ and I is minimal.

Let us observe, that /1b/ is the hardest part of our algorithmic problem. Assumption /1a/ is void when $K_\phi=1$. The other extreme case is when $f_\ell=1$, in this case /1a/-/1c/ imply that all tasks of J_ℓ can be scheduled in arbitrary manner (fulfilling /0a/-/0c/ and /2/ of course).

Naturally all task $T_{\ell,\phi,k}$ we have to schedule, is a member of exactly one block.

The problem could be described and solved using Integer Linear Programming, too, but the number of variables and equations grows exponentially ([2],[3]).

2 The algorithms

Recall, that the input is (in fact) the *one-dimensional sequence* of the tasks

$$T_{1,1,1}, \dots, T_{\ell,\phi,k}, \dots, T_{L,\phi_{\text{fl}},K_{\text{fl}}} \quad (5)$$

where the terms $T_{\ell,\phi,k}$ are for $1 \leq \ell \leq L$, $1 \leq \phi \leq f_{\ell}$ and $1 \leq k \leq K_{\phi}$. Of course the input contains also the *delimiters* for determining the jobs and blocks in (4) according to /0a/ and /0b/. Single tasks without delimiters are considered one-element blocks: $B_{\ell,\phi} = (T_{\ell,\phi,1})$, i.e. $K_{\phi} = 1$.

Our **problem** is the *online scheduling*: we have to schedule each task $T_{\ell,\phi,k}$ *immediately* after reading it:

/5/ receiving $T_{\ell,\phi,k}$ we have to choose M_i and the starting time d such that M_i can solve $T_{\ell,\phi,k}$ in the interval $(d, d + \text{Ido}[i, \tau])$ without a break, where $\tau = \varphi(\ell, \phi, k)$.

Clearly, when deciding /5/, we have no information on the further tasks or on the length of the job or block we are working on, even not the number of jobs. Our scheduling in /5/ can not be altered later, of course. (One illustrative example can be found in the 3rd section.)

In our research we implemented, tested and compared the following variants (numberings refer to our developing):

Variante 3 : Give the next task $T_{\ell,\phi,k}$ to the machine M_i **if**

M_i can *finish* $T_{\ell,\phi,k}$ the soonest,

i.e. $d + \text{Ido}[i, \tau]$ has the possible smallest value, where $\tau = \varphi(\ell, \phi, k)$ and d is the starting time.

Variante 4 : Give the next task $T_{\ell,\phi,k}$ to the machine M_i **if**

M_i can *start* $T_{\ell,\phi,k}$ the soonest,

i.e. $d + \text{Ido}[i, \tau]$ has the possible smallest value.

Variante 5 : Give the next task $T_{\ell,\phi,k}$ to the machine M_i **if**

M_i can *solve* $T_{\ell,\phi,k}$ in the shortest time,

i.e. $\text{Ido}[i, \tau]$ has the possible *smallest* value.

All the three variants are clearly greedy ones, but in the great level of lack of information, we have no other simple idea for solving PMS.

3 Running experiments

So far we have tested the above algorithms on large size input datasets, in which each block contained *only one task*, i.e. when the scheduling has *no* possibility for solving tasks parallel. Further research for the solution of the general problem is in progress, will be summarized in [5].

Let us start with a short illustrative example:

Input (Ls4.dat):

```

8 = t = number of Task-types
3 = H = number of Machines
2 = K = number of tasks per Jobs
10 = L = number of Jobs
-----+
2 5 1 9 5 6 7 2 / Machine 1. | = I[h,τ]
3 4 8 7 4 1 3 1 / Machine 2. | table of times
4 6 9 8 3 4 6 3 / Machine 3. | necessary for
-----+ Mh for Tτ

1 2 / Job 1.
3 4 / Job 2.
1 3 / Job 3.
2 3 / Job 4.
3 3 / Job 5.
2 6 / Job 6.
5 2 / Job 7.
3 5 / Job 8.
1 6 / Job 9.
4 8 / Job 10.
0 / END

```

The solution of variant)3(can be seen on Figure 1.

Before any running experiments or theoretical investigations one might have the following *prejudices* (only few of them were justified during the runs):

" Variant)3(may give the *best* scheduling solution but it runs slowest (since it is the most precise?) ",

" Variant)4(may be the *quickest* but it provides bad solution (since it deals with the starting time only?) ",

" Variant)5(may be *quickest* but it provides not so good + not so bad solution (since it may give many tasks to some M_{h_0}) ".

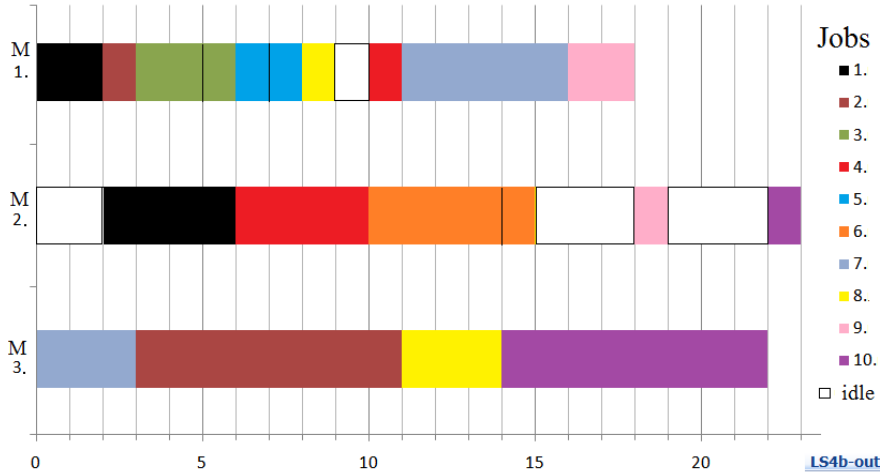


Figure 1: A solution of the example Ls4.dat

We tested the variants $\mathcal{J}3(-)\mathcal{J}5($ with *many* (more than 600) *large*, randomly (uniformly) generated datasets, each of them containing 500-1000 jobs. We used a personal computer with Intel® Core™ **Quad** CPU Q6600 @ 2.40GHz, 4GB RAM, Windows 7 and Delphi 7 language.

The summary of one of the runs can be seen in [4], which includes and compares schedulig- and running times, too.

Let us first serve the details of running of *medium size* datasets. N stands for the number of the datasets. Determining

$$1 \leq t \leq 100, 1 \leq m \leq 100, 1 \leq \mathbf{I}(h, \tau) \leq 100, 1 \leq L \leq 200, 1 \leq K \leq 20, N=500, \quad (6)$$

the average *running time* (for each dataset) were some minues, the *sizes* of the ouput files were 10kb - 1Mb separately, 100Mb total. These data are valid for all the three variants.

There were some differences in *running times*:

variant $\mathcal{J}5($ is highly faster than $\mathcal{J}3($, in detail :

in **14%** cases $0.5 * \mathcal{J}3(< \mathcal{J}5(\leq 0.9 * \mathcal{J}3($

in **75%** cases $\mathcal{J}5(< 0.5 * \mathcal{J}3($

However, the resulted *scheduling times* (solutions) were totally different (see [4] for details):

variant $\mathcal{J}4($ has extremaly bad solutions in almost all cases,

variant $\mathcal{J}3($ has **the best** solution with few exceptions:

in **6%** cases $\mathcal{J}5($ is better, than $\mathcal{J}3($,

in **17%** cases \mathcal{A} is as good as \mathcal{B} ,
 variant \mathcal{A} has **not so bad** solutions:
 in **29%** cases $1.0 * \mathcal{B} < \mathcal{A} < 1.2 * \mathcal{B}$
 in **42%** cases $1.2 * \mathcal{B} \leq \mathcal{A} < 2.0 * \mathcal{B}$

As we observed, the above data *slightly depend* upon the size of the problem.

Now, let us turn to the experiment of *large* datasets. Relying on the bad experiences in the previous subsection, we *excluded* variant \mathcal{A} from our further experiments. Our settings were:

$$1 \leq t, m, \mathbf{I}(h, \tau), L, K \leq 1000, \quad N=650. \quad (7)$$

The average *running times* (for all the N=650 datasets, total) :

variant \mathcal{B} runned for 81 hours (!),

variant \mathcal{A} runned for 2 hours,

i.e. the difference is large: variant \mathcal{A} is tremendously faster than \mathcal{B} :

in **23%** cases $0.5 * \mathcal{B} < \mathcal{A} \leq 0.50 * \mathcal{B}$,

in **75%** cases $\mathcal{A} < 0.05 * \mathcal{B}$!

The *sizes* of output files (separately) were between 1Mb and 50Mb, 4Gb (!) total for both these variants.

The rate of resulted *scheduling times* showed opposite direction:

variant \mathcal{B} has **always the best** solution, \mathcal{A} is **fairly bad**

in **11%** cases $1.0 * \mathcal{B} < \mathcal{A} < 1.2 * \mathcal{B}$,

in **29%** cases $1.2 * \mathcal{B} \leq \mathcal{A} < 2.0 * \mathcal{B}$,

in **33%** cases $2.0 * \mathcal{B} \leq \mathcal{A} < 3.0 * \mathcal{B}$,

in **25%** cases $3.0 * \mathcal{B} \leq \mathcal{A}$.

Conclusions

In general, but especially in large size datasets, version \mathcal{A} was exponentially faster than version \mathcal{B} , but considering the resulted schedulings, version \mathcal{A} gave worse results than version \mathcal{B} , in *moderate manner*. This shows again the old dilemma: "shorter running time" *versus* "better results " !

We have *no* comparison with the *absolute* (offline) optimum, but \mathcal{B} **might be optimal** (in the offline sense) **in many cases**, since there are *very few idle time* (pause) of the machines, and they finish almost at the same "moment".

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