Non-Cooperative Scheduling of Multiple Bag-of-Tasks Applications TU Wien Seminar

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Motivation

Large-scale distributed platforms result from the collaboration of many users:

- Multiple applications execute concurrently on heterogeneous platforms and compete for CPU and network resources.
- Sharing resources amongst users should somehow be fair.
- Large scale platform require distributed scheduling.

In a grid or Volunteer Computing context, sharing is generally done in the "low" layers (network, OS).

- We analyze the behavior of K non-cooperative schedulers that use the optimal strategy to maximize their own utility while fair sharing is ensured at a system level ignoring applications characteristics [INFOCOM'07].
- We use Lagrangian optimization and distributed gradient descent to propose a fair and distributed scheduling algorithm for this framework [JPDC'13?].

Outline

Non-cooperative Scheduling

- Framework and Notations
- Non-cooperative Scheduling
- Measuring Efficiency
- Fairness, Pricing, Lagrangian Optimization and Distributed Gradient
 - Model and Notations
 - Lagrangian Optimization
 - A Carefull Investigation



Hypotheses :

- ► N processors with processing capabilities W_n (in Mflop.s⁻¹)
- using links with capacity B_n (in Mb.s⁻¹)

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Hypotheses :

Multi-port

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Communications to P_i do not interfere with communications to P_j .



Hypotheses :

- Multi-port
- No admission policy but an ideal local fair sharing of resources among the various requests

- ► N processors with processing capabilities W_n (in Mflop.s⁻¹)
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Hypotheses :

Definition.

We denote by platform-system a triplet (N, B, W) where N is the number of machines, and B and W the vectors of size N containing the link capacities and the computational powers of the machines.

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bilities W_n (in Mflop.s⁻¹)

▶ Multiple applications (*A*₁,...,*A*_{*K*}):



- each consisting in a large number of same-size independent tasks
- Different communication and computation demands for different applications. For each task of Ak:
 - processing cost w_k (MFlops)
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- Such applications are typical Desktop Grid or Volunteer Computing applications (SETI@home, Einstein@Home, ...)

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Definition.

We define an application-system as a triplet (K, b, w) where K is the number of applications, and b and w the vectors of size K representing the size and the amount of computation associated to the different applications.

Steady-State scheduling

In the following our K applications run on our N workers and compete for network and CPU access:

Definition.

A system S is a sextuplet (K, b, w, N, B, W), with K, b, w, N, B, W defined as for a user-system and a platform-system.

- ► Task regularity ~> steady-state scheduling.
- Maximize throughput (average number of tasks processed per unit of time)

$$\rho_k = \lim_{t \to \infty} \frac{done_k(t)}{t}$$

Similarly: $\rho_{n,k}$ is the average number of tasks of type k performed per time-unit on the processor P_n .

$$\rho_k = \sum_n \rho_{n,k}.$$

• ρ_k is the utility of application k.

The scheduler of each application thus aims at maximizing its own throughput, i.e. ρ_k .

However, as applications use the same set of resources, we have the following general constraints:

Computation
$$\forall n \in [\![0, N]\!] : \sum_{k=1}^{K} \rho_{n,k} \cdot w_k \leqslant W_n$$

Communication $\forall n \in [\![1, N]\!] : \sum_{k=1}^{K} \rho_{n,k} \cdot b_k \leqslant B_n$

Applications should decide:

- which worker to use,
- when to send data from the master to a worker,
- when to use a worker for computation.

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Optimal strategy for a single application

Single application

This problem reduces to maximizing $\sum_{n=1}^{N} \rho_{n,1}$ while: $\begin{cases} \forall n \in \llbracket A, N \rrbracket : \rho_{n,1} \cdot w_1 \leqslant W_n \\ \forall n \in \llbracket 1, N \rrbracket : \rho_{n,1} \cdot b_1 \leqslant B_n \\ \forall n, \quad \rho_{n,1} \ge 0. \end{cases}$

The optimal solution to this linear program is obtained by setting

$$\forall n, \ \rho_{n,1} = \min\left(\frac{W_n}{w_1}, \frac{B_n}{b_1}\right)$$

In other words

The master process should saturate each worker by sending it as many tasks as possible.

A simple acknowledgment mechanism enables the master process to ensure that it is not over-flooding the workers, while always converging to the optimal throughput. We suppose a purely non-cooperative game where no scheduler decides to "ally" to any other (i.e. no coalition is formed).

As the players constantly adapt to each others' actions, they may (or not) reach some equilibrium, known in game theory as Nash equilibrium [Nash51].

 $\rho_{n,k}^{(nc)}$ denotes the rates achieved at such stable states (if any).

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A simple example

Two computers 1 and 2: $B_1 = 1$, $W_1 = 2$, $B_2 = 2$, $W_2 = 1$. Two applications: $b_1 = 1$, $w_1 = 2$, $b_2 = 2$ and $w_2 = 1$.



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Cooperative Approach:

Application 1 is processed exclusively on computer 1 and application 2 on computer 2. The respective throughput is $\rho_1^{(\text{coop})} = \rho_2^{(\text{coop})} = 1.$



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Non-Cooperative Approach: $\rho_1^{(nc)} = \rho_2^{(nc)} = \frac{3}{4}$





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Resource Usage



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Computation



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Computation



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We assume $c_1 \leq c_2 \leq \cdots \leq c_K$. Let us denote by \mathcal{W}_n the set of players that are computation-saturated and by \mathcal{B}_n the set of players that are communication-saturated on a given arbitrary worker n.

• If
$$\sum_{k} \frac{C_n}{c_k} \leqslant K$$
 then $\mathcal{W}_n = \emptyset$ and $\forall k, \rho_{n,k}^{(nc)} = \frac{B_n}{K.b_k}$.
• Else, if $\sum_{k} \frac{c_k}{C_n} \leqslant K$ then $\mathcal{B}_n = \emptyset$ and $\forall k, \rho_{n,k}^{(nc)} = \frac{W_n}{K.w_k}$.

● Else, \mathcal{B}_n and \mathcal{W}_n are non-empty and there exists an integer $m \in$ [[1; K - 1] such that

$$\frac{c_m}{C_n} < \frac{m - \sum_{k=1}^m \frac{c_k}{C_n}}{K - m - \sum_{k=m+1}^K \frac{C_n}{c_k}} < \frac{c_{m+1}}{C_n}.$$

Then, we have $\mathcal{W}_n=\{1,\ldots,m\}$ and $\mathcal{B}_n=\{m+1,\ldots,K\}$ and

$$\begin{pmatrix}
\rho_{n,k}^{(nc)} = \frac{B_n}{b_k} \frac{|\mathcal{W}_n| - \sum_{p \in \mathcal{W}_n} \frac{c_p}{C_n}}{|\mathcal{W}_n||\mathcal{B}_n| - \sum_{p \in \mathcal{W}_n} c_p \sum_{p \in \mathcal{B}_n} \frac{1}{c_p}} & \text{if } k \in \mathcal{B}_n \\
\rho_{n,k}^{(nc)} = \frac{W_n}{w_k} \frac{|\mathcal{B}_n| - \sum_{p \in \mathcal{B}_n} \frac{C_n}{c_p}}{|\mathcal{W}_n||\mathcal{B}_n| - \sum_{p \in \mathcal{W}_n} c_p \sum_{p \in \mathcal{B}_n} \frac{1}{c_p}} & \text{if } k \in \mathcal{W}_n
\end{cases}$$

Theorem 1.

For a given system (N, B, W, K, b, w) there exists exactly one Nash Equilibrium and it can be analytically computed.

Proof.

Under the non-cooperative assumption, on a given worker, an application is either communication-saturated or computation-saturated. Puting schedules in some canonical form enables to determine for each processor, which applications are communication-saturated and which ones are computation-saturated and then to derive the corresponding rates.

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Definition: Pareto optimality.

An allocation is said to be Pareto-optimal if it is impossible to strictly increase the throughput of an application without strictly decreasing the one of another.

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When is our Nash Equilibrium Pareto-optimal ?

Theorem 2.

The allocation at the Nash equilibrium is Pareto inefficient if and only if there exists two workers, namely n_1 and n_2 such that all applications are communication-saturated on n_1 and computation-saturated on n_2 (i.e. $\sum_k \frac{B_{n_1}}{W_{n_1}} \frac{w_k}{b_k} \leq K$ and $\sum_k \frac{b_k}{w_k} \frac{W_{n_2}}{B_{n_2}} \leq K$).

Corrolary: on a single-processor system, the allocation at the Nash equilibrium is Pareto optimal.

Braess-like Paradox

Pareto-inefficient equilibria can exhibit unexpected behavior.

Definition: Braess Paradox [Braess68].

There is a Braess Paradox if there exists two systems $ini \mbox{ and } aug$ such that

 $ini < aug \text{ and } \rho^{(nc)}(ini) > \rho^{(nc)}(aug).$

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Theorem 3.

In the non-cooperative multi-port scheduling problem, Braess like paradoxes cannot occur.

Proof.

- Defining an equivalence relation on sub-systems.
- Defining an order relation on equivalent sub-systems.

Measuring Pareto-Inefficiency Price of Anarchy

Definition [Koutsoupias.Papadimitriou'98] Price of Anarchy:

$$\phi_{\Sigma} = \max_{S} \frac{\sum_{k} \rho_{k}^{(\Sigma)}(S)}{\sum_{k} \rho_{k}^{(nc)}(S)} \ge 1.$$

Measuring Pareto-Inefficiency Price of Anarchy

Definition [Koutsoupias.Papadimitriou'98] More generally:

$$I_f(S) = \frac{f\left(\rho_1^{(f)}(S), \dots, \rho_K^{(f)}(S)\right)}{f\left((\rho_1^{(nc)}(S), \dots, \rho_K^{(nc)}(S)\right)} \ge 1.$$
Measuring Pareto-Inefficiency Price of Anarchy

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Problem measures the "distance" to a particular point... Illustration 1 machine $(B_1 = W_1 = 1)$ and K applications b = $(\frac{1}{M}, 1...1)$ and $w = (\frac{1}{M}, 1...1)$. Utility Set $I_{\Sigma}(S_{M,K}) = \frac{M}{\frac{M}{K} + \frac{K-1}{K}}$ Max-min Allocation Nash Equilibrium ρ_k $\xrightarrow{M \to \infty} K$ Profit Allocation Utility set and allocations 0 M0 $S_{M,K}$ (K = 3, M = 2). ρ_1

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Measuring Pareto-Inefficiency

Selfishness Degradation Factor



Degradation Factor is related to ε -approximation of Pareto-curves [Papadimitriou.Yannakakis'00].

In our context, Selfishness Degradation Factor is at least 2.

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Pareto Optimality and Monotonicity of Performance Measures



Most classical performance measures decrease with resource

augmentation!

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Recap

Conclusion

- Applying fair and optimal sharing on each resource does not ensure any fairness nor efficiency when users do not cooperate.
- Being "locally Pareto optimal" (i.e on each single machine) does not help being Pareto optimal.
- Even with Pareto optimal situations, classical performance measures can be non monotonic.

 \sim either applications cooperate or new complex and global access policies should be designed

Other Future Work

- Measuring Pareto-inefficiency is an open question.
- In the one-port communication model, Braess-like paradoxes seem to arise.

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- Communications and computations can be overlapped.
- Multi-port communication model.

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Application Model

Multiple applications:

• A set A of K applications A_1, \ldots, A_K



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- ► Each consisting in a large number of same-size independent tasks ~> each application is defined by a communication cost w_k (in MFlops) and a communication cost b_k (in MB)
- Different communication and computation demands for different applications





Each application originates from a master node P_{m(k)} that initially holds all the input data necessary for each application A_k

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Therefore if an application k wants to use a node P_n , all its data will use a single path from $P_{m(k)}$ to P_n denoted by $(P_{m(k)} \sim P_n)$



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All deployment trees may be different

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 - Variables need to respect the following constraints

$$\begin{split} \forall n, \quad \sum_{k} \rho_{n,k} w_k \leqslant W_n \\ \forall (P_i \rightarrow P_j), \quad \sum_{k}^{k} \sum_{\substack{(P_i \rightarrow P_j) \in (P_{m(k)} \sim P_n)}} \rho_{n,k} b_k \leqslant B_{i,j} \end{split}$$

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We would like to maximize all throughputs ρ_k .

"Defining" fairness is one way to go from a multi-criteria problem to a more classical mono-criteria problem.

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Utility Function

In a general context, each application is characterized by a utility function u_k defined on $(\rho_{n,k})_{1 \le k \le K, 1 \le n \le N}$.

In our context, the utility is simply defined by

$$u_k(\rho) = \sum_n \rho_{n,k} = \rho_k$$

But we could perfectly imagine other utility functions:



How can fair be defined? Does it always mean "give the same thing to everyone"? How can efficiency be defined?



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Fairness can be seen as the trade-off between individual satisfaction and global satisfaction.



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• A few problems with max-min fairness:

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Not "efficient" when applications are very different



A few problems with max-min fairness:

- Not "efficient" when applications are very different
- Seems to be hard to reach on such platforms [TPDS'07]



A few problems with max-min fairness:

- Not "efficient" when applications are very different
- Seems to be hard to reach on such platforms [TPDS'07]
- Let's try proportional fairness!

MAXIMIZE
$$\left(\sum_{k \in K} \log \rho_k\right)$$

A new optimization problem

$$\begin{split} \text{Maximize } \sum_{k} \log \left(\sum_{n} \rho_{n,k} \right) \text{ under the constraints:} \\ \forall n, \quad \sum_{k} \rho_{n,k} w_k \leqslant W_n \\ \forall (P_i \to P_j), \quad \sum_{k}^{k} \sum_{\substack{(P_i \to P_j) \in (P_m(k) \rightsquigarrow P_n)}} \rho_{n,k} b_k \leqslant B_{i,j} \end{split}$$

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A new optimization problem

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$$\sum_{k} \log \left(\sum_{n} \rho_{n,k} \right)$$
 under the constraints:
 $\forall n, \quad \sum_{k} \rho_{n,k} w_k \leq W_n$
 $\forall (P_i \rightarrow P_j), \quad \sum_{k}^{k} \sum_{\substack{(P_i \rightarrow P_j) \in (P_m(k) \sim P_n)}} \rho_{n,k} b_k \leq B_{i,j}$

Can be solved in polynomial time e.g., with semi-definite programming. It is very centralized though.

Can we solve it in a distributed way?

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Designed to solve non linear optimization problems:

- Let $\rho \to f(\rho)$ be a function to maximize.
- Let $(C_i(\rho) \ge 0)_{i \in [1..n]}$ be a set of n constraints.
- We wish to solve:

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- ► The Lagrangian function: $\mathcal{L}(\rho, \lambda) = f(\rho) \sum_{i \in [1..n]} \lambda_i C_i(\rho).$
- Under mild assumptions, there is no duality gap [Bertsekas-Tsitsiklis]:

$$\max_{\rho \ge 0} \min_{\lambda \ge 0} \mathcal{L}(\rho, \lambda) = \min_{\substack{\lambda \ge 0 \ \rho \ge 0}} \max_{\rho \ge 0} \mathcal{L}(\rho, \lambda) \stackrel{\text{def}}{=} \min_{\lambda \ge 0} d(\lambda)$$

Primal problem (P) Dual problem (D)



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So what?..

- ► Two coupled problems with simple constraints.
- \mathcal{L} is concave in ρ and convex in λ
- ► The structure of constraints is transposed to (*D*) and a gradient descent algorithm is a natural way to solve these two problems.
- This technique has been used successfully for network resource sharing [Kelly.98], TCP analysis [Low.03], flow control in multi-path network [Hang.et.al.03], ...

Tsitsiklis]:

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Primal problem (P)

Trying to use Lagrangian optimization

What does the Lagrangian function look like ?

$$\mathcal{L}(\rho,\lambda,\mu) = \sum_{k \in K} \log\left(\sum_{i} \rho_{i,k}\right) + \sum_{i} \lambda_{i} \left(W_{i} - \sum_{k} \rho_{i,k} w_{k}\right) + \sum_{(P_{i} \to P_{j})} \mu_{i,j} \left(B_{i,j} - \sum_{k} \sum_{\substack{n \text{ such that} \\ (P_{i} \to P_{j}) \in (P_{m(k)} \rightsquigarrow P_{n})} \rho_{n,k} b_{k}\right)$$

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$$+ \sum_{(P_{i} \to P_{j})} \mu_{i,j} \left(B_{i,j} - \sum_{k} \sum_{\substack{n \text{ such that} \\ (P_{i} \to P_{j}) \in (P_{m(k)} \rightsquigarrow P_{n})} \rho_{n,k} b_{k}\right)$$

► Remember, we want to compute $\min_{\lambda,\mu \ge 0} \max_{\rho \ge 0} \mathcal{L}(\rho, \lambda, \mu)$.

 \mathcal{L} is concave-convex \rightsquigarrow simple "alternate" gradient descent (I'm skipping a few details here to keep it simple and just present the general idea)

Update equations:
$$\begin{cases} \rho_{i,k} & \leftarrow \rho_{i,k} + \gamma \frac{\partial \mathcal{L}}{\partial \rho_{i,k}} \\ \lambda_i & \leftarrow \lambda_i - \gamma \frac{\partial \mathcal{L}}{\partial \lambda_i} \\ \mu_{i,j} & \leftarrow \mu_{i,j} - \gamma \frac{\partial \mathcal{L}}{\partial \mu_{i,j}} \end{cases}$$

Toward a Distributed Algorithm...

- $\rho_{i,k}$ is "private" to the agent of application k running on node i
- ▶ λ_i is attached to node *i* and $\mu_{i,j}$ is attached to $(P_i \rightarrow P_j)$

Toward a Distributed Algorithm...

- $\rho_{i,k}$ is "private" to the agent of application k running on node i
- ▶ λ_i is attached to node i and µ_{i,j} is attached to (P_i → P_j) λ_i and µ_{i,j} are called shadow variables or shadow prices. They can naturally thought of as the price to pay to use the corresponding resource.
- ► A gradient descent algorithm on the primal-dual problem can thus be seen as a bargain between applications and resources.
Toward a Distributed Algorithm...



Toward a Distributed Algorithm...



Toward a Distributed Algorithm...

Prices and rates can thus be propagated and aggregated to perform the following updates:

$$p_{k}^{i}(t+1) \leftarrow b_{k}\eta_{k}^{i}(t) + w_{k}\lambda_{i}(t)$$

$$\rho_{k}(t+1) \leftarrow \sigma_{k}^{m(k)}(t+1)$$

$$\rho_{i,k}(t+1) \leftarrow \left[\rho_{i,k}(t) + \gamma_{\rho}(U_{k}'(\rho_{k}(t)) - p_{k}^{i}(t))\right]^{+}$$

$$\lambda_{i}(t+1) \leftarrow \left[\lambda_{i}(t) + \gamma_{\lambda}\left(\sum_{k} w_{k}\rho_{i,k} - W_{i}\right)\right]^{+}$$

$$\mu_{i,j}(t+1) \leftarrow \left[\mu_{i,j}(t) + \gamma_{\mu}\left(\sum_{k} b_{k}\sigma_{k}^{i} - B_{i,j}\right)\right]^{+}$$

- This algorithm is fully distributed and converges to the optimal solution provided a good choice of γ_ρ, γ_λ and γ_μ is done.
- This algorithm should seamlessly adapts to application/node arrival and to load variations.

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Non-Cooperative Scheduling

Illustration of convergence on a toy platform



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There are three main steps to come up with such an algorithm:

- **O Modeling:** concave non-linear maximization problem
- **2** Partial derivatives: differentiate \mathcal{L}
- Algorithm design: ascent on primal and descent on dual exploiting the structure of derivatives

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"Technical" issues for convergence:

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"Technical" issues for convergence:

- Beware of divisions by 0
- Strict convexity of the objective function (\rightsquigarrow proximal variables $\tilde{\rho}$)

 $\max_{\widetilde{\rho} \geqslant 0} \max_{\rho \geqslant 0} \min_{\lambda \geqslant 0} L(\widetilde{\rho}, \rho, \lambda),$

Nested structure needs to be broken in practice!!!

And in practice ?

In 2006, we stumbled on [Hang.et.al.03]: "Optimal Flow Control and Routing in Multi-path Networks", where we learnt about these techniques

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Effectiveness of the approach was illustrated on a few simple cases



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And in practice ?

In 2006, we stumbled on [Hang.et.al.03]: "Optimal Flow Control and Routing in Multi-path Networks", where we learnt about these techniques

Using the same formulas did not work out well...



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Outline

1 Non-cooperative Scheduling

- Framework and Notations
- Non-cooperative Scheduling
- Measuring Efficiency

Fairness, Pricing, Lagrangian Optimization and Distributed Gradient

- Model and Notations
- Lagrangian Optimization
- A Carefull Investigation

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- SimGrid based simulation
- Check correctness with semi-definite programming
- Adjusting 4 different steps at a time seems tricky:
 - use designed experiments and Analysis of Variance to assess the effectiveness of step modification
 - Use the same 30 platforms for each step configuration and distinguish platform variability from algorithm variability

Experimental Setting

A typical experiment

-								
		$\gamma_{ ho}$	$\gamma_{\widetilde{ ho}}$	γ	μ	γ_{λ}		
low (-	1)	0.001 0.	001	1.00	e-15	1.00e-1	5	
high (1)́	0.1 ().1	1.00	e-13	1.00e-1	3	
	, ,		<i>c</i>	<u> </u>		1 .		
(a)	Par	rameters	tor	fact	orial	desigr	۱ I	
• • •						-		
	Dt	Sum Sq		Mean Sq	F value	Pr(>F)		
ρ	1	634854.4356	6348	354.4356	75.1771	0.0000	***	
$\tilde{\rho}$	1	79293.1822	792	293.1822	9.3896	0.0023	**	
λ	1	39882.0712	398	382.0712	4.7227	0.0303	*	
μ	1	32497.4344	324	197.4344	3.8482	0.0504		
platform	29	470172.3060	162	212.8381	1.9199	0.0032	**	
$\rho: \tilde{\rho}$	1	67441.2012	674	41.2012	7.9861	0.0049	**	
$\rho:\lambda$	1	27584.2533	275	584.2533	3.2664	0.0714		
$\tilde{\rho}: \lambda$	1	330.0890	1 3	330.0890	0.0391	0.8434		
ρ:μ	1	38450.8702	384	150.8702	4.5532	0.0334	*	
õ: u	1	31651.5862	316	551.5862	3.7481	0.0535		
$\lambda : \mu$	1	14136.1931	141	136.1931	1.6740	0.1964		
$\rho: \tilde{\rho}: \lambda$	1	1489.7248	14	189.7248	0.1764	0.6747		
0: õ: u	1	35856.2101	358	356.2101	4.2460	0.0399	*	
ο:λ:μ	1	16345.9826	163	345.9826	1.9356	0.1649		
$\tilde{\rho}: \lambda: \mu$	1	16.1140		16.1140	0.0019	0.9652		
$\rho: \tilde{\rho}: \lambda: \mu$	1	158 4895	1	58 4895	0.0188	0.8911		
Residuals	435	3673480.0717	84	44 7818	0.0100			

(c) ANOVA results: * means that parameter is significant



Conclusion of the investigation

Here is what we found out:

The algorithm works great when using identical applications, even with 100 nodes

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Primal $\Delta \rho_{n,k} = -\alpha.L$ $\neq -\alpha \Delta \lambda_n$ Dual

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(3) Division by small values and discontinuities: $[x(t)]^+ = \max(0, x(t)).$

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- Slow convergence: use quasi-Newton scheme
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 Combining these three techniques provides a distributed algorithm that converges even for platforms with 500 nodes.

Open remarks

Similitude between fair steady-state scheduling and flow control in multi-path networks motivated the Lagrangian approach.

All the convergence issues would have been overlooked if we had been less stubborn

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Theory vs. "Practice"

Theory difficulty lies in non strict convexity. Practice scaling and borders are the most important issues. Using designed experiments we can find "robust" step sizes Similitude between fair steady-state scheduling and flow control in multi-path networks motivated the Lagrangian approach.

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Theory vs. "Practice"

Theory difficulty lies in non strict convexity. Practice scaling and borders are the most important issues. Using designed experiments we can find "robust" step sizes

► The exponential decay reminds of a barrier function ~> interior point methods ? I still lack perspective on these problems but such optimization techniques are still ongoing research.

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