



Sizing Tasks for Scientific Workflows with Reinforcement Learning

Nicolas Zunker

n.zunker@campus.tu-berlin.de

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Mobile Cloud Computing Chair

Technische Universität Berlin

Examiner 1: Prof. Dr. Odej Kao

Advisor: Jonathan Bader

Declaration

I hereby declare that I have created the present work independently and by my own without illicit assistance and only utilizing the listed sources and tools.

Hiermit erkläre ich, dass ich die vorliegende Arbeit selbstständig und eigenhändig sowie ohne unerlaubte fremde Hilfe und ausschliesslich unter Verwendung der aufgeführten Quellen und Hilfsmittel angefertigt habe.

Die selbstständige und eigenständige Anfertigung versichert an Eides statt:

DATE

NAME

Acknowledgment

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INITIALS

Abstract

The growth of cloud computing and big data has lead to increased use of computing clusters to process large datasets. Particularly in the sciences this has become an enormous advantage for analysis and research and it has become common for scientists to break down their computing needs into a sequence of smaller tasks, a so called workflow. These workflows can then be run on computer clusters and cloud computing platforms. One of the most pertinent fields for this is Bioinformatics. Another result of the rise of cloud computing is that there has also been a paradigm shift towards designing programs to be as best suited as possible for cluster environments and this has also led to a shift towards containerization due to the inherent advantage of their 'write once, deploy anywhere' philosophy for such clusters. The result of this combination of factors is the increased popularity of scientific workflows and the frameworks or scripting languages which can be used to make such workflows scalable and reproducible, through containerization, and prepare them for deployment to various cloud computing platforms and management softwares (e.g. Kubernetes, AWS etc.). Nextflow is one such bioinformatics workflow manager. It includes both a domain specific language to compose tasks into portable workflows as well as the workflow management software for deployment to the various different execution platforms with which it is compatible.

Since workflows are composed of segregated inter-dependent tasks which can run in their own containers, the individual tasks which make up a workflow can be assigned differing computational resources and doing so can speed up the entire process and/or make it more efficient. This is naturally of particular interest to the scientists running these workflows since it will either lower the costs or speed up the process or both. This thesis aims to investigate the allocation of resources to individual tasks and specifically how reinforcement learning can be applied to improve the allocation to be as optimal as possible. Since the tasks within these scientific workflows will reoccur quite often the use of an algorithm which learns "on the go" presents a potential for near optimal resource allocation in the long run. The implementation will be integrated into the nextflow source code.

Differing approaches and reinforcement learning algorithms will be compared and their performance will be judged firstly by how they minimize runtime and secondly by their maximization of resource usage. The primary approach will use a simple gradient bandit and may be compared to other approaches such as TD-learning.

Should such an algorithm prove useful and provide an improvement in resource usage it would naturally indicate this is an area which should be explored further and that the use of these methods will help make scientific workflows more efficient. This would be helpful to both the scientists which use workflow managers as well as the administrators of the execution platforms on which such workflow managers will run.

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Chapter 1

Introduction

This chapter gives a general introduction into the content of and the motivation for this thesis. This is done by initially outlining the motivation and problem in section [1.1](#), defining the goal in section [1.2](#), and explaining the structure of the thesis in section [1.3](#).

1.1 Motivation & Problem Description

Thus, many executives feel fear to become unconscious [**TheInquisitiveMind**]. Also, I want to buy **smart home devices** such as the ones listed in table ??.

1.2 Goal of the Thesis

The goal of this thesis is to XXXXX. The contributions are:

1. Analysis ...
2. Design ...

3. Prototypical implementation of the design.

1.3 Structure of the Thesis

The remainder of the thesis is structured as follows

Chapter 2

Related Work

This chapter discusses other works in related areas and discusses the nuances of their approaches and results.

2.1 Reinforcement Learning to Schedule Tasks/Jobs

In the LearningToSchedule... paper the researchers attempted to use reinforcement learning to schedule data-parallel processing jobs in shared clusters. In particular they considered resource management systems such as YARN or Mesos. Their approach was to schedule the jobs based on co-location goodness, which is a metric for how well a job can share the resources of the machine which it is scheduled on with another job. This metric was based on cpu usage and disk and network IO since these two activities tend to be very compatible, while the job waits on the disk or network another job can be free to use the cpu and vice versa. The paper considered three different scheduling approaches: FIFO, a gradient bandit which starts new each time and a gradient bandit which starts with the data from previous runs. In the end the best performing approach was the bandit which had the benefit of the data from previous runs. This was because that bandit skipped the exploration phase during

which bandit sacrifice performance for learning by virtue of having the previous runs' data. This gradient bandit approach was then incorporated into three different schedulers called Mary, Hugo and Hugo*. Mary schedules based on co-location goodness, Hugo extends this by aggregating jobs into groups based on their similarities and scheduling the groups, and Hugo* does the same as Hugo but also considers how long a task has been waiting.

In the SCARL paper researchers use reinforcement learning for scheduling jobs among machines with heterogeneous resources. Their approach was based on combining attentive representation and the policy-gradient method. Attentive representation is a technique for focusing attention more quickly within a neural network. The model which they used to represent the problem used the allocation of jobs to machines as the state whilst the available actions at any given point were the scheduling of jobs. The reward function was based on a slowdown metric: $(elapsed_time * penalty_factor) / computation_time$. Ultimately the researchers found that for high levels of heterogeneity SCARL outperformed the shortest job first (SJF) metric by 10% and for lower levels of heterogeneity it outperformed it by 4 %.

2.2 Reinforcement Learning to Allocate Resources

SmartYARN applied a Q-learning approach to balance resource usage optimization against application performance. This is one of the central considerations for any client of a cloud computing platform- the need to balance minimizing cost by using less resources against the need to increase runtime by using more resources. In this paper the researchers used the performance of the app under a certain resource configuration as the state-space and the actions available to the agent were increasing or decreasing one unit of cpu or memory or keeping the previous allocation. In the end the researchers found that the agent was able to achieve 98% of the optimal cost reduction and generally performed at the optimal level, finding the optimal allocation the vast majority of the time.

VMConf tackled a similar problem- configuring the resource allocations of virtual machine's (VM's) using reinforcement learning. Their approach was a continuing discounted MDP with Q-learning. The state space was a triple of CPU Time credits (used for scheduling cpu time), virtual cpu's (vCPU's), and memory. The agent's available actions were increasing, decreasing or leaving the allocations, with only one resource allowed to be changed per action. As a reward the ratio of the achieved throughput to a reference throughput was used. One key trick used by the researchers was to use model based reinforcement learning. By modeling states the agent is able to simulate or predict the reward it can expect from previously unseen action-state pairs, whereas the classic approach requires the agent to experiment with each action-state pair. This means the model based agent is able to learn much faster and can enter its exploitation phase earlier than the static agent which must explore the action-state space for a long time when there are large state spaces. Ultimately however the researchers found that the static agent performed quite well and achieve high levels of throughput in its own regard, although the model-based approach consistently outperformed the static one.

Finally there is the DeepRM paper on resource management using deep reinforcement learning. These researchers used a policy gradient reinforcement learning algorithm combined with a neural network. The state-space consisted of the resources of the cluster and the resource requirements of arriving jobs, represented as images which could be fed into the neural network. The available actions were simply to allocate the jobs. In order to speed up the process the agent was given the option to schedule up to M jobs in a row, thus enabling it to complete leap forward by as many as M timesteps instead of always progressing by a single timestep. The reward given to the agent was the negative average slowdown. In the end the shortest job first metric and a very strong, heuristic based algorithm called Tetris were both outperformed by the researchers' approach. A key takeaway was that the agent achieved this increased performance by learning a policy to maximize the number of small jobs it completed. While the approach taken in the paper was still better with large jobs it was distinctly better with small jobs because the network learned to always keep a small amount of resources available to quickly service any small jobs which might arrive, whereas

the other approaches inevitably scheduled all the available resources so that small jobs also had to wait.

2.3 Takeaways

All of the papers discussed above managed to utilize reinforcement learning in an administrative fashion to improve the performance of jobs in their respective computing environments. A common theme among the papers concerned with resource utilization was to limit the action space to increasing, decreasing or keeping the current allocation of resources. The majority of the papers also utilized neural networks or deep reinforcement learning. It is also worth noting that most of these papers used runtime in their reward functions. A key difference between all of the above papers and this thesis is that these papers handled the allocation of resources among machines in their system whereas in nextflow the jobs are given a resource allocation and then passed on to a platform or cluster management software which handles the actual allocation of the job to a machine.

Appendix A

Technologies and Concepts of a Larger Context

This appendix chapter contains definitions for technologies and concepts that are mentioned in the thesis, but are not of higher importance for it. Section [A.1](#) introduces a consensus algorithm for distributed systems.

A.1 Finding Consensus in a Distributed System

The Paxos algorithm has first been described by Leslie Lamport in his work about the parliament on the island of Paxos [[paxos1998](#)]. The parliament's part-time legislators had been able to maintain consistent copies of their records by following the algorithm protocol. While the original work's main focus has not been the application in computer science, the author simplified his explanations later in [[paxos2001](#)] and described how Paxos can be used to find consensus in fault-tolerant distributed systems.

Even though the Paxos algorithm cannot mitigate Byzantine failures (see appendix [C](#)), it

can mitigate the effects of different processing speeds of participants and reordered, delayed, or lost messages. The only requirement is that at least half of the participants can somehow communicate.

In Paxos, four different roles exist. The **client** issues requests and waits for responses. Based on the client's request, multiple **proposer** propose a value. All these proposals are processed by the **acceptors** which will agree upon one at the end. Finally, there are the **learners** which guarantee persistence of agreed decisions and respond to client reads. The usual setup used in practice, in which every machine participating fulfills each of the roles besides the client role, is shown in figure A.1.

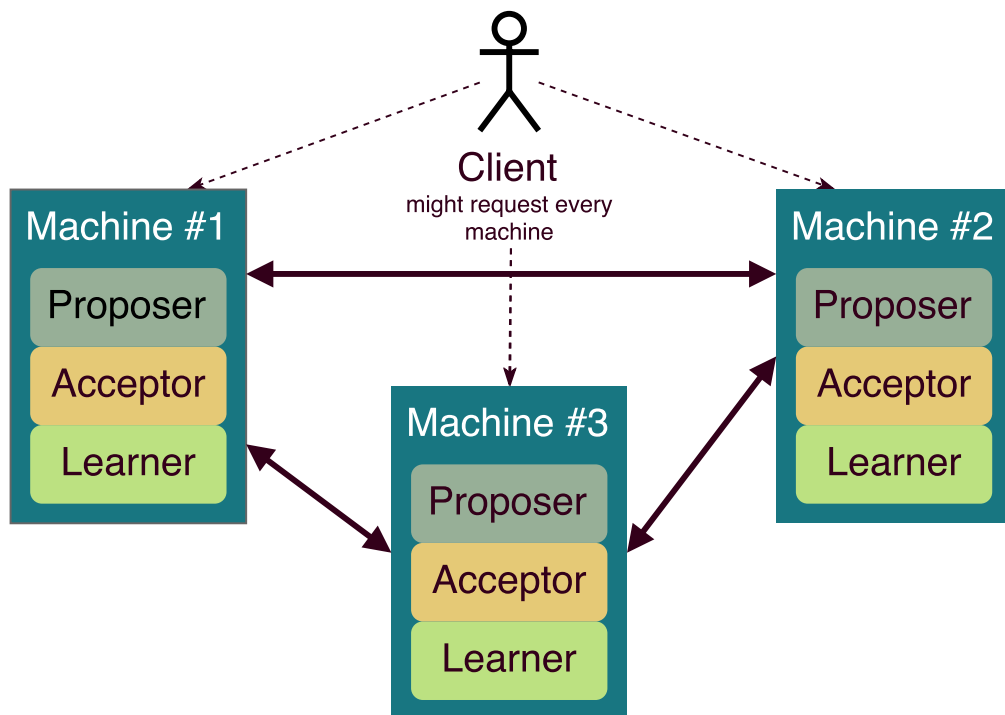


Figure A.1: Example Setup of Three Machines Agreeing on Values Based on the Paxos Algorithm

A.2 Detecting Mutual Inconsistency

Parker et al. claim that a system must ensure the mutual consistency of data copies by applying all changes made to one copy to all others correspondingly [**versionVektor**]. Each time two copies of the same original data item have a different set of modification applied to them, they become incompatible and this conflict must be detected. However, this is not trivial, because “network partitioning can completely destroy mutual consistency in the worst case”. Nevertheless, Parker et al. state that the efficient detection of conflicts that lead to mutual inconsistency can be done by a concept they call *version vectors*. They define a version vector for an item f as “a sequence of n pairs, where n is the number of sites at which f is stored. The i th pair $(S_i: v_i)$ gives the index of the latest version of f made at site S_i ”. An example vector for an item stored at the sites A, B and D is $\langle A:2, B:4, D:3 \rangle$, which translates in a file that has been modified twice on site A, four times on site B, and thrice on site D.

A set of vectors is “compatible when one vector is at least as large as any other vector in every site component for which they have entries”. Otherwise the vectors conflict and are incompatible. E.g., the two vectors $\langle A:2, B:4, D:3 \rangle$ and $\langle A:4, B:5, D:3 \rangle$ are compatible because the second one dominates the other one. $\langle A:3, B:4, D:3 \rangle$ and $\langle A:2, B:5, D:3 \rangle$ conflict, because the first one indicates that the data item was modified one more time on node A, while the second one indicates one more modification occurred on node B. However, if we add a third vector $\langle A:3, B:5, D:4 \rangle$, no conflict exists anymore, because it dominates the two others. The consequences of an operation performed on a data item for a data item’s vector is depicted in table [A.1](#).

By using version vectors, one can detect version conflicts and initiate (automatic) reconciliation. However, Parker et al. warn that two identical updates made on separate partitions will result in a conflict, even though none is present. Thus, they recommend to additionally check two data items on differences before a conflict is raised in certain applications. Furthermore, the

Operation related to data item	Consequence for vector of data item
Update on site S_i	Increment v_i by one
Delete or rename on site S_i	Keep vector and increment v_i by one, remove data item value
Reconcile version conflict	Set each v_i to maximum v_i from all vectors used for reconciliation. In addition, increment v_i of site that initiated reconciliation by one
Copy to new site	Augment vector to include new site

Table A.1: Influence of Operations on a Data Item's Version Vector

reconciliation is most times not trivial, that's why tools such as Cassandra delegate this task to the application layer [cassandra2010].

Appendix B

Acronyms

AWS Amazon Web Services

IoT Internet of Things

Appendix C

Lexicon

Byzantine failure A malfunction of a component that leads to the distribution of wrong/conflicting information to other parts of the system is called Byzantine failure [**lamport1982byzantine**]. The name is based on the Byzantines Generals Problem, in which three Byzantine generals need to agree on a battle plan while one or more of them might be a traitor trying to confuse the others.

Appendix D

Listings

This is the appendix for code, that does not need to be provided directly inside the thesis.

D.1 Configuration for Node A

