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Computational Complexity of some Optimization Problems in Planning

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ABSTRACT

Automated planning is known to be computationally hard in the general case. Propositional planning is $PSPACE$ -complete and first-order planning is undecidable. One method for analyzing the computational complexity of planning is to study restricted subsets of planning instances, with the aim of differentiating instances with varying complexity. We use this methodology for studying the computational complexity of planning.

Finding new tractable (i.e. polynomial-time solvable) problems has been a particularly important goal for researchers in the area. The reason behind this is not only to differentiate between easy and hard planning instances, but also to use polynomial-time solvable instances in order to construct better heuristic functions and improve planners.

We identify a new class of tractable cost-optimal planning instances by restricting the causal graph. We also study the computational complexity of oversubscription planning (such as the net-benefit problem) under various restrictions and reveal strong connections with classical planning. Inspired by this, we present a method for compiling oversubscription planning problems into the ordinary plan existence problem. We further study the parameterized complexity of cost-optimal and net-benefit planning under the same restrictions and show that the choice of numeric domain for the action costs has a great impact on the parameterized complexity.

We finally consider the parameterized complexity of certain problems related to partial-order planning. In some applications, less restricted plans than total-order plans are needed. Therefore, a partial-order plan is being used instead. When dealing with partial-order plans, one important question is how to achieve optimal partial order plans, i.e. having the highest degree of freedom according to some notion of flexibility. We study several optimization problems for partial-order plans, such as finding a minimum deordering or reordering, and finding the minimum parallel execution length.

The research presented in this thesis has been partially funded by the National Graduate School of Computer Science in Sweden (CUGS).

Populärvetenskaplig sammanfattning

Huvudtemat i denna avhandling är studiet av beräkningskomplexitet för automatisk planering. På en hög nivå kan man betrakta planeringsproblemet så här; man utgår från en värld eller ett system, som ett obemannat fordon eller en autonom robot, och antar att det finns vissa handlingar som ändrar tillståndet i systemet. Man antar dessutom att det finns ett initial- och ett måltillstånd. Planering är problemet att hitta en sekvens av handlingar som transformerar initialtillståndet till måltillståndet. En sådan sekvens av handlingar kallas en plan. En planerare är ett datorprogram som löser planeringsproblem, och planerare baseras ofta på sökning genom tillstånden i systemet. Vanligtvis vill man optimera sökningen och hitta en plan som är optimal med avseende på något kriterium. Vad man vill optimera beror på tillämpningen men det är ofta att hitta en kortaste eller billigaste plan.

En viktig fråga när man studerar automatisk planering är hur snabbt man kan hitta lösningar. Ett sätt att analysera denna fråga är att studera beräkningskomplexiteten för olika planeringsproblem. Komplexitetsteori är den gren av datavetenskapen som försöker klassificera hur svåra olika beräkningsproblem är, det vill säga analysera hur mycket resurser (såsom tid och minne) som behövs. Man likställer ofta effektivt lösbara problem med de som kan lösas i *polynomisk tid*: ett problem kan lösas i polynomisk tid om det finns en algoritm vars tidsåtgång begränsas av ett polynom i indatas storlek. Det grundläggande planeringsproblemet är exempelvis känt att vara **PSPACE**-fullständigt. **PSPACE** är den komplexitetsklass som innehåller alla de beräkningsproblem som kan lösas genom att använda polynomiskt mycket minne och de **PSPACE**-fullständiga problemen utgör de svåraste problemen i **PSPACE**. Man tror att det inte finns några **PSPACE**-fullständiga problem som kan lösas i polynomisk tid—detta är ett viktigt olöst problem inom teoretisk datalogi—och man förväntar sig därför inte att det generella planeringsproblemet kan lösas i polynomisk tid.

En metod för att analysera beräkningskomplexitet av planering är att studera begränsade delmängder av planeringsinstanser i syfte att särskilja fall med varierande komplexitet. Vi använder denna metod genomgående i denna avhandling. Så kallad "oversubscription planning" är en typ av planering där man mäter hur nära ett givet måltillstånd

man kan komma. Vi studerar beräkningskomplexiteten för denna typ av planering under olika restriktioner och avslöjar starka kopplingar med vanlig planering. Kostnadsoptimal planering är planering där varje handling har en kostnad och målet är att hitta en plan med minimal kostnad. Dessa handlingskostnader kan hämtas från olika numeriska domäner såsom de positiva heltalen eller de rationella talen. Vi studerar effekter av valet av numeriska domäner och noterar intressanta komplexitetsmässiga skillnader. För kostnadsoptimal planering identifierar vi dessutom en ny klass av instanser som kan lösas i polynomisk tid. Slutligen studerar vi komplexiteten för olika problem inom så kallad partialordningsplanering där man tillåter planer som inte nödvändigtvis är sekvenser av handlingar.

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MEYSAM AGHIGHI
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May 2017

List of Papers

The thesis is based on the following papers:

1. Meysam Aghighi and Peter Jonsson.
Oversubscription planning: Complexity and compilability.
In *Proceedings of the Twenty-Eighth AAAI Conference on Artificial Intelligence, Québec City, Québec, Canada, July 27 -31, pages 2221-2227, 2014.*
2. Meysam Aghighi, Peter Jonsson and Simon Ståhlberg.
Tractable cost-optimal planning over restricted polytree causal graphs.
In *Proceedings of the Twenty-Ninth AAAI Conference on Artificial Intelligence, Austin, Texas, USA, January 25-30, pages 3225-3231, 2015.*
3. Meysam Aghighi and Christer Bäckström.
Cost-optimal and net-benefit planning - A parameterised complexity view.
In *Proceedings of the Twenty-Fourth International Joint Conference on Artificial Intelligence (IJCAI), Buenos Aires, Argentina, July 25-31, pages 1487-1493, 2015.*
4. Meysam Aghighi and Christer Bäckström.
A multi-parameter complexity analysis of cost-optimal and net-benefit planning.
In *Proceedings of the Twenty-Sixth International Conference on Automated Planning and Scheduling (ICAPS), London, UK, June 12-17, pages 2-10, 2016.*
5. Meysam Aghighi and Christer Bäckström.
Plan reordering and parallel execution – a parameterized complexity view.
In *Proceedings of the Thirty-First AAAI Conference on Artificial Intelligence, San Francisco, California, USA, February 4-9, pages 3540-3546, 2017.*

In addition to the above papers, the author has also contributed to the following publications:

1. Meysam Aghighi, Christer Bäckström, Peter Jonsson and Simon Ståhlberg.
Analysing Approximability and Heuristics in Planning Using the

Exponential-time Hypothesis.

In *Proceedings of the Twenty-second European Conference on Artificial Intelligence (ECAI)*, the Hague, Holland, August 29th - September 2nd, pages 184-192, 2016.

2. Meysam Aghighi, Christer Bäckström, Peter Jonsson and Simon Ståhlberg.

Refining Complexity Analyses in Planning by Exploiting the Exponential-time Hypothesis.

In *Annals of Mathematics and Artificial Intelligence (AMAI)*, pages 157-175, 2016.

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Part I

Introduction



1 Thesis Overview

1 About this Thesis

The main topic of this thesis is the study of the computational complexity of automated planning problems. Automated planning has been studied by the artificial intelligence community over the past fifty years. The intersection of complexity theory and planning has been studied since the early work by Chapman [21] which proved that first-order planning is undecidable. Since then, the complexity of planning has attracted a lot of interest.

This thesis is a compilation of five papers. It consists of two parts. The first part contains four chapters. This chapter gives an overview of the thesis. Chapters 2 and 3 give an introduction to computational complexity theory and automated planning, respectively. In Chapter 4, the background of each paper and its contributions are summarized. The second part consists of five peer-reviewed conference papers published between the years 2014 and 2017.

In order to read and understand the contents of the thesis, a basic knowledge in discrete mathematics and graph theory is expected from the reader. Also, familiarity with planning and computational complexity theory is very helpful, in order to understand the contributions of each paper, the motivation of the discussed problems and the intuition behind the proofs. Some good references for complexity theory are Arora and Barak [4] and Garey and Johnson [31], for parameterized complexity Flum and Grohe [30] and Downey and Fellows [27], and for planning Ghallab, Nau, and Traverso [33] and Geffner and Bonet [32].

2 Brief Summary of Papers

This section gives a brief summary of the five papers. More information about the background and contributions of each paper can be found in Chapter 4.

Paper 1

This paper provides a complexity analysis of the oversubscription planning problems, many sub-classes are introduced under the PUBS and Bylander restrictions and a detailed complexity map is presented for them. Moreover, a polynomial-time compilation of net-benefit planning problem to plan existence is presented using a novel counter-based reduction.

Paper 2

This paper presents a new class of tractable cost-optimal planning instances. Our algorithms uses a novel concept of variable isomorphism and shows that every planning instance with: (1) bounded domain size, and (2) polytree causal graph with bounded diameter, is polynomial-time solvable.

Paper 3

This paper gives a parameterized complexity analysis of the cost-optimal and net-benefit planning problems. The two problems are studied with different cost domains for action costs and under Bylander and PUBS restrictions. Our results show that cost-optimal planning for positive integers (\mathbb{Z}_+) is not harder than length-optimal planning while it becomes harder if we use non-negative integers (\mathbb{Z}_0) or positive rationals (\mathbb{Q}_+).

Paper 4

This paper continues paper 3, but this time using a multi-parameter complexity analysis. The paper identifies the parameters that affect complexity the most, such as plan length, maximum denominator of action costs and sum of goal utilities.

Paper 5

This paper continues the work of Bäckström [5] on reordering partial-order plans by applying parameterized complexity theory and introducing parameters like the height and width of a partial-order. The paper analyzes plan optimization using two criteria: order size and parallel length of a plan. It also introduces a new problem: finding the minimum parallel length of a

partial-order plan when there is a bounded number of processors. We show that this problem is significantly harder than the unbounded case.

The papers appear exactly as in the conference proceedings. However, there has been some minor formatting adjustments. The papers of this thesis can be divided into categories from several points of view. From the goal perspective, the first four papers consider cost-optimal and net-benefit planning while paper 5 is on plan optimization. From the complexity analysis point of view, papers 1 and 2 use traditional complexity while papers 3, 4 and 5 use parameterized complexity. From the planning domain perspective, papers 1, 3 and 4 use syntactic restrictions on classical planning, paper 2 uses both syntactic and semantic restrictions and paper 5 goes beyond classical planning by considering partial-order plans.



2

Computational Complexity Theory

1 Computational Complexity

Computational Problems

Computational complexity theory is a branch of computer science that classifies computational problems according to their inherent difficulty. There are several types of computational problems, such as decision, counting, search, optimization, etc. An optimization problem asks for best solution among the set of all possible solutions and a decision problem asks for a yes or no answer, i.e. is there a solution for a given problem instance? Formally, a problem instance and its solution are considered to be strings of bits (i.e. members of the set $\{0, 1\}^*$). Simple encodings can represent general mathematical objects such as integers, graphs, vectors, etc. A **computational problem** is a set of problem instances with a solution for each one of them, i.e. a function $f : \{0, 1\}^* \rightarrow \{0, 1\}^*$. Note that this solution can be an arbitrary mathematical object, like a list of numbers or a set of shortest paths in a graph. A *decision problem* is a special case of computational problem where f is a boolean function $f : \{0, 1\}^* \rightarrow \{0, 1\}$ which is often identified with the set $L_f = \{x : f(x) = 1\}$ that is called a *language*. In complexity theory we are mostly interested in decision problems. We give two concrete examples.

Example I: Satisfiability

Let Φ be a formula in conjunctive normal form (CNF), e.g.

$$\Phi(x_1, x_2, x_3) = (x_1 \vee \neg x_2) \wedge (\neg x_1 \vee x_2 \vee x_3)$$

Φ is satisfiable if there exists a truth assignment to the variables x_1, x_2 and x_3 such that the value of the formula $\Phi(x_1, x_2, x_3)$ is true. For example, $x_1 \mapsto \text{TRUE}, x_2 \mapsto \text{FALSE}$ and $x_3 \mapsto \text{TRUE}$ is a satisfiable assignment for the above formula.

The decision version of SATISFIABILITY (SAT) problem is defined as follows:

INPUT: A CNF formula Φ with variables in $X = \{x_1, x_2, \dots, x_n\}$.

QUESTION: Is there a satisfying truth assignment for Φ ?

A truth assignment can be viewed as a function $\alpha : X \rightarrow \{\text{TRUE}, \text{FALSE}\}$ and the decision problem as a function $f : \{0, 1\}^* \rightarrow \{0, 1\}$. f takes a string in the input that represents a CNF formula and the set X . Then it returns 1 if the formula is satisfiable, otherwise it returns 0. L_f is the set of all strings on which f returns 1, i.e. the set of all satisfiable CNF formulas.

Next, we present an example based on an optimization problem:

Example II: Traveling Salesperson Problem

The optimization version of TRAVELING SALESPERSON PROBLEM (TSP) is defined as follows:

INPUT: A graph $G = (V, E)$ and a function $c : E \rightarrow \mathbb{N}$.

QUESTION: Find the minimum value $k \in \mathbb{N}$ for which a Hamiltonian cycle of cost k exists, i.e. there is a permutation (v_1, v_2, \dots, v_n) of vertices in V such that $(v_n, v_1) \in E$, for every $1 \leq i \leq n - 1, (v_i, v_{i+1}) \in E$ and

$$\left(\sum_{i=1}^{n-1} c(v_i, v_{i+1}) \right) + c(v_n, v_1) \leq k$$

The optimization problem can be viewed as a function $f : \{0, 1\}^* \rightarrow \{0, 1\}^*$. f takes a string in the input that represents a TSP instance, which is a graph and a cost function for the edges. Then it returns a string in the output that represents the minimum integer k for which a Hamiltonian cycle of cost k exists, or a special designated character that means no Hamiltonian cycles exist in the graph.

Model of Computation

In order to measure the complexity of computational problems and analyze the resources required for algorithms and computers, we need to define a model for computation. There are several models for computation such as the random-access machine, the finite-state machine, the Turing machine, the logic circuit, etc. The Turing machine was proposed by Alan Turing in 1936 and is the most commonly used model in computation. This is because of the Church-Turing thesis; a hypothesis stating that every "reasonable" model of

computation can be simulated by a Turing machine with only a polynomial slowdown in run-time in terms of the input size. Please note that this was not the original Church-Turing thesis, but a result of the extended Church-Turing hypothesis [4]. It is also believed that for any computer algorithm, a Turing machine can be constructed to simulate the algorithm's logic [70].

A **Turing machine** consists of an infinite memory tape divided into cells, a set of symbols and a finite table of instructions. The machine's head is positioned over a cell and reads the symbol there. Then, based on the read symbol and the current position of the machine in the table of instructions, the head writes a symbol in the cell and then moves either one cell to the left or one cell to the right. Then, it either proceeds to the next instruction or halts the computation.

Formally, a deterministic (one-tape) Turing machine M is described by a 7-tuple $(Q, \Gamma, b, \Sigma, \delta, q_0, F)$ where [45]

- Q is a finite, non-empty set of states that M can be in, including a designated initial state (q_0) and a set of final or accepting states ($F \subseteq Q$).
- Γ is a finite, non-empty set of symbols that M 's tape can contain, including a designated blank symbol b (the only symbol allowed to occur on the tape initially often at any step during the computation). Γ is called the *alphabet* of M .
- $\Sigma \subseteq \Gamma \setminus \{b\}$ is the set of input symbols.
- $\delta : (Q \setminus F) \times \Gamma \rightarrow Q \times \Gamma \times \{L, R\}$ is a function describing the instruction that M uses in every step. L and R denote the left and right movement by the machine's head. δ is called the *transition function* and denotes the table of instructions. If δ is not defined in the current state and the current tape symbol, then the machine halts.

A Turing machine accepts the initial tape contents, if it eventually halts in a state from F .

There are several variations of the Turing machine. The two most common are *deterministic* Turing machine (DTM) and *non-deterministic* Turing machine (NTM). In a non-deterministic Turing machine, the instructions' table allows more than one action in each situation and the transition function is replaced by a transition relation. For instance δ may contain both $((q, a), (q_1, b, L))$ and $((q, a), (q_2, c, R))$. This is to be interpreted as: if the machine is in state q and its head reads the symbol a , it has the choice of going to state q_1 , write symbol b and move left, or go to state q_2 , write symbol c and move right. An NTM accepts an input if there exists some choice of actions that leads to the designated halting state. These variations are in principle equally powerful but when the resources (such as time and space) are restricted this equality breaks down.

Complexity Measures

The two most common complexity measures are time and space. The running time of an algorithm, intuitively, is the number of basic operations performed by it, where a basic operation takes a fixed amount of time to perform. The run-time of algorithm A on input x is formally defined as the number of state transitions of the corresponding deterministic Turing machine for the algorithm, on input x . The time complexity of an algorithm is the run-time as a function of the input size (number of bits needed to represent the input). Similar to time, the space complexity is the amount of working storage an algorithm needs. The complexity is often reported *asymptotically*, as the input size goes to infinity. For instance, if an algorithm takes $3n^5 + 4n^2 \log n - 2n$ units of time for an input of size n , we say that the time complexity of the algorithm is $O(n^5)$, using the big O notation. This is because for large values of n , the value other terms are insignificant compared to the term $3n^5$. Since the time complexity of an algorithm may vary with different inputs of the same size, we usually consider the complexity as the maximum value over all inputs of size n . This is referred to as the *worst-case* complexity analysis.

Reduction and NP-completeness

Reduction is a way to show that one problem is at least as hard as another problem. Intuitively, problem A is reducible to problem B , if an algorithm that solves problem B , can be modified to solve problem A . There are several types of reductions and the most common form is the polynomial-time many-one (Karp) reduction which is usually referred to as the polynomial-time reduction.

Formally, we say that a language $A \subseteq \{0,1\}^*$ (equivalently, a decision problem) is polynomial-time reducible to a language $B \subseteq \{0,1\}^*$ denoted by $A \leq_p B$ if there exists a polynomial-time computable function $f : \{0,1\}^* \rightarrow \{0,1\}^*$ such that for every $x \in \{0,1\}^*$, $x \in A$ if and only if $f(x) \in B$.

A complexity class is a set of computational problems that can be computed within a given resource. For instance, $\mathbf{DTIME}(f(n))$ is the class of all decision problems that can be solved by some DTM running in time $O(f(n))$ where n is the input size and f is some function. The two most well-known complexity classes are \mathbf{P} , the class of problems solvable in polynomial time by a DTM, and \mathbf{NP} , the class of problems solvable in polynomial time by an NTM. Informally, a decision problem belongs to the class of \mathbf{NP} if and only if its instances for which the answer is “yes” have efficiently (polynomial-time) verifiable proofs. The two definitions are equivalent because the first part in both contains a *guess* about the solution, which is generated in a non-deterministic way, while the second part contains a deterministic algorithm that verifies or rejects the guess as a valid solution to the problem.

There are many other complexity classes such as $\mathbf{EXPTIME}$, the class of problems solvable in exponential time by a DTM, \mathbf{PSPACE} , the class of prob-

lems solvable in polynomial space by a DTM, **NL**, the class of problems solvable in logarithmic space by an NTM, **#P**, the class of counting problems associated with decision problems in **NP**, etc. The complexity classes are usually distinguished by: (1) The type of the computational problem (decision, counting, etc.), (2) The model of computation (DTM, NTM, Boolean circuits, etc.), and (3) The resource constraints (polynomial time, logarithmic space, etc.).

Let **C** be a complexity class. Problem *A* is **C**-hard if there is a polynomial-time reduction from every member of **C** to *A*. This means if there is an efficient algorithm for *A*, then this algorithm can also solve all problems in **C** with only a polynomial blowup in time. Moreover, *A* is said to be **C**-complete if it is **C**-hard and additionally in **C**. One may say that the **C**-complete problems are, intuitively, the hardest problems in **C**. Both SAT and TSP decision problems from Examples I and II are proved to be **NP**-complete [31, LO1, GT37]. This results in the optimization of TSP to be **NP**-hard, because a solution to the optimization version of TSP will immediately solve its decision version.

The **P** vs **NP** problem ($\mathbf{P} \stackrel{?}{=} \mathbf{NP}$) is a major open problem in computer science with a million dollar prize by the Clay Mathematics Institute. It asks if every problem whose solution can be verified in polynomial time by a computer can also be solved in polynomial time by a computer. Most researchers in the area believe that the answer to this question is negative (i.e. $\mathbf{P} \neq \mathbf{NP}$) and this would imply that many common problems cannot be solved efficiently. However, a positive answer would have enormous practical consequences; for instance, many cryptography algorithms rely on the difficulty of certain problems [60].

2 Parameterized Complexity

In traditional complexity analysis, the time complexity is measured only in terms of the input size. However, many problem instances consist of two or more parameters that contribute differently to the complexity of the problem.

Example: Vertex Cover vs Dominating Set

The problems VERTEX COVER and DOMINATING SET are defined as follows:

VERTEX COVER

INPUT: A graph $G = (V, E)$ and a positive integer k .

QUESTION: Does G have a vertex cover of size at most k , i.e. is there a set of vertices $V' \subseteq V$ such that $|V'| \leq k$ and for every edge $(u, v) \in E$, either $u \in V'$ or $v \in V'$?

DOMINATING SET

INPUT: A graph $G = (V, E)$ and a positive integer k .

QUESTION: Does G have a dominating set of size at most k , i.e.

is there a set of vertices $V' \subseteq V$ such that $|V'| \leq k$ and for every vertex $u \in V \setminus V'$, $(u, v) \in E$ for some $v \in V'$?

Although both problems are **NP**-complete, the input parameter k contributes differently to the complexity of the two problems. Let n be the number of vertices. The best known algorithm for VERTEX COVER solves the problem in time $O(1.2738^k + kn)$ [22], while the best known algorithm for DOMINATING SET is the brute force algorithm of trying all k -subsets which takes $O(n^{k+1})$ time. This is a significant difference; the ratio of n^{k+1} to $1.2738^k + kn$ is more than 10^{16} for $n = 150$ and $k = 10$. \square

The observation of such differences between **NP**-complete problems motivated a new and more fine-grained form of complexity analysis: parameterized complexity analysis. Parameterized complexity tries to find how different parameters of a problem contribute to its hardness. For instance, parameter k in VERTEX COVER usually has a much smaller value compared to n in practice. Therefore, even an exponential algorithm that runs in time $O(1.2738^k + kn)$ performs successfully on many real world examples.

A *parameterized problem* is a language $L \subseteq \Sigma^* \times \Sigma^*$ over some finite alphabet Σ . The *instances* of L are tuples $\langle \mathbb{I}, k \rangle$, where k is called the *parameter*. The parameter is often a non-negative integer, but it can be anything, e.g. a rational number, three integers or a graph. For simplicity, we first assume the parameter is a non-negative integer, i.e. $L \subseteq \Sigma^* \times \mathbb{Z}_0$. A parameterized problem is *fixed-parameter tractable (fpt)* if there exists an algorithm that solves every instance $\langle \mathbb{I}, k \rangle$ of size $n = |\mathbb{I}|$ in time $f(k) \cdot n^c$ where f is an arbitrary computable function and c is a constant independent of both n and k . **FPT** is the class of all fixed-parameter tractable decision problems. In contrast to classical tractability, some exponentiality is allowed, but confined to the parameter only, thus better reflecting reality. For instance, the VERTEX COVER problem parameterized by the size of the vertex cover, k , is in **FPT**. Note that $1.2738^k + kn \leq 1.2738^k n + kn = (1.2738^k + k)n = f(k)n$.

Hardness for parameterized classes is proven in the usual way, but using *fpt* reductions instead of ordinary polynomial-time reductions. An *fpt reduction* from a parameterized language $L \subseteq \Sigma^* \times \mathbb{Z}_0$ to another parameterized language $L' \subseteq \Sigma^* \times \mathbb{Z}_0$ is a mapping $R : \Sigma^* \times \mathbb{Z}_0 \rightarrow \Sigma^* \times \mathbb{Z}_0$ such that: (1) $\langle \mathbb{I}, k \rangle \in L$ if and only if $\langle \mathbb{I}', k' \rangle = R(\mathbb{I}, k) \in L'$; (2) there is a computable function f and a constant c such that R can be computed in time $f(k) \cdot n^c$, where $n = |\mathbb{I}|$; and (3) there is a computable function g such that $k' \leq g(k)$.

Parameterized complexity offers a completeness theory, similar to the theory of **NP**-completeness. It contains the following hierarchy of parameterized complexity classes

$$\mathbf{FPT} \subseteq \mathbf{W}[1] \subseteq \mathbf{W}[2] \subseteq \mathbf{W}[3] \subseteq \dots \subseteq \mathbf{W}[P],$$

known as the *W hierarchy*. $\mathbf{W}[i]$ is defined using WEIGHTED CIRCUIT SATISFIABILITY problem: a parameterized problem is in $\mathbf{W}[i]$ if its every instance $\langle \mathbb{I}, k \rangle$ is *fpt*-reducible to a boolean circuit of weft i , such that $\langle \mathbb{I}, k \rangle$ has a solution if and only if the circuit has a truth assignment of weight (number

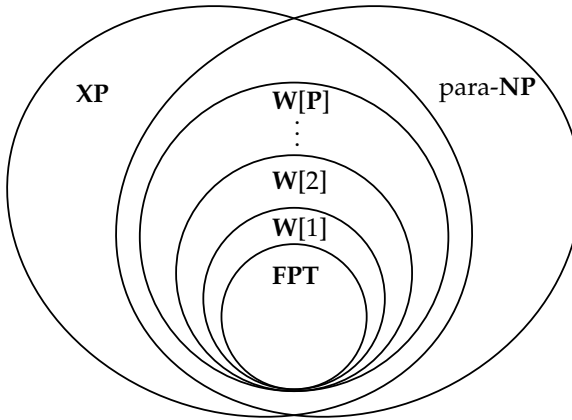


Figure 2.1: Parameterized Complexity Classes

of 1s in the input) at most k . Weft is the maximum number of gates with fan-in more than two on any path from the inputs to the output. The class $W[P]$ corresponds to arbitrary circuits. For instance, the DOMINATING SET problem parameterized by the size of the dominating set is $W[2]$ -complete [25]. There are also classes outside of the W -hierarchy; para-NP is the class of problems solvable in *non*-deterministic time $f(k)n^c$ and XP is the class of problems solvable in time $n^{f(k)}$; where f is an arbitrary computable function and c is a constant independent of both n and k . It is known that $W[P]$ is a subset of both para-NP and XP [30]. Figure 2.1 shows the known relation between the mentioned parameterized complexity classes.

It is known that $P \subseteq \text{FPT}$ and $\text{NP} \subseteq \text{para-NP}$, but otherwise there is not much known about the relations between parameterized complexity classes and the classical ones. For instance, there are NP -complete problems that are $W[P]$ -complete and there are PSPACE -complete problems that are in FPT . The hardness of a parameterized problem depends a lot on the chosen parameter(s). For instance, consider the NP -complete problem GRAPH k -COLORING: this problem is para-NP -hard if parameterized by k , the number of available colors (mentioned in Paper 5 of this thesis) while it is in FPT if parameterized by n , the number of vertices (using a brute-force algorithm).



3

Automated Planning

1 What is Planning?

Planning is the art and practice of thinking before acting.

Patrik Haslum

In this section we give a high-level description of planning. Imagine a world or a system, like a car, a robotic spacecraft or a classroom with all the students and the teacher in it. There are some actions that change the state of the world. By executing the actions we can move between different states of the world. For instance, entering and leaving a student, giving a lecture or taking a break are all actions that change the state of the classroom. Planning may be defined as the problem of finding a sequence of such actions that lead to a specified goal state from a given initial state. A more abstract definition is given by Ghallab, Nau, and Traverso [34]:

Planning is the reasoning side of acting. It is an abstract, explicit deliberation process that chooses and organizes actions by anticipating their expected outcomes. This deliberation aims at achieving as best as possible some pre-stated objectives. Automated planning is an area of Artificial Intelligence (AI) that studies this deliberation process computationally.

We will continue with an example, Blocks world, which is one of the most famous planning domains in Artificial Intelligence.

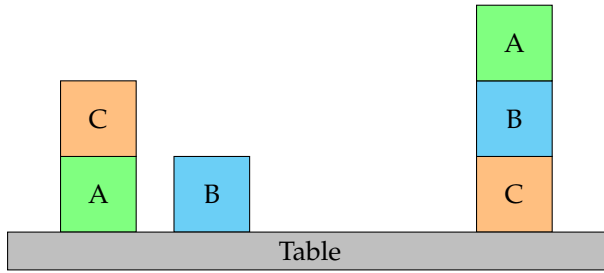


Figure 3.1: A Blocks world example; initial state (left) and goal state (right)

Example: Blocks world

In the Blocks world domain, we are given a set of identical blocks with different letter identifiers on them. The blocks are positioned in several stacks on a table. We have an arm to move the blocks. This arm can only pick blocks from the top of the stacks and put them on top of another stack or on the table. The goal is to reach another specified configuration. This version of Blocks world is known as the Elementary Blocks World (EBW). There are also other variations of this problem [38].

To define the state of the world, we use the following predicate types:

- **ON(A,B)**: Block A is on block B.
- **ONTABLE(A)**: Block A is on the table.
- **CLEAR(A)**: Block A has nothing on it.
- **HOLDING(A)**: The arm holds block A.
- **ARMEMPTY**: The arm holds nothing.

We also define the following actions:

- **UNSTACK(A,B)**: Pick up clear block A from block B.
- **STACK(A,B)**: Place block A using the arm onto clear block B.
- **PICKUP(A)**: Lift clear block A with the empty arm from the table.
- **PUTDOWN(A)**: Place the held block A onto a free space on the table.

Figure 3.1 shows an example with three blocks. In this example, the initial (*I*) and goal (*G*) states contain the following predicates

$$I = \{\text{CLEAR}(C), \text{CLEAR}(B), \text{ONTABLE}(A), \text{ONTABLE}(B), \text{ON}(C,A), \text{ARMEMPTY}\}$$
$$G = \{\text{CLEAR}(A), \text{ONTABLE}(C), \text{ON}(A,B), \text{ON}(B,C), \text{ARMEMPTY}\}$$

UNSTACK(A,B):	
<i>precondition list</i>	= {ON(A,B), CLEAR(A), ARMEMPTY}
<i>add list</i>	= {CLEAR(B), HOLDING(A)}
<i>delete list</i>	= {ON(A,B), ARMEMPTY}
STACK(A,B):	
<i>precondition list</i>	= {HOLDING(A), CLEAR(B)}
<i>add list</i>	= {ARMEMPTY, ON(A,B)}
<i>delete list</i>	= {HOLDING(A), CLEAR(B)}
PICKUP(A):	
<i>precondition list</i>	= {CLEAR(A), ONTABLE(A), ARMEMPTY}
<i>add list</i>	= {HOLDING(A)}
<i>delete list</i>	= {ONTABLE(A), ARMEMPTY}
PUTDOWN(A):	
<i>precondition list</i>	= {HOLDING(A)}
<i>add list</i>	= {ONTABLE(A), ARMEMPTY}
<i>delete list</i>	= {HOLDING(A)}

Figure 3.2: The actions of the Blocks world domain

Every action is defined by three lists of predicates; a *precondition list*, the list of predicates that need to hold for the action to be applicable in a state, an *add list*, the list of predicates that are added after the action is executed in a state, and a *delete list*, the list of predicates that are deleted after the execution of the action. The actions are defined in Figure 3.2.

The objective is to find the shortest sequence of actions (a plan) that bring the initial state (I) to the goal state (G). In this example, the unique optimal plan, ω , is the following list of actions:

$$\omega = (\text{UNSTACK}(C,A), \text{PUTDOWN}(C), \text{PICKUP}(B), \\ \text{STACK}(B,C), \text{PICKUP}(A), \text{STACK}(A,B))$$

□

Given a Blocks world example with n blocks, one can produce in time $O(n)$ a plan (if one exists) that moves no block more than twice. However, the problem of finding the optimal plan is NP-hard [38].

Planning is closely connected to search. The state space of the blocks world example can be seen as a directed graph where states (combination of predicates) are the nodes and there is an edge from node u to node v if there is an action that transforms state u to state v . Hence, the problem of finding an optimal plan becomes the problem of finding the shortest path between two nodes. Although this may seem easy using shortest path algorithms, the main challenge is that the number of nodes of this graph is exponential in

the input size. For instance, a blocks world instance with n blocks has more than 2^{n^2} states, which means even for $n = 10$ and checking 2 billion states per second, a linear search takes more than 10 billion years. However, the problem can easily be solved by a human in a few seconds.

2 Planning Model and Formalism

There are different ways to model the planning world. The differences come from how the following questions (motivated by Ghallab, Nau, and Traverso [34]) are answered:

- Is the planning process offline (static) or is it online (dynamic)?
- Is the state of the plan fully, partially or not observable?
- Is there only a single agent or are there several agents to execute actions?
- Do variables have discrete or continuous domain? Do they have finite or infinite domain?
- Are the actions deterministic, non-deterministic or stochastic?
- Do the actions have cost or duration?
- Can the actions be executed at the same time or do we only have sequential plans?
- Is the objective to satisfy a goal or to find the optimal plan under some criteria?

The most simple planning model is called **classical planning**. In classical planning the following assumptions are imposed:

- planning is static
- variable domains are discrete and finite
- actions are deterministic
- states are fully observable
- there is a single agent (sequential plan)
- actions have no cost or duration

There are many models and representations for planning in the literature [34]. Two common formalisms that are used to study the complexity of planning are STRIPS and SAS⁺. This is because they are simple to work with and at the same time expressive enough to be able to represent many complex planning domains.

STRIPS

STRIPS (Stanford Research Institute Problem Solver) was defined by Fikes and Nilsson [29]. It is a planner and a formalism for describing planning instance that is based on first-order logic: the state of the world is viewed as a set of ground and function-free first-order literals. We will not go into the details, but merely note that the Blocks world example in this chapter is an example of a STRIPS planning instance. In many cases, one is content with the simplified variant of STRIPS where the state of the world is viewed as a set of propositional literals. This variant is referred to as propositional STRIPS.

SAS⁺

SAS⁺ (Simplified Action Structures) is very similar to propositional STRIPS, with the difference that the two-valued propositions are replaced by variables that take their values from a finite domain. It was introduced by Bäckström and Klein [10]. Another difference is that instead of predicates the domain values are assignments to variables in preconditions, effects, initial and goal state.

We have mostly used SAS⁺ formalism in the papers of this thesis and will formally define it here (Although depending on the application, SAS⁺ is defined in slightly different ways in the papers of this Thesis.). A SAS⁺ planning instance is a 4-tuple $\langle V, A, I, G \rangle$, where

- $V = \{v_1, v_2, \dots, v_n\}$ is a set of variables over some finite domain \mathcal{D} (in some definitions each variable v has an individual domain $D(v)$). Define $\mathcal{D}_+ = \mathcal{D} \cup \{\mathbf{u}\}$, where \mathbf{u} is an extra value that stands for undefined. \mathcal{D}^n is the set of *total states* where \mathcal{D}_+^n is the set of *partial states*. The value of variable v in state s (partial or total) is denoted by $s[v]$.
- A is a set of actions and each action $a \in A$ is of the form $a : \text{pre}(a) \rightarrow \text{eff}(a)$, where $\text{pre}(a), \text{eff}(a) \in \mathcal{D}_+^n$ and are called precondition and effects (or post-conditions), respectively. An action a is applicable in total state s if and only if every value with a defined value ($\neq \mathbf{u}$) in $\text{pre}(a)$ has the same value in s . If a is applicable in s , the result of a in s is a total state t where $t[v] = s[v]$ if $\text{eff}(a)[v]$ is undefined and $t[v] = \text{eff}(a)[v]$ otherwise.
- $I \in \mathcal{D}^n$ is the initial state.
- $G \in \mathcal{D}_+^n$ is the goal state.

For two total states s_I and s_G , a sequence of actions (a_1, a_2, \dots, a_k) is called a *plan* from s_I to s_G if and only if there exist total states $\{s_0, s_1, \dots, s_k\}$ such that $s_0 = s_I, s_k = s_G, a_i$ is applicable in s_{i-1} and s_i is the result of a_i in s_{i-1} . A plan from I to a total state s is a *solution plan* if $s[v] = G[v]$ for every defined variable in G .

3 Planning Problems

There are many planning problems described in the literature, with respect to different planning models, formalisms, constraints and objectives. In this thesis we investigate the following problems:

PSAT Plan satisfiability problem (also known as plan existence problem (PE)) is the most basic planning problem. It asks if a planning instance (e.g. a SAS⁺ instance) has a solution.

LOP Length-optimal planning takes an integer ℓ and asks if the planning instance has a solution of length at most ℓ .

COP Cost-optimal planning is a generalization of LOP. It assigns a cost to each action ($c : A \rightarrow \mathbb{D}$), takes a cost threshold w and asks if there is plan of cost at most w .

PSP Partial-satisfaction problem is a generalization of PSAT. It takes an integer k in the input and asks if there is a plan that satisfies at least k of the goals.

NBP Net-benefit planning assigns a utility to each goal ($u : G \rightarrow \mathbb{D}$) and a cost to each action. It takes a net-benefit threshold b and asks if there is a plan with net-benefit at least b . The net-benefit of a plan is the sum of utilities of the satisfied goals minus the plan cost.

We also investigate some problems beyond the boundaries of classical planning. To continue, we first need to define *partial-order* and *parallel* plans. For a more detailed explanation please refer to Paper 5.

Let Π be a planning problem instance. A partial-order plan for Π is a tuple $P = \langle A, < \rangle$, where A is a set of action occurrences and $<$ is a partial order on A . P is a valid plan for Π if all linearizations of it are valid plans. The size of an order is the number of relation tuples in it. A parallel plan is a tuple $P = \langle A, <, \# \rangle$ where $\langle A, < \rangle$ is a partial-order plan and $\#$ is an irreflexive and symmetric relation on A . The relation $\#$ is called the *non-concurrency* relation. For every two actions a and b , $a\#b$ indicates that they cannot be executed in parallel. Let $d : A \mapsto \mathbb{N}$ denote the duration of each action. A *parallel execution* of a parallel plan P is a function $r : A \mapsto \mathbb{N}$, denoting the release time for every action in A , such that: 1) If $a < b$ then $r(a) + d(a) \leq r(b)$, and 2) If $a\#b$ then either $r(a) + d(a) \leq r(b)$ or $r(b) + d(b) \leq r(a)$. The length of a parallel execution or $\max_{a \in A} \{r(a) + d(a)\}$ denotes the latest finishing time of any action. A *k-processor parallel execution* of a parallel plan is a parallel execution such that for each time $t \in \mathbb{N}$ there are at most k actions a such that $r(a) \leq t < r(a) + d(a)$.

PPL Parallel plan length takes a parallel plan $P = \langle A, <, \# \rangle$, a length threshold ℓ and asks if there is a parallel execution for P of length at most ℓ .

PPL_k This problem takes a parallel plan $P = \langle A, <, \# \rangle$, an integer k , a length threshold ℓ and asks if there is a k -processor parallel execution for P of length at most ℓ .

MCR Minimum constrained reordering (MCR), takes a partial-order plan, a value k and asks if there is another valid order for the plan with size k .

MCD Minimum constrained deordering (MCD) is similar to MCR but requires the new order to be a subset of the initial order.

MPR Minimum parallel reordering asks if a partial-order plan has a reordering of parallel length ℓ .

MPD Minimum parallel deordering is the deordering version of MPR.

Complexity of Planning

Planning problems are known to be computationally hard in the general case. PSAT for propositional STRIPS and SAS⁺ is PSPACE-complete [11, 18] and for STRIPS with predicates over infinite domains is undecidable [21], i.e. it is impossible to construct an algorithm that always generates a correct “yes” or “no” answer.

One method to reduce the computational complexity of planning is to apply restrictions on the planning instance. The complexity of STRIPS planning varies from constant time to undecidable depending on which restrictions we make. These restrictions can be syntactic restrictions, such as restrictions on the number of preconditions and effects [2, 9, 18], or restrictions on action types [7, 10, 11, 36, 49]; semantic restrictions, such as restrictions on the variable dependencies [42, 43, 46, 48, 49].

Finding new tractable problems in planning has received a special attention not only because it helps to distinguish between easy and hard planning instances, but also because one can use the polynomial-time solvable instances in order to construct better heuristic functions and improve planners [42, 43]. Many tractable classes of planning problems have been identified [3, 11, 15, 36, 49, 52].

A recent method to tackle the hardness of planning problems is to get a more fine-grained complexity analysis by using parameterized complexity. Downey, Fellows, and Stege [26] showed that STRIPS planning is W[1]-hard parameterized by the plan length (Later, Bäckström, Jonsson, Ordyniak, and Szeider [9] strengthened this result and showed that LOP parameterized by the plan length is W[2]-complete). Since then, the parameterized complexity of PSAT, LOP, COP and NBP has been studied under various restrictions and different parameterizations [1, 6, 9, 12, 39, 56].

There are also other complexity results on variants of planning that are out of the scope of this thesis [13, 19, 28, 37, 41, 57, 58, 67].



4

Background and Contributions

This section highlights the background and contributions of each paper in this thesis.

1 Paper 1: Oversubscription Planning

Meysam Aghighi and Peter Jonsson

Oversubscription planning: Complexity and compilability

In Proceedings of the Twenty-Eighth AAAI Conference on Artificial Intelligence, Québec City, Québec, Canada, July 27 -31, 2014.

In classical planning all goals need to be satisfied at the same time. However, in many real-world applications there are a larger number of possible goals and the aim is to find a feasible subset of them [71]. This type of planning is called partial satisfaction or oversubscription planning. Several custom-made planners and heuristics have been suggested for oversubscription planning [14, 17, 61].

In this paper we study two different problems: the partial satisfaction problem (PSP) where the goal is to maximize the number of achieved goals, and the net-benefit problem (NBP) where the goal is to maximize the total weight of the achieved goals minus the plan cost. The net-benefit problem was first formalized by Briel, Nigenda, Do, and Kambhampati [17] and was shown to be \mathbf{PSPACE} -complete. After that, there has not been a more detailed complexity analysis of this problem until paper 1 of this thesis. We consider two types of restrictions in our complexity analysis: (1) restrictions on the

number of preconditions and effects that were first proposed by Bylander [18], and (2) the PUBS restrictions proposed by Bäckström and Klein [10].

The results indicate that there is a high correlation between PSAT, COP, PSP and NBP. For instance, if PSAT for some set of instances X (under mild additional assumptions) is NP-complete, then PSP for X is NP-complete, too. We also propose a polynomial-time reduction from NBP into PSAT, using a novel counter to keep track of plan cost during the plan execution and calculate the net-benefit after plan execution. The importance of this compilation is that the number of variables increase slowly in the size of the original instance, and this means that the algorithms for PSAT are able to solve NBP instances with a limited slow-down. Compiling different planning problems into each other [55, 64, 72] and into other problems [16, 20, 54] have been active research areas for quite some time.

At the end of the paper we suggested a continuation of the work by imposing restrictions on the causal graph. Later, Katz and Mirkis [53] discovered several tractable fragments of net-benefit planning by restricting the causal graph.

2 Paper 2: Causal Graphs for Tractability

Meysam Aghighi, Peter Jonsson, and Simon Ståhlberg

Tractable cost-optimal planning over restricted polytree causal graphs

In Proceedings of the Twenty-Ninth AAAI Conference on Artificial Intelligence, Austin, Texas, USA, January 25-30, 2015.

A causal graph is a directed graph with the domain variables as nodes. Node x in the graph is connected to node y if there exists some action a where $\text{eff}(a)[y] \neq \mathbf{u}$ and either $\text{eff}(a)[x] \neq \mathbf{u}$ or $\text{pre}(a)[x] \neq \mathbf{u}$. There has been many studies on the causal graph of a planning problem [15, 36, 51]. Giménez and Jonsson [35] showed that plan generation is NP-hard for instances whose causal graph is a directed path and have domain size at least 5. Katz and Domshlak [48, 50] have studied planning instances whose causal graphs are forks and inverted-forks. Katz and Keyder [52] have shown that cost-optimal planning instances with bounded domain size and hourglass causal graphs are in P. In this paper, we generalize this result and prove that cost-optimal planning instances with bounded domain size and polytree causal graphs with bounded diameter, are in P.

The algorithm in this paper is built on a novel notion of isomorphic variables. Intuitively, two variables are isomorphic if the structure of their domain transition graphs are isomorphic with regard to the initial value, goal value and actions. This results in the same behavior of these variables in every optimal solution, which allows us to combine them into a single variable. By repeating this process, we gradually reduce the size of the initial instance and then solve it in polynomial time.

Brafman and Domshlak [15] have presented a polynomial-time algorithm for instances with binary variables and where the causal graph is a

polytree such that the in-degree of every vertex is bounded by a constant. Giménez and Jonsson [36] later generalized this result by giving an algorithm that, instead of restricting the in-degree, restricts the number of prevail-conditions (a bound on the in-degree implies a bound on the number of prevail-conditions but not the other way round). Our algorithm works differently: it allows not only binary domains, but any domain size bounded by a constant and it allows the actions to have an unbounded number of prevail-conditions. However, this flexibility comes at the cost of bounding the diameter by a constant.

3 Papers 3, 4: The Effect of Cost Domain in Cost-optimal and Net-benefit Planning

Meysam Aghighi and Christer Bäckström

Cost-optimal and net-benefit planning - A parameterised complexity view

In Proceedings of the Twenty-Fourth International Joint Conference on Artificial Intelligence, IJCAI 2015, Buenos Aires, Argentina, July 25-31, 2015.

Meysam Aghighi and Christer Bäckström

A multi-parameter complexity analysis of cost-optimal and net-benefit planning

In Proceedings of the Twenty-Sixth International Conference on Automated Planning and Scheduling, ICAPS 2016, London, UK, June 12-17, 2016.

Although cost-optimal and net-benefit planning have been vastly studied in the literature, very little attention has been paid to motivate and discuss the choice of numeric domain for the action costs. For instance, Katz and Domshlak [49] and Helmert, Haslum, Hoffmann, and Nissim [44] use non-negative reals; Bäckström and Jonsson [8] and Yang, Culberson, Holte, Zahavi, and Felner [77] use non-negative integers; Cooper, Roquemaurel, and Régnier [24] use positive integers; Coles, Fox, Long, and Smith [23] use non-negative costs with no specified type; and Thayer, Stern, Felner, and Ruml [73] disregard cost types.

A traditional computational complexity analysis does not differentiate between different cost types; planning remains **PSPACE**-complete regardless of whether action costs are positive or non-negative, integer or rational. In paper 3, we use parameterized complexity to draw a line between different cost domains: cost-optimal planning is **W[2]**-complete for positive integer (\mathbb{Z}_+) costs while it is para-**NP**-hard if the costs are non-negative integers (\mathbb{Z}_0) or positive rationals (\mathbb{Q}_+). This is a strong indication that the latter cases are substantially harder. The paper continues by providing a detailed complexity map of both cost-optimal and net-benefit planning problems, with **PUBS** restrictions and restrictions on the number of preconditions and effects.

This result may seem in contradiction with the usual assumption that the use of rational costs is the same as integer costs: by multiplying all action costs and plan cost threshold by a sufficiently large number. By using

this transformation, from any rational-cost planning instance we can get an integer-cost instance and solve it by using an integer-cost planner. Although this method is correct and the new instance with integer costs has a solution if and only if the initial rational-cost instance has a solution, the transformation does not preserve the parameterized complexity (it is not an fpt-reduction).

The work is continued by a multi-parameter complexity analysis of the same problems in paper 4. In this work we consider a total of thirteen parameters including the plan length, ℓ , the plan cost, k , the largest denominator of the action costs, d , and the sum of variable utilities, t . The latter parameter only applies to NBP. Two new cost domains were also considered in this paper, the non-negative rationals (\mathbb{Q}_0) and rationals greater than or equal to one (\mathbb{Q}_1). Our results show that COP remains **W[2]**-hard for all cost domains even when combining all parameters. We also show the **W[2]**-membership of COP for the following parameter combinations and domains:

- $\{k, \ell\}$ for \mathbb{Z}_0
- $\{k, d\}$ for \mathbb{Q}_+
- $\{k, d, \ell\}$ for \mathbb{Q}_0

Another interesting and important result of paper 4 is the effect of parameter t , the sum of all utilities. We show, through an fpt reduction, that NBP parameterized by t is reducible to COP parameterized by k over the same numeric domain. This reduction is both related to the method used by Keyder and Geffner [55] to compile away soft goals, and to the compilation of NBP to PSAT by Aghighi and Jonsson [2] in paper 1 of this thesis. However, if we use the second method and apply the counter reduction from paper 1, we have to limit ourselves only to integer costs.

4 Paper 5: Plan Optimization

Meysam Aghighi and Christer Bäckström

Plan reordering and parallel execution – a parameterized complexity view
In Proceedings of the Thirty-First AAAI Conference on Artificial Intelligence, San Francisco, California, USA, February 4-9, 2017.

Some real-world applications require a more flexible structure than a sequential plan. For instance the applications that involve real-time execution or multi-agent taskability [75, 76]. One method to model this flexibility is to use partial-order plans. This allows not only different ways to linearize a plan, but also the possibility to postpone making decisions for linearization to run-time. Another important feature of a partial-order plan is its ability to be used as a parallel plan, where mutually unordered actions can be executed in parallel. Some planners directly produce a partial-order plan. At the same time, there has been attempts to produce a partial-order plan from the output of a total-order planner. The core idea is to remove the unnecessary ordering relations and several algorithms have been proposed to do this, both in

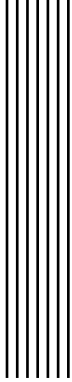
the early 1990s and recently [47, 62, 63, 66, 74]. These algorithms have been applied to various planning domains including workflow planning [68], automatic generation of narratives [40], composition of web services [59], plan execution and monitoring [62], plan repair [69], and distributed control systems [65].

Bäckström [5] theoretically approached the problem and defined deordering and reordering of a partial-order plan. Deordering of a plan is to replace the plan order by a subset of the order such that the new plan remains a valid partial-order plan. Reordering is to replace the plan order with any other order such that the result is a valid partial-order plan. He showed that it is **NP**-complete to optimize the size (number of relation tuples) of both deordering and reordering. There can be other criteria for plan optimization as well. One example is the length of the parallel execution of the plan. Bäckström [5] has also looked into this problem and showed that both finding the shortest parallel execution of a parallel plan (PPL) and optimization of the parallel length by reordering (MPR) are **NP**-complete problems.

In paper 5 of this thesis, we initiate a parameterized complexity investigation of the problems proposed by Bäckström [5] in order to shed more light on them and get more refined results. Two important parameters of a partial-order are its width (the size of the largest anti-chain) and height (the length of the longest chain). We study the effect of these parameters on the parameterized complexity of finding the optimal deordering (MCD) and reordering (MCR). We show that both MCD and MCR are **W**[2]-hard and in **W**[P] if parameterized by the desired order size, and MCD is in **FPT** if parameterized by the original order size.

We further show that the length of the parallel execution of a plan is highly related to both the height of the partial order and the chromatic number of the non-concurrency graph. Moreover, we prove that finding the minimum parallel execution of a plan is in **FPT** if parameterized by the size of the non-concurrency relation.

Finally, we propose a more realistic criterion for plan optimization: the length of the k -processor parallel execution. We study this criterion through the problem PPL_k . We prove that PPL_k is in **FPT** if parameterized by the combination of: (1) the size of the non-concurrency relation, (2) the length of the k -processor parallel execution and (3) the number of processor respectively. On the other hand, the problem remains para-**NP**-hard if only parameterized by $\{n_{\#}, l_p\}$; i.e. significantly harder than the case with unlimited number of processors (PPL). This is consistent with the intuition that the bound on the number of processors makes the problem much harder.



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