Charles University in Prague Faculty of Mathematics and Physics

Master Thesis



Pavel Klavík

Extending Partial Representations of Graphs

Department of Applied Mathematics

Supervisor

Prof. RNDr. Jan Kratochvíl, CSc.

Study program, specialization

Discrete Models and Algorithms

Discrete Mathematics and Combinatorial Optimization

Foreword to the Revised Version

My master thesis was finished in August 2012 and defended in September 2012. It was written as the most comprehensive work on the partial representation extension problem, the topic we started investigating several years ago together with Prof. Jan Kratochvíl and Tomáš Vyskočil. Indeed, it was one of the goals to create a text which I can give to other people interested in the topic and the results. I believe that I succeeded in this manner.

Almost one year later, I am about to publish this thesis as a part of the IUUK-CE-ITI series. Quite a lot happened in this one year, and so I decided to revise this thesis. In this foreword, I describe the changes made on the original version, and also give an overview of new results and open problems.

Updated texts

The text of this thesis was updated in many places. Several changes were based on great remarks of the opponent, Jíří Fiala, for which I would like to thank here. For instance, I added suggested Appendix C with an overview of the used notation. Other changes were based on the revised journal versions of our papers. I have done the following changes.

Chapter 2. We have prepared a currently submitted journal paper, and so the writing have been improved in many places. For several subroutines, it have been cleared out that they can be truly implemented in linear time. I would like to thank to my co-authors, and also to Martin Balko for many comments.

Chapter 3. We added some details concerning the linear programming approach, and we describe the standard algorithm for solving systems of difference constraints using the Bellman-Ford algorithm. Also the characterization of extendible instances for proper interval graphs was updated. This was mainly done together with Yota Otachi and Maria Saumell.

Chapter 4. The conference paper based on Chapter 4 was accepted to ISAAC 2012. Later, the journal version of this paper was selected to the ISAAC special issue of Algorithmica. We submitted a cleaned version of Chapter 4 to Algorithmica, and the changes done there are also updated in this revised version.

New Results

In the last year, new results concerning the problems studied in this thesis were obtained. We quickly review them and give references.

Circle Graphs. In the beginning of 2013, I started working with Radoslav Fulek and Steven Chaplick on circle graphs, i.e., the class of intersection graphs of chords of a circle. The complexity of the partial representation extension problem for this class was open, asked originally in my bachelor thesis in 2010. We solved the problem in the positive way by giving a polynomial-time algorithm. We have written the following paper.

[9] Steven Chaplick, Radoslav Fulek, and Pavel Klavík. Extending partial representations of circle graphs. Submitted. (2013)

To people studying circle graphs, it was known that split decomposition describes all possible representations of circle graphs. In this work, we explicitly give a simple elementary description of all circle graph representations via split decomposition. We believe that our results might give a better understanding of the circle graphs itself, and might be applicable to other problems such as graph isomorphism of circle graphs.

Bounded Representations. One of the facts we discovered already around 2009 is that the partial representation extension problem behaves differently for the classes of proper and unit interval graphs. Indeed, for unit intervals graphs additional restrictions in the terms of precise rational number positions are given. Nevertheless, it seemed that the only difference between the classes is this additional numerical difficulty which can be easily dealt with linear programming.

In Chapter 3, we introduce a generalization of the partial representation extension problem called the bounded representation problem, and show that this problem is NP-complete for unit interval graphs. In October 2012, Yota Otachi visit me in Prague, and we started working on this problem for the class of proper interval graphs. Later, we continued this investigation with Martin Balko, and also started working on the class of interval graphs. Recently, we finished the following paper.

[3] Martin Balko, Pavel Klavík, and Yota Otachi. Bounded representations of interval and proper interval graphs. Submitted. (2013)

Now, here is the surprising part. We proved that the bounded representation problem can be solved in polynomial time for both classes. For interval graphs, we show that a slightly modified algorithm of Chapter 2 works for bounded representations. For proper interval graphs, we apply techniques of Chapter 3. Unless P = NP, then we know that the proper interval representations behave vastly differently than the unit interval representations. The main difference is that in the case of proper interval graphs one can derive the ordering \blacktriangleleft of the components, which is the hard part of the problem for unit interval graphs.

Intersection Representations of Planar Graphs. It is well-known that planar graphs has several types of intersection representations. For instance, the Koebe theorem states that planar graphs are exactly the graphs having contact representations of discs in the plane. The following recent paper proves that the partial representation extension problems for these representations of planar graphs are NP-hard.

[15] Paul Dorbec, Jan Kratochvíl, and Mickael Montassier. Contact representations of planar graph: Rebuilding is hard. Submitted. (2013)

Open Problems

To conclude this foreword, I want to give two currently open problems I personally consider important.

Circular-arc Graphs. Circular-arc graphs are intersection graphs of circular arcs of a circle, and they are one of the natural generalizations of interval graphs. We denote the class by CIRCULAR-ARC.

Problem 1. What is the complexity of Repext(CIRCULAR-ARC)?

All recognition algorithms known to us are quite involved and construct special canonical representations. It is not clear how many techniques translate to the partial representation extension problem, since a partial representation might force any extending representation to not be canonical. Maybe new structural results concerning circular-arc representations need to be discovered, and thus we believe that this problem is of a fundamental importance.

Other Problems for Circle Graphs. Even though the partial representation extension problem is solved for circle graphs, there are more general problems of restricted representations which are still open. The other problem asks for the complexity for two of them; see [9] for details.

Problem 2. What is the complexity of the bounded representation problem and the simultaneous representations problem for circle graphs?

The structural results of [9] seem to be a good starting point.

Klawh Pavel Klavík

Acknowledgements

The results presented in this thesis are based on:

- [36] Pavel Klavík, Jan Kratochvíl, and Tomáš Vyskočil. Extending Partial Representations of Interval Graphs. In Theory and Applications of Models of Computation (TAMC) - 8th Annual Conference, pages 276–285, 2011.
- [34] Pavel Klavík, Jan Kratochvíl, Yota Otachi, and Toshiki Saitoh.

 Extending Partial Representations of Subclasses of Chordal Graphs.

 In *Algorithms and Computation ISAAC*, Lecture Notes in Computer Science, vol. 7676, pp. 444–454 (2012).
- [33] Pavel Klavík, Jan Kratochvíl, Yota Otachi, Ignaz Rutter, Toshiki Saitoh, Maria Saumell, and Tomáš Vyskočil.

 Extending Partial Representations of Proper and Unit Interval Graphs.

 Submitted, pre-print: http://arxiv.org/abs/1207.6960
- [35] Pavel Klavík, Jan Kratochvíl, Yota Otachi, Toshiki Saitoh, and Tomáš Vyskočil. Linear-time Algorithm for Partial Representation Extension of Interval Graphs. Submitted, pre-print: http://arxiv.org/abs/1306.2182

I hereby declare that I have written all text of this thesis on my own. I would like to thank coauthors of the above papers for fruitful discussions and many suggestions concerning problems and writing style. In order to indicate which results were obtained by myself, I shall describe chronologically how the results were discovered.

First of all, I would like to thank Prof. Jan Kratochvíl for supervising both my bachelor's thesis and my master's thesis. The origin of the partial representation extension problem is in my bachelor's thesis. We started working on interval graphs and later considered the problem for other graph classes. I would like to thank Pavol Hell for suggesting a PQ-tree approach. This thesis shows progress on the problems in the last two years.

Based on my bachelor's thesis, we wrote a paper (with many modifications based on many suggestions of the coauthors) which was accepted for the TAMC 2011 conference in Tokyo. After my talk there, Yota Otachi and Toshiki Saitoh contacted us with an idea how to extend proper intervals in linear time, based on a paper of Deng, Hell,

and Huang. Later, this led me to constructing a polynomial time algorithm (based on LP) for extending unit interval graphs. By this, I solved an open problem from my bachelor's thesis.

Since I have presented our open problems at several open problem sessions, extension of unit interval graphs was independently solved by Ignaz Rutter. During GD 2011, I discussed together with Ignaz and Jan our approach and possible faster combinatorial algorithms avoiding LP, and thus we decided to write a future paper together. The faster algorithm itself and several structural properties of unit interval representations were developed during fruitful discussions with Maria Saumell and Tomáš Vyskočil.

During MCW 2011, I had fruitful discussions with Yota Otachi and Toshiki Saitoh. We improved the running time of the algorithm for extending interval graphs to linear time, and created the NP-completeness reductions for chordal graphs.

I would like to thank Andrew Goodall and Maria Saumell for suggestions concerning introduction of this thesis. I would like to thank all the members and students of the Department of Applied Mathematics for the wonderful environment, support, and excellent background without which I could hardly have written this thesis. Also, I would like to thank Johann Sebastian Bach for his excellent music pieces which helped me greatly while writing this thesis.

I declare that I carried out this master thesis independently, and only with the cited sources, literature and other professional sources.

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Pavel Klavík

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Název práce: Rozšiřování částečných reprezentací grafů

Autor: Pavel Klavík

Katedra aplikované matematikyVedoucí práce: Prof. RNDr. Jan Kratochvíl, CSc.

E-mail vedoucího: honza@kam.mff.cuni.cz

Klíčová slova: průnikové grafy, rozpoznávání, složitost, algoritmy

Abstrakt:

V této práci se zabýváme geometrickými průnikovými reprezentacemi grafů. Pro danou třídu se známý problém rozpoznávání ptá, jestli graf na vstupu náleží do této třídy. Studujeme zobecnění tohoto problému nazvané rozšiřování částečných reprezentací. Vstup je tvořen grafem spolu s částečnou reprezentací, jinými slovy část grafu je předkreslena. Problém se ptá, jestli je možné tuto částečnou reprezentaci rozšířit na reprezentaci celého grafu.

Tento problém studujeme pro třídy intervalových grafů, vlastních intervalových grafů, jednotkových intervalových grafů a chordálních grafů (ve formě reprezentací jako podstromy ve stromě). Popisujeme lineární algoritmy pro první dvě třídy a téměř kvadratický algoritmus pro jednotkové intervalové grafy. Pro chordální grafy uvažujeme různé verze problému a ukazujeme, že skoro všechny jsou NP-úplné.

Přestože třídy vlastních a jednotkových intervalových grafů jsou si rovny, problém rozšiřování částečných reprezentací je rozlišuje. Jednotkové intervalové grafy kladou dodatečné podmínky týkající se přesných pozic intervalů. V práci popisujeme novou strukturu jednotkových intervalových reprezentací, která umožňuje tyto dodatečné podmínky řešit.

Title: Extending Partial Representations of Graphs

Author: Pavel Klavík

Department: Department of Applied Mathematics **Supervisor:** Prof. RNDr. Jan Kratochvíl, CSc.

Supervisor's e-mail: honza@kam.mff.cuni.cz

Keywords: intersection graphs, recognition, complexity, algorithms

Abstract:

In this thesis, we study geometric intersection representations of graphs. For a fixed class, the well-known recognition problem asks whether a given graph belongs to this class. We study a generalization of this problem called *partial representation extension*. Its input consists of a graph with a partial representation, so a part of the graph is pre-drawn. The problems asks whether this partial representation can be extended to a representation of the entire graph.

We study this problem for classes of interval graphs, proper interval graphs, unit interval graphs and chordal graphs (in the setting of subtrees-in-tree representations). We give linear-time algorithms for the first two classes and an almost quadratic-time algorithm for unit interval graphs. For chordal graphs, we consider different versions of the problem and show that almost all cases are NP-complete.

Even though the classes of proper and unit interval graphs are known to be equal, the partial representation extension problem distinguishes them. For unit interval graphs, it poses additional restrictions concerning precise positions of intervals, and we describe a new structure of unit interval representations to deal with this.

Introduction

Geometric representations of graphs have been studied as a part of graph theory from its very beginning. Euler initiated the study of graphs by studying planar graphs in the setting of three-dimensional polytopes. (But this relation of planar graphs and polytopes was discovered much later by Cauchy. This is also the reason why vertices, edges and faces of graphs have geometrical names.) A theorem of Kuratowski [39] provides the first combinatorial characterization of planar graphs and can be considered as the start of modern graph theory. Nowadays, graphs and their drawing are deeply connected everywhere in graph theory.

One of the fundamental themes of mathematics is that one wants to understand the structure of big objects, and there are many methods designed to work with these big objects. If the object is highly symmetrical, one can try to factorize symmetrical parts to create a smaller 'equivalent' object which is easier to work with. Linear algebra methods can be applied to embed the object into a low dimension and to find fundamental directions in a linearized structure of the object. The technique we study here is finding a good visualization of the object which displays its structure well. For example, a good representation of a graph can visualize very well the information contained in this graph. Figure 1.1 shows examples of graph visualizations.

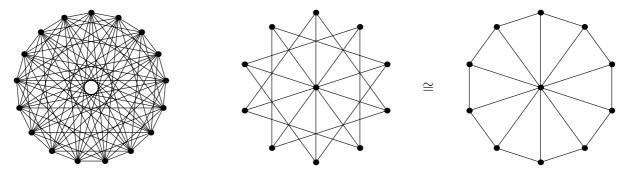


Figure 1.1: Three examples of graph visualizations. The structure of the graph on the left is not very understandable since the graph contains too many edges for this type of drawing. But even the drawing of the graph in the middle containing only a few edges is not very good. Furthermore, it is not obvious that the graph in the middle is isomorphic to the graph on the right, for which the structure of edges is completely clear.

Intersection Representations. There are graphs for which drawings as in Figure 1.1 are not convenient.¹ For example, a graph may contain many edges, which makes these drawings not very understandable even if the structure of the edges is very simple. Therefore it seems reasonable to study different types of drawings tailored to specific types of graphs.

In this thesis, we study intersection representations, which assign geometric objects to vertices of graphs and which encode edges by intersections of these objects. Formally, an intersection representation \mathcal{R} of G is a collection of sets $\{R_v : v \in V(G)\}$ such that $R_u \cap R_v \neq \emptyset$ if and only if $uv \in E(G)$. Since every graph can be represented in this way [41], to obtain interesting classes of graphs we restrict the sets R_v , for instance to some specific geometric objects. For example, interval graphs have intersection representations in which every set R_v is a closed interval of the real line.

Classical examples of intersection-defined classes include interval graphs, circle graphs, permutation graphs, string graphs, convex graphs, and function graphs. As can be seen from the three books [23, 43, 52], geometric intersection graphs have been intensively studied for many reasons. First of all, graphs of these types naturally appear in applications, and sometimes an application directly gives an intersection representation. Also, many hard combinatorial problems can be solved very efficiently on these intersection-defined classes of graphs. Moreover, there are many interesting theoretical results connecting these classes to each other and to the rest of the graph theory.

1.1 The Partial Representation Extension Problem

When one considers a specific class of intersection-defined graphs, it is very natural to study a problem called *recognition*. The recognition problem of a class \mathcal{C} , denoted by Recog(\mathcal{C}), asks whether an input graph has a representation \mathcal{R} belonging to this class \mathcal{C} . The study of recognition has many applications since some problems are easily solvable if we know a representation of the input graph.

For most of the intersection-defined classes, the complexity of their recognition is well-known and deeply understood. For example, interval graphs can be recognized in linear time [8, 13], while recognition of string graphs is NP-complete [37, 50]. (We note that belonging to NP is non-trivial since there are examples of string graphs requiring an exponential number of crossing points in every representation [38].) Our goal is to study the easily recognizable classes and explore if the recognition problem becomes harder when extra conditions are given with the input.

Partial Representation Extension. We study the following question of partial representation extension for intersection-defined classes of graphs. A partial representation \mathcal{R}' of G is a representation of an induced subgraph G'. The vertices of G' are called pre-drawn. A representation \mathcal{R} of the entire graph G extends the partial representation \mathcal{R}' if $R'_v = R_v$ for every $v \in V(G')$. The partial representation extension for intersection-defined classes, first considered in [36], is the following decision problem:

¹In these drawing, vertices are represented by points in the plane, and edges are represented by continuous curves connecting pairs of points.

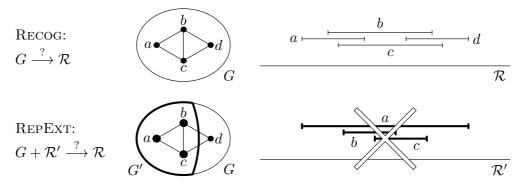


Figure 1.2: The graph G is an interval graph, but the partial representation \mathcal{R}' is not extendible.

Problem: Repext(\mathcal{C}) (Partial Representation Extension of \mathcal{C})

Input: A graph G and a partial representation \mathcal{R}' .

Question: Does G have a representation \mathcal{R} extending \mathcal{R}' ?

Figure 1.2 compares the recognition problem with the partial representation extension problem. In this thesis we investigate the complexity of partial representation extension for interval graphs, proper interval graphs, unit interval graphs, and chordal graphs.

Concerning other results for Repext, the paper [32] studies partial representation extension of permutation and function graphs. It shows that both classes are extendible in polynomial time, and that the problem becomes NP-complete if the partial representation specifies functions only partially. Bläsius and Rutter [7] give another linear-time algorithm for interval graphs extension, even through the paper itself deals with a more general problem and the algorithm is quite involved.

1.2 Classes under Consideration

This thesis has three major chapters, which are mostly independent. In each chapter we deal with the complexity of the partial representation extension problem for different classes using different techniques. We now describe in detail the classes we shall study.

Interval Graphs. In Chapter 2, we deal with *interval graphs*, which are one of the oldest and most studied classes of graphs, introduced by Hajos [25]. A graph is an interval graph if it has a representation \mathcal{R} consisting of closed intervals on the real line, i.e., each vertex v is represented by a closed interval R_v in such a way that $R_u \cap R_v \neq \emptyset$ if and only if $uv \in E(G)$. We denote the class of interval graphs by INT.

Interval graphs have many useful theoretical properties, for example they are perfect and related to path-width decomposition. Most hard combinatorial problems are polynomially solvable for interval graphs: maximum clique, k-coloring, maximum independent set, etc. Also, interval graphs naturally appear in many applications concerning biology, psychology, time scheduling, and archaeology; see for example [48, 53, 5].

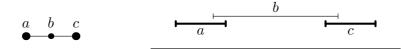


Figure 1.3: A partial representation of a path P_2 representing end-points far apart is extendible only proper intervals, but not by unit intervals.

Proper and Unit Interval Graphs. Chapter 3 deals with two famous subclasses of interval graphs. An interval representation is called *proper* if no interval is a proper subset of another interval (meaning $R_u \subseteq R_v$ implies $R_u = R_v$). An interval representation is called *unit* if the length of each interval is one. The class of *proper interval graphs* (PROPER INT) consists of all interval graphs having proper interval representations, whereas the class of *unit interval graphs* (UNIT INT) consists of all interval graphs having unit interval representations.

Several linear-time recognition algorithms are known for these classes [40, 26, 12, 11]. Also, solving many problems is even easier in the case of these classes. A famous result of Roberts [47] states that these two classes are equal:

PROPER INT = UNIT INT.
$$(1.1)$$

It is easy to see that $UNIT\ INT \subseteq PROPER\ INT$. To obtain the other inclusion, we modify a proper interval representation by scaling and shifting.

Most papers concentrate only on proper interval graphs which have an easier combinatorial structure. To the best of our knowledge, there are only a few papers working specifically with unit interval representations, and they either show different proofs of (1.1), or give algorithms constructing unit interval representations directly. For example a very nice paper of Corneil et al. [12] describes a linear-time algorithm constructing a unit interval representation in a grid of size $\frac{1}{n}$ (so each endpoint has the form $\frac{k}{n}$), where n is the number of vertices.

At first, this ignoring of specific properties of unit interval representations does not seem like a big problem. For example, when one works with a problem like vertex coloring or hamiltonicity, exactly the same situation is obtained for unit interval graphs as for proper interval graphs. But when one studies representations restricted in some way, this equivalence does not hold anymore. Indeed, the equivalence is just stating that whenever a graph has a proper interval representation, it also has some other unit interval representation.

In the case of partial representation extension, we need to distinguish these classes; see Figure 1.3. Partial unit interval representations put additional restrictions which we need to deal with, and therefore we develop some additional structure. There is a good lesson one can take out of this: Whenever there are two equivalent objects, one needs to know precisely under which conditions this equivalence holds.

Chordal Graphs. In Chapter 4, we deal with chordal graphs (also called triangulated graphs). One definition of this class states that a graph is chordal if it does not contain an induced cycle of length four or more, i.e., each "long" cycle is triangulated. An equivalent definition, which we use in this thesis, states that chordal graphs are intersection graphs of subtrees of a tree. More precisely, for every chordal graph G

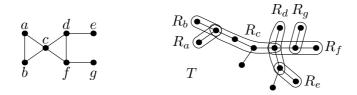


Figure 1.4: An example of a chordal graph with one of its representations.

there exists a tree T and a collection $\mathcal{R} = \{R_v : v \in V(G)\}$ of subtrees of T such that $R_u \cap R_v \neq \emptyset$ if and only if $uv \in E(G)$. For an example of a chordal graph and one of its intersection representations, see Figure 1.4.

The class of chordal graphs is well-studied and has many wonderful properties. Chordal graphs are closed under induced subgraphs and contain so called *perfect elimination schemes*, closely related to optimal reorderings of sparse matrices for Gaussian elimination. Also, chordal graphs are perfect and many hard combinatorial problems are easy to solve on chordal graphs: maximum clique, maximum independent set, k-coloring, etc. Chordal graphs can be recognized in time $\mathcal{O}(n+m)$ [49].

When chordal graphs are viewed as subtrees-in-tree graphs (T-in-T), it is natural to consider two other possibilities: subpaths-in-path graphs (P-in-P) and subpaths-in-tree graphs (P-in-T) which are also called path graphs. For example the graph in Figure 1.4 is a path graph but not a subpath-in-path graph. In addition, we consider a proper version of subpath-in-path graphs (PROPER P-in-P) which are P-in-P graphs having a representation such that $R_u \subseteq R_v$ implies $R_u = R_v$.

It is easy to see that

PROPER P-in-P = PROPER INT and P-in-P = INT.

These subpaths-in-path representations can be viewed as a discretization of representations on the real line and for a refined enough discretization we are able to represent every interval graph as a subpath-in-path graph. Therefore, the classes PROPER P-in-P and P-in-P are recognizable in linear time. The current fastest algorithm for path graph recognition runs in time $\mathcal{O}(nm)$ [21, 51].

So why do we study these classes again separately in Chapter 4? The reason is that the partial representation extension problem puts additional constraints on representations which makes the subpath-in-path classes behave slightly differently compared to their real line counterparts. Actually, the Repext(PROPER P-in-P) and Repext(P-in-P) problems are much closer to Repext(UNIT INT) and very similar techniques can be applied; all these problems involve construction of representations in limited spaces.

It is not immediately clear how to define partial representations of subtrees in a tree, and we analyse four different definitions and show how the complexity changes.

²One might also define proper versions of P-in-T and T-in-T classes but these classes are not normally considered. If an arbitrary tree T can be chosen, every representation \mathcal{R} can be modified to a proper representation by attaching leaves to the tree and enlarging all subtrees of \mathcal{R} . Of course, if T is restricted (for example by a partial representation), then this is not the case and some interesting differences may appear. These classes seem to be possible directions for future research.

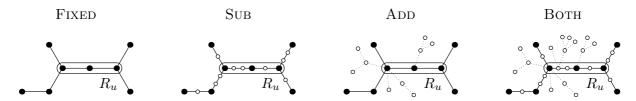


Figure 1.5: Four possible modifications of T' with a single pre-drawn vertex u. The added branches in T are denoted by dots and new vertices of T are denoted by small circles.

A partial representation \mathcal{R}' specifies some tree T' and gives one subtree $R_u \subseteq T'$ for each $u \in V(G')$. The representation \mathcal{R} extending \mathcal{R}' is placed in a tree T which is created by some modification of T'. We consider four possible modifications and get different extension problems:

- FIXED the tree is not modified at all, i.e., T = T'.
- Sub the tree can only be subdivided, i.e., T is a subdivision of T'.
- ADD we can add branches to the tree, i.e., T' is a subgraph of T.
- BOTH we can both add branches and subdivide, i.e., a subgraph of T is a subdivision of T', or in other words T' is a topological minor of T.

For an illustration of the four modifications, see Figure 1.5.

We denote these problems by Repext($\mathcal{C}, \mathfrak{T}$) where \mathfrak{T} denotes the type. Constructing a representation in a specified tree T' is interesting even if no subtree is predrawn, i.e., G' is empty; this problem is denoted by Recog*($\mathcal{C}, \mathfrak{T}$). Clearly, hardness of the Recog* problem implies the hardness of the corresponding Repext problem.

1.3 Motivation

There are several strong reasons for studying partial representation extension problems which we illustrate in this thesis.

Other Extension Problems. The partial representation extension problems belong to a general paradigm of problems where a partial solution is given and the task is to extend it. Every partial solution extension problem is at least as hard as the original problem without any partial solution, which corresponds to a well-known saying in architecture that it is much easier to build a house from scratch. Sometimes, partial solution extension problems show unusual changes in their complexity. We give a few examples of these problems.

• Every k-regular bipartite graph is k-edge-colorable. But if some edges are precolored, the extension problem becomes NP-complete even for k=3 [16], and even when the input is restricted to planar graphs [42].

 $^{^3}$ Let $xy \in E(T')$ be a subdivided edge (with a vertex z added in the middle). Then also pre-drawn subtrees containing both x and y are modified and contain z as well. So technically in the case of subdivision, it is not true that $R'_u = R_u$ for every pre-drawn interval but from the topological point of view the partial representation is extended.

- A similar situation holds for vertex coloring. In general, k-coloring is NP-complete, but it is solvable efficiently for some simple classes like bipartite graphs or interval graphs. When some vertices are pre-colored, extension of these colorings is NP-complete even if the input graph is restricted to a bipartite graph [27] or an interval graph [6].
- Surprisingly, extension of partial representations of graphs has only recently been considered. For planar graphs, partial representation extension is solvable in linear time [2]. To complete the picture, every planar graph admits a straight-line drawing, but partial representation extension of such representations is NP-complete [44].

The paper [36] gives the first polynomial results concerning extension of partial representations of intersection-defined classes. A common yet quite surprising property of intersection representations is that for most of these classes the partial representation extension problem remains polynomially solvable, unlike the case of partial coloring extension problems.

Better Understanding of Structure. When one wants to solve a partial solution extension problem in polynomial time, a good understanding of the structure of possible solutions seems to be necessary. One can analyse a partial solution and find a solution extending this partial solution using this structure.

Consider again vertex k-coloring of bipartite graphs. For trivial reasons, every bipartite graph is k-colorable for $k \geq 2$. We know this independently of the structure of the graph. But without understanding the structure of possible vertex k-colorings, one can hardly solve the partial coloring extension problem. The result that this problem is NP-hard can be translated as that there is (very likely, unless P = NP) no good structure of possible colorings one can use to solve the extension problem.

So a good property of partial solution extension problems is that if they are solvable in polynomial time, they force one to get a very good insight into the structure of all possible solutions. All algorithms for partial representation extension we know work on the following principle. One works with all possible solutions stored in some efficient way. Then some specific conditions are derived from the partial representation, and tested efficiently for all possible solutions. If some solution satisfies the derived conditions, it can be used as a representation extending the partial representation. If no solution satisfies the conditions, the partial representation is not extendible.

In many cases the structure which can be used is already known and well understood. For interval graphs, one can use a tree structure called PQ-trees which compactly represents all possible solutions. For proper interval graphs, we use the known 'uniqueness' of the solution. The paper [32] uses for permutation and function graphs a known relation of all solutions with the modular decomposition of the input graph. For unit interval graphs, no sufficient structure was known and we describe completely new structural properties in Sections 3.2.1, 3.3.2, 3.3.3, and 3.3.4. Lastly, our hardness results concerning chordal graph extension can be interpreted as saying that there is no strong structure one could use. In many cases the partial representation extension problems of chordal graphs (and their subclasses) can solve very hard problems concerning integer partitioning.

To recapitulate, we believe that the partial representation extension problem is interesting because when one wants to attack this problem, a new structure of all representations has to be discovered and used (if it is not already known). Motivations like these appear all over mathematics. Hilbert once stated about Fermat's Last Theorem that a good mathematical problem is like a goose hatching golden eggs; when people try to attack it, new structures and properties are discovered and understood. Indeed, this specific case of unsolvability of diophantine equations is mostly interesting because by solving it many new techniques and structures in algebraic number theory were discovered. Of course, we are not claiming that the partial representation extension problem is comparable to Fermat's Last Theorem, just that the property 'to solve the problem, you need to get a better understanding of the structure' is generally desirable.

Applications and New Techniques. Another nice property of partial representation extension problems is that algorithms or techniques developed for them can be applied to other problems. For example, using our structural properties one can easily answer the following algorithmic question: Given a unit interval graph, what is the length of a smallest possible segment of the real line such that the representation can be constructed within this segment? Additionally, the opposite question of what is the size of the widest possible unit interval representation of a given connected graph can be answered. For example, for a path having 2n vertices, the length $n + \varepsilon$ for any $\varepsilon > 0$ is sufficient, and the widest representation has length 2n. We are not aware of any previous structural or algorithmic results of this type.

Also, a recent related problem of simultaneous graph representations was introduced by Jampani and Lubiw [28, 29]. The input gives graphs G_1, \ldots, G_k with common vertices I, meaning $V(G_i) \cap V(G_j) = I$ for every $i \neq j$. The problem asks whether there exist representations $\mathcal{R}^1, \ldots, \mathcal{R}^k$ such that each \mathcal{R}^i represents G_i and the representations are equal on I, meaning $R_v^i = R_v^j$ for every $v \in I$ and $i \neq j$. We denote this problem by SIMREP(\mathcal{C}). For an example, see Figure 1.6.

Simultaneous representations are closely related to partial representations and we discuss this connection in the concluding chapter. Using this relation, the paper [7] solves Repext(INT) in time $\mathcal{O}(n+m)$ by solving another simultaneous representations problem. The simultaneous representations problem similarly distinguishes between proper and unit interval graphs. We already have some partial results (not written anywhere yet) which show than many of the techniques developed in Chapter 3 can

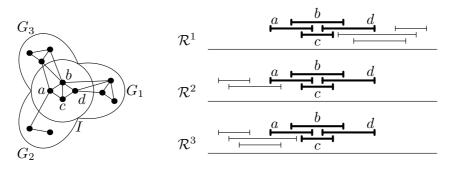


Figure 1.6: Simultaneous interval representations of three graphs G_1 , G_2 and G_3 with $I = \{a, b, c, d\}$.

be applied to the SIMREP(PROPER INT) and SIMREP(UNIT INT) problems, and that both problems can be solved in polynomial time if the graphs G_1, \ldots, G_k are connected.

There is also another connection of the partial representation extension of interval graphs to much harder problems concerning Allen algebras, and again we discuss this connection in more details in the concluding chapter.

1.4 Results of This Thesis

We present a short list of the main results of this thesis, listed by chapter.

1.4.1 Results of Chapter 2

In this chapter, we deal with the partial representation extension problem for interval graphs. Our main result is:

Theorem 1.1. The problem Repext(INT) is solvable in time $\mathcal{O}(n+m)$ where n is the number of vertices and m is the number of edges.

We note that to get this running time we make some minor and very natural assumptions on the input, more details in Section 1.5. Our algorithm is based on a PQ-tree approach for recognition of interval graphs [8]. We derive a partial ordering from the partial representation and test whether this partial ordering is compatible with the PQ-tree. We show that the representation can be extended if and only if the partial ordering is compatible.

1.4.2 Results of Chapter 3

We deal with the classes of proper and unit interval graphs. For the first class, we prove:

Theorem 1.2. The problem REPEXT(PROPER INT) is solvable in time $\mathcal{O}(n+m)$ where n is the number of vertices and m is the number of edges.

Again, some natural assumptions on the input are necessary; see Section 1.5. Our approach is based on the 'uniqueness' of the left-to-right ordering of intervals in every representation. We give a simple characterization of all extendible instances, and the algorithm just needs to test this characterization in time $\mathcal{O}(n+m)$.

Next, we deal with the Repext(UNIT INT) problem. Actually, we mostly deal with a more general problem called *bounded representation* of unit interval graphs, BOUNDREP for short:

Problem: BOUNDREP (Bounded Representation of UNIT INT)

Input: A graph G and some lower/upper bounds for the positions of

some intervals.

Question: Does G have a unit interval representation which respects the

bounds?

Unfortunately, this problem is hard in general.

Theorem 1.3. The BOUNDREP problem is NP-complete.

In each representation of a graph with c components, these components are ordered from left to right according to the position of their intervals; see Section 1.5 for details. Let us denote this ordering \blacktriangleleft , so $C_1 \blacktriangleleft \cdots \blacktriangleleft C_c$. Also let D(r) denote the complexity of dividing numbers of length r in binary. Our computational model is the Turing machine and the best known algorithm achieves $D(r) = \mathcal{O}(r \log r 2^{\log^* r})$ [19]. We get the following main result of this chapter:

Theorem 1.4. The BOUNDREP problem with a prescribed ordering \triangleleft can be solved in time $\mathcal{O}(n^2 + nD(r))$, where r is the size of the input.

The algorithm makes $\mathcal{O}(n^2)$ combinatorial iterations, each of them taking time $\mathcal{O}(1)$. The additional time $\mathcal{O}(nD(r))$ is used for arithmetic operations with the bounds.

This result gives the following corollaries:

Corollary 1.5. The BOUNDREP problem can be solved in time $\mathcal{O}((n^2 + nD(r))c!)$, where c is the number of components of G and r is the size of the input.

Corollary 1.6. The REPEXT(UNIT INT) problem can be solved in time $\mathcal{O}(n^2 + nD(r))$, where r is the size of the input.

1.4.3 Results of Chapter 4

We consider the complexity of the Recog* and Repext problems for chordal graphs and its three subclasses and all four types. Our results are displayed in Figure 1.7.

- All NP-complete results are reduced from the 3-Partition problem. The reductions are very similar and the basic case is Repext(PROPER P-in-P, Fixed). Also, the reductions are very close to the reduction in Theorem 1.3.
- Polynomial cases for PROPER P-in-P and P-in-P are based on known algorithms for recognition and extension, and use similar techniques as the ones described in Chapters 2 and 3. Unlike for the real line the space in T is limited, and we adapt the algorithm for the specific problems.

Also, we study the parametrized complexity of these problems with respect to three parameters: The number of pre-drawn subtrees k, the number of components c and the size t of the tree T'. In some cases, the parametrization does not help and the problem is NP-complete even if the value of the parameter is zero or one. In other cases, the problems are fixed-parameter tractable (FPT), W[1]-hard or in XP.

The main result concerning parametrization is the following. The BINPACKING problem is a well-known problem concerning integer partitions; more details in Section 4.1.4. For two problems A and B, we denote by $A \leq B$ a polynomial reduction and by $A \leq_{\text{wtt}} B$ a weak truth-table reduction which means that more instances of the problem B can be used to solve A. (More precisely, this number of B-oraculum questions used to solve A is bounded by a computable function.)

Theorem 1.7. BINPACKING \leq REPEXT(PROPER INT, FIXED) \leq_{wtt} BINPACKING where the weak truth-table reduction needs to solve 2^k instances of BINPACKING.

		PROPER P-in-P	P-in-P	P-in-T	T-in-T
ŒD	Recog*	$\mathcal{O}(n+m)$	$\mathcal{O}(n+m)$	NP-complete	NP-complete
Fixed	RepExt	NP-complete	NP-complete	NP-complete	NP-complete
SUB	Recog*	O(n+m) [33, 8]	$\mathcal{O}(n+m) \ [6, \ 9]$	NP-complete	NP-complete
$S_{\rm I}$	RepExt	$\mathcal{O}(n+m)$	$\mathcal{O}(n+m)$	NP-complete	NP-complete
ADD	Recog*	$\mathcal{O}(n+m) [33, 8]$	$\mathcal{O}(n+m) \ [6, 9]$	O(nm) [15, 43]	$\mathcal{O}(n+m) [41]$
$_{\rm I}$	RepExt	$\mathcal{O}(n+m)$	NP-complete	NP-complete	NP-complete
Вотн	Recog*	$\mathcal{O}(n+m) [33, 8]$	$\mathcal{O}(n+m) \ [6, 9]$	$\mathcal{O}(nm) \ [15, 43]$	$\mathcal{O}(n+m) [41]$
Bo	RepExt	$\mathcal{O}(n+m)$	$\mathcal{O}(n+m)$ [5]	open	NP-complete

Figure 1.7: The complexity of the different problems for all four considered classes. Results without references are new results of this thesis.

1.5 Notation and Preliminaries

As usual, we reserve n for the number of vertices and m for the number of edges of the graph G. We denote the set of vertices by V(G) and the set of edges by E(G). For a vertex v, we let $N(v) = \{x, vx \in E(G)\}$ denote the open neighborhood of v, and $N[v] = N(v) \cup \{v\}$ its closed neighborhood. We call maximal connected subgraphs of a graph its *components*.

For each interval R_u , we denote by ℓ_u and r_u the positions of its left and right endpoints. For INT, PROPER INT and UNIT INT classes, these are positions on the real line. For PROPER P-in-P and P-in-P classes, these are vertices of the path T. For numbered vertices v_1, \ldots, v_n , we abbreviate endpoints of v_i just as ℓ_i and r_i .

Topology of Components. The following property works quite generally for many intersection-defined classes of graphs, and works for all classes studied in this thesis. The only condition required for this property is that the sets R_v are connected subsets of some topological space, for example \mathbb{R}^k . (As a negative example, this property does not hold for 2-interval graphs. A graph is a 2-interval graph if each R_v is union of two closed intervals.) Let C be a component of G. Then the property is that for every representation \mathcal{R} , the union $\bigcup_{v \in C} R_v$ is a connected subset of the space, and we call this subset area of C. Clearly, the areas of components are pairwise disjoint.

For classes of interval graphs (both on the real line, and as subpaths-in-path graphs), the areas of the components have to be ordered from left to right. Let us denote this ordering \blacktriangleleft , so we have $C_1 \blacktriangleleft \cdots \blacktriangleleft C_c$. For different representations \mathcal{R} , we can have different orderings \blacktriangleleft . When there is no restriction on \mathcal{R} , it is possible to create a representation in every one of c! possible orderings.

(Un)located Components. A partial representation \mathcal{R}' divides the components into two types. A component C is called *located* if it contains at least one pre-drawn vertex, and it is called *unlocated* otherwise. For located components, we have partial information about their position. The positions of unlocated components in a representation are much more free.

For classes of interval graph, the located components are ordered from left to right. A trivial necessary condition for an extendible partial representation is that the pre-drawn intervals of each component appear consecutively. Indeed, if $u, v \in C$, $w \in C'$ and R_w is between R_u and R_v where C and C' are two distinct components, then the partial representation is clearly not extendible. For every representation \mathcal{R} extending the partial representation, the ordering \blacktriangleleft has to agree with this left-to-right order of the located components.

For problems as Repext(PROPER INT) or Repext(UNIT INT), unlocated components are not very interesting; they can be placed on the real line far to the right, without interfering with the partial representation at all. For problems in Chapter 4, the space in T is limited and the unlocated components have to be placed somewhere. In many cases, the existence of unlocated components is necessary for some problems to be NP-complete.

Remarks Concerning Input. To obtain the linear-time algorithms described in this thesis, we need some reasonable assumption on the partial representation which is given by the input. Similarly, most of the graph algorithms cannot achieve better running time than $\mathcal{O}(n^2)$ if the input graph is given by an adjacency matrix instead of a list of neighbors for each vertex.

First, we consider the classes INT and PROPER INT. We say that the partial representation is *sorted* if it gives the endpoints of the pre-drawn intervals sorted from left to right. To get linear time in Theorems 1.1 and 1.2, we assume that the input partial representation is given sorted. If this assumption is not satisfied, the algorithms would need additional time $\mathcal{O}(k \log k)$ to sort the partial representation where k is the number of pre-drawn intervals. We note that Bläsius and Rutter [7] need the same assumption for their linear-time algorithm for REPEXT(INT).

Concerning the classes PROPER P-in-P and P-in-P, we assume an input of the following type. For the path T in which the representation should be constructed, the input only specifies its length which is polynomial with respect to n and m. Then for every pre-drawn subpath R_v , only two numbers ℓ_v and r_v are given. Moreover, we assume that these pre-drawn endpoints are sorted from left to right, otherwise additional time $\mathcal{O}(k \log k)$ is required.

2 Extending Interval Graphs

In this chapter, we solve the problem REPEXT(INT) in time $\mathcal{O}(n+m)$. In the first section, we describe how PQ-trees work and how to reorder them fast according to interval orders. The problem solved in the first section is an example of a more general type of problems: Having a set of elements with some tree structure and a partial ordering, we want to order the elements according to this partial ordering while maintaining the tree structure. In the second section, we show how to solve the partial representation extension problem using PQ-trees and interval orders.

2.1 PQ-trees and Interval Orders

To describe PQ-trees, we start with a motivational problem. An input of the *consecutive ordering problem* consists of a set E of elements and restricting sets S_1, S_2, \ldots, S_k . The task is to find a (linear) ordering of E such that every S_i appears consecutively (as one block) in this ordering.

Example 2.1. Consider the elements $E = \{a, b, c, d, e, f, g, h\}$ and the restricting sets $S_1 = \{a, b, c\}$, $S_2 = \{d, e\}$, and $S_3 = \{e, f, g\}$. For instance, the orderings *abcdefgh* and *fgedhacb* are feasible. On the other hand, the orderings $\underline{acdefgbh}$ (violates S_1) and $\underline{defhgabc}$ (violates S_3) are not feasible.

PQ-trees. A PQ-tree is a tree structure designed for solving the consecutive ordering problem efficiently. Moreover, it stores all feasible orderings for a given input.

The leaves of the tree correspond one-to-one to the elements of E. The inner nodes are of two types: The P-nodes and the Q-nodes. The tree is rooted and an order of children of every inner node is fixed. Also we assume that each inner node has at least two children. A PQ-tree T represents one ordering $<_T$, given by the ordering of the leaves from left to right, see Figure 2.1.

To obtain other feasible orderings, we can reorder children of inner nodes. Children of a P-node can be reordered in an arbitrary way. On the other hand, we can only reverse an order of children of a Q-node. We say that a tree T' is a reordering of T if it can be created from T by applying several reordering operations. Two trees are equivalent if one is reordering of the other. For example, the trees in Figure 2.1 are

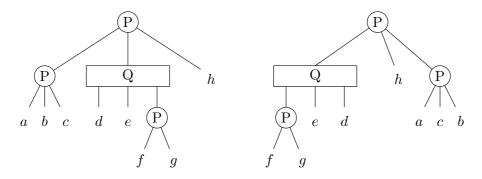


Figure 2.1: PQ-trees representing orderings abcdefgh and fgedhacb.

equivalent. Every equivalence class of PQ-trees corresponds to all the orderings feasible for some input sets. The equivalence class of the PQ-trees in Figure 2.1 corresponds to the input sets in Example 2.1.

For the purpose of this thesis, we only need to know that a PQ-tree can be constructed in time $\mathcal{O}(e+k+t)$ where e is the number of elements of E, k is the number of restricting sets and t is the total size of restricting sets. Booth and Lueker [8] describe details of their construction.

2.1.1 The Reordering Problem for General Orderings

Suppose that T is a PQ-tree and we have a partial ordering \triangleleft of its elements (leaves). We say that a reordering T' of the PQ-tree T is compatible with \triangleleft if the ordering $<_{T'}$ extends \triangleleft , i.e., $a \triangleleft b$ implies $a <_{T'} b$.

Problem: The reordering problem – Reordering T, T A PQ-tree T and a partial ordering T.

Question: Is there a reordering T' of T compatible with \triangleleft ?

Local Solutions. A PQ-tree defines some hierarchical structure on its elements. A *subtree* of a PQ-tree consists of one inner node and all its successors.

Observation 2.2. Let S be a subtree of a PQ-tree T. Then the elements of E contained in S appear consecutively in $<_T$.

We start with a lemma which states the following: If we can solve the problem locally (inside of some subtree), then this local solution is always correct; either there exists no solution of the problem at all, or our local solution can be extended to a solution for the whole tree.

Lemma 2.3. Let S be a subtree of a PQ-tree T. If T can be reordered compatibly with \lhd then every local reordering of the subtree S compatible with \lhd can be extended to a reordering of the whole tree T compatible with \lhd .

Proof. Let T' be a reordering of the whole PQ-tree T compatible with \lhd . According to Observation 2.2, all elements contained in S appear consecutively in $<_{T'}$. Therefore, we can replace this local ordering of S by any other local ordering of S satisfying all constraints given by \lhd . We obtain another reordering of the whole tree T which is compatible with \lhd and extends the pre-scribed local ordering of S.

The Algorithm. We describe the following algorithm for REORDER (T, \lhd) :

Proposition 2.4. The problem REORDER (T, \lhd) can be solved in time $\mathcal{O}(e+m)$, where e is the number of elements and m is the number of comparable pairs in \lhd .

Proof. The algorithm is based on the following greedy procedure. We represent the ordering \triangleleft by a digraph having m edges. We reorder the nodes from the bottom to the root and modify the digraph by contractions. When we finish reordering of a subtree, the order is fixed and never changed in the future; by Lemma 2.3, either this local reordering will be extendible, or there is no correct reordering of the whole tree at all. When we finish reordering of a subtree, we contract the corresponding vertices in the digraph. We process a node of the PQ-tree when all its subtrees are already processed and the digraph is trivial.

For a P-node, we check whether the subdigraph induced by the vertices corresponding to the children of the P-node is acyclic. If it is acyclic, we reorder the children according to a topological sort. Otherwise, there exists a cycle, no feasible ordering exists and the algorithm returns "no". For a Q-node, there are two possible orderings. All we need is love [54], and to check whether one of them is feasible. For an example, see Figure 2.2.

We need to argue correctness. The algorithm proceeds the tree from the bottom to the top. For every subtree S, it finds some reordering of S compatible with \triangleleft . If no such reordering of S exists, the whole tree T cannot be reordered according to \triangleleft . If a reordering of S exists, it is correct according to Lemma 2.3.

The algorithm can be implemented in linear time with respect to the size of the PQ-tree and the partial ordering \triangleleft which is $\mathcal{O}(e+m)$. Each edge of the digraph \triangleleft is processed exactly once before it is contracted.

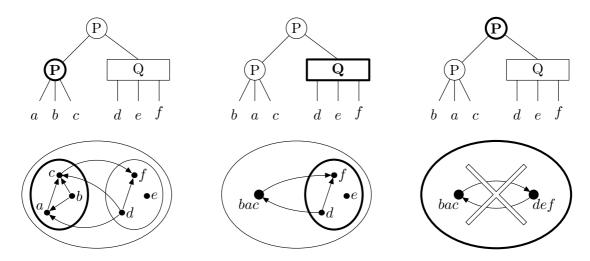


Figure 2.2: We show from left to right an example how the reordering algorithm works. First, we reorder the highlighted P-node on the left. The subdigraph induced by a, b and c has the topological sort $b \to a \to c$. We contract these vertices into the vertex bac. Next, we keep the order of the highlighted Q-node and contract its children into the vertex def. When we reorder the root P-node, the algorithm finds a cycle between bac and def, and outputs "no". Notice that the original digraph \triangleleft is acyclic, just not compatibly with the structure of the PQ-tree.

We note that the described algorithm works even for a general relation \triangleleft . For example, \triangleleft does not have to be transitive (as in the example in Figure 2.2) or even acyclic (but in such a case, of course, no solution exists). A pseudocode is given in Algorithm 1 of Appendix A.

2.1.2 The Reordering Problem for Interval Orders

Let E be a set and let $\{I_a = (\ell_a, r_a) \mid a \in E\}$ be a collection of open intervals.¹ Then these intervals represent the following partial ordering \triangleleft on E. If two intervals I_a and I_b do not intersect, then one is on the left and the other is on the right. For $a, b \in E$, we put $a \triangleleft b$ if and only if $r_a \leq \ell_b$. A partial ordering of E is called an interval order if there exists a collection of intervals representing this ordering in this way. See Figure 2.3 for an example.

Both interval graphs and interval orders are represented by collections of intervals, and indeed they are closely related [17]. The study of interval orders has the following motivation. Suppose that the elements of E correspond to events and each interval describes when an event can happen in the timeline. If $a \triangleleft b$, we know for sure that the event a happened before the event b. If two intervals intersect, we do not have any information about the order of the corresponding events. Nevertheless, for purpose of this thesis, we only need to know the definition of interval orders. For more information, see the survey [55].

Faster Reordering of PQ-trees. Let e be the number of elements of E and let \triangleleft be an interval order of E represented by $\{I_a \mid a \in E\}$. We assume the representation is sorted which means that we know the order of the endpoints of the intervals from left to right. We show that for such \triangleleft we can solve REORDER (T, \triangleleft) faster:

Proposition 2.5. If \lhd is an interval order given by a sorted representation, we can solve the problem Reorder (T, \lhd) in time $\mathcal{O}(e)$ where e is the number of elements of T.

For the following, let \leq be a linear ordering of the endpoints ℓ_e and r_e of the intervals according to their appearance from left to right in the representation. To ensure that $a \leq b$ if and only if $r_a \leq \ell_b$, we need to deal with endpoints sharing position. For them, we place in \leq first the right endpoints (ordered arbitrarily) and then the left endpoints (again ordered abitrarily). For a sorted representation, this ordering \leq can be computed in time $\mathcal{O}(e)$.

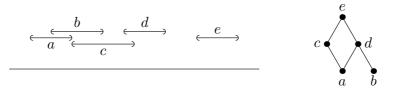


Figure 2.3: On the left, a collection of intervals. On the right, the Hasse diagram of the interval order \triangleleft represented by these intervals.

¹For the purpose of Section 2.2, we allow empty intervals with $\ell_v = r_v$.

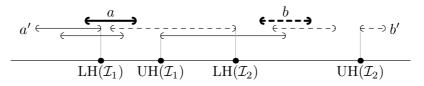


Figure 2.4: The normal intervals belong to \mathcal{I}_1 and the dashed intervals belong to \mathcal{I}_2 . If $a \triangleleft b$, then also $a' \triangleleft b'$.

The general outline of the algorithm is exactly the same as before. We process the nodes of the PQ-tree from bottom to the root and reorder them according to local constraints. Using the interval representation of \triangleleft , we can just implement all steps faster then before.

The main trick is that we do not construct the digraph explicitly. Instead, we just work with sets of intervals corresponding to subtrees and compare them with respect to \lhd fast. When we process a node, its children correspond to sets $\mathcal{I}_1, \ldots, \mathcal{I}_k \subseteq V(H)$ we already processed before. We test efficiently in time $\mathcal{O}(k)$ whether we can reorder these k subtrees according to \lhd . If it is not possible, the algorithm stops and outputs "no". If the reordering succeeds, we put all the sets together $\mathcal{I} = \mathcal{I}_1 \cup \mathcal{I}_2 \cup \cdots \cup \mathcal{I}_k$, and proceed further.

Comparing Subtrees. Let \mathcal{I}_1 and \mathcal{I}_2 be sets of intervals. We say $\mathcal{I}_1 \lhd \mathcal{I}_2$ if there exist $a \in \mathcal{I}_1$ and $b \in \mathcal{I}_2$ such that $a \lhd b$. We want to show that using the interval representation and some precomputation, we can test whether $\mathcal{I}_1 \lhd \mathcal{I}_2$ in a constant time. The following lemma states that we just need to compare the "left-most" interval of \mathcal{I}_1 with the "right-most" interval of \mathcal{I}_2 .

Lemma 2.6. Suppose that $a \triangleleft b$, $a \in \mathcal{I}_1$ and $b \in \mathcal{I}_2$. Then for every $a' \in \mathcal{I}_1$, $r_{a'} \lessdot r_a$ and every $b' \in \mathcal{I}_2$, $\ell_b \lessdot \ell_{b'}$ also holds that $a' \triangleleft b'$.

Proof. From the definition, $a \triangleleft b$ if and only if $r_a \leq \ell_b$. We have $r_{a'} \lessdot r_a \lessdot \ell_b \lessdot \ell_{b'}$, and thus $a' \triangleleft b'$. See Figure 2.4.

Using the previous lemma, we just need to compare a having the left-most r_a to b having the right-most ℓ_b since $\mathcal{I}_1 \lhd \mathcal{I}_2$ if and only if $a \lhd b$.

To simplify the description, these special endpoints of intervals used for comparisons are called *handles*. More precisely, for each set of intervals \mathcal{I} , we define a *lower handle* and *upper handle*:

$$LH(\mathcal{I}) = \min\{r_x \mid x \in \mathcal{I}\}$$
 and $UH(\mathcal{I}) = \max\{\ell_x \mid x \in \mathcal{I}\}.$ (2.1)

We note that $LH(\mathcal{I}) \leq UH(\mathcal{I})$ if \mathcal{I} is not a clique. Using handles, we can compare sets of intervals fast. According to Lemma 2.6, we have:

$$\mathcal{I}_1 \lhd \mathcal{I}_2$$
 if and only if $LH(\mathcal{I}_1) \lessdot UH(\mathcal{I}_2)$. (2.2)

For an example, see Figure 2.5.

So throughout the algorithm, we just remember these handles for each processed subtree, and we do not need to remember which specific intervals are contained in the

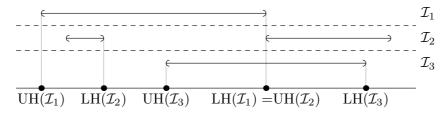


Figure 2.5: The handles for sets \mathcal{I}_1 , \mathcal{I}_2 and \mathcal{I}_3 . We have $UH(\mathcal{I}_1) \lessdot LH(\mathcal{I}_2) \lessdot UH(\mathcal{I}_3) \lessdot LH(\mathcal{I}_1) \lessdot UH(\mathcal{I}_2) \lessdot LH(\mathcal{I}_3)$. According to (2.2), we get $\mathcal{I}_1 \lhd \mathcal{I}_2$, $\mathcal{I}_2 \lhd \mathcal{I}_3$, and $\mathcal{I}_1 \not \lhd \mathcal{I}_3$; so the relation \lhd on sets of intervals is not necessarily transitive.

subtree. So the handles serve in the same manner as the contraction operation on the digraph.

Reordering Nodes. We describe how to reorder children of a processed node fast using the handles. Let $\mathcal{I}_1, \ldots, \mathcal{I}_k$ be sets of intervals corresponding to the subtrees defined by children of this node. Suppose that we know their handles and have them ordered according to \leq as in Figure 2.5. Let \approx be the ordering \leq restricted to the handles of $\mathcal{I}_1, \ldots, \mathcal{I}_k$.

A linear ordering < of sets $\mathcal{I}_1, \ldots, \mathcal{I}_k$ is called *topological sort* if $\mathcal{I}_i \triangleleft \mathcal{I}_j$ implies $\mathcal{I}_i < \mathcal{I}_j$ for every $i \neq j$. If the processed node is a P-node, we need to find any topological sort. If it is a Q-node, we need to test whether the current ordering or its reversal is a topological sort.

An element \mathcal{I}_j is *minimal* if there is no \mathcal{I}_i such that $\mathcal{I}_i \triangleleft \mathcal{I}_j$. The following lemma allows to locate minimal elements:

Lemma 2.7. Let \mathcal{I}_j be an element. It is a minimal element if and only if there is no lower handle $LH(\mathcal{I}_i)$ for $i \neq j$ such that $LH(\mathcal{I}_i) \approx UH(\mathcal{I}_j)$.

Proof. According to (2.2), $\mathcal{I}_i \lhd \mathcal{I}_j$ if and only if $LH(\mathcal{I}_i) \lessdot UH(\mathcal{I}_j)$. If these is no such \mathcal{I}_i , then \mathcal{I}_i is minimal.

We can use this lemma to identify all minimal elements:

- If the ordering \leq starts with two lower handles $LH(\mathcal{I}_i)$ and $LH(\mathcal{I}_j)$, there exists no minimal element. The reason is that all upper handles are larger, and so both \mathcal{I}_i and \mathcal{I}_j are smaller than everything else; specifically, we get $\mathcal{I}_i \triangleleft \mathcal{I}_j \triangleleft \mathcal{I}_i$.
- If the first element of the ordering \leq is $LH(\mathcal{I}_i)$ then \mathcal{I}_i is the unique candidate for a minimal element. We just need to check whether there is some other $LH(\mathcal{I}_j)$ smaller than $UH(\mathcal{I}_i)$, and if so, no minimal element exists.²
- If \leq starts with a consecutive group of upper handles, we have several candidates for a minimal element. First, all \mathcal{I}_i 's of these upper handles are minimal elements. Second, if the lower handle following the group of upper handles is $LH(\mathcal{I}_j)$, then \mathcal{I}_j is a candidate for a minimal element. As above, \mathcal{I}_j is minimal if there is no other lower handle smaller than $UH(\mathcal{I}_j)$.

²This can be done in constant time if we remember in each moment positions of the two left-most lower handles in the ordering, and update this information after removing one of them from \leq .

For every topological sort, the ℓ -th element is minimal when restricted to the elements $\{\ell, \ell+1, \ldots, k\}$. So we can construct all topological sorts as follows. We repeatedly detect all minimal element \mathcal{I}_i and always pick one of them. (For different choices we get different topological sorts). We remove the handles of the picked minimal element \mathcal{I}_i from \approx and append \mathcal{I}_i to the constructed topological sort. We stop when all elements are placed in the topological sort. If in some step no minimal element exists, we know that no topological sort exists.

For a P-node, we just need to find any topological sort by repeated removing of minimal elements. For a Q-node, we test whether the current ordering or its reversal is a topological sort. We iterate through each of the prescribed sorts, check whether each element is a minimal element, and then removing its handles from $\tilde{\leq}$. In both cases, if we find a correct topological sort, we use it reorder the children of the node. Otherwise, the reordering is not possible and the algorithm outputs "no". We are able to do the reordering of the node in time $\mathcal{O}(k)$.

The Algorithm. Now, we are ready to show that our algorithm allow us to find a reordering of the PQ-tree T according to an interval order \triangleleft with a sorted representation in time $\mathcal{O}(e)$:

Proposition 2.5. We first deal with details of the implementation. First, we precompute the handles for every set of intervals corresponding to the subtree of every inner node of T. For each leaf, the two handles correspond for the two endpoints. We proceed the tree from bottom to the root. Suppose that we have an inner node corresponding to the set \mathcal{I} of intervals and it has k children corresponding to $\mathcal{I}_1, \ldots, \mathcal{I}_k$ for which we already know their handles. Then we can calculate handles of \mathcal{I} using

$$LH(\mathcal{I}) = \min\{LH(\mathcal{I}_i)\}$$
 and $UH(\mathcal{I}) = \max\{UH(\mathcal{I}_i)\}.$ (2.3)

This can clearly be computed in time $\mathcal{O}(e)$, and we also note for each endpoint a list of nodes for which it is a handle. Using these list, we can iterate the sorted representation and compute the orderings \approx for every inner node of T, again in $\mathcal{O}(e)$.

Now we test each inner node of T with the given ordering \leq whether its subtrees can be reordered according to \triangleleft . The algorithm is correct since it works in the same way as in Proposition 2.4, based on lemmas 2.6 and 2.7.

Concerning the time complexity, we already discussed that we are able to compare sets of intervals using handles in a constant time, by Lemma 2.6. The precomputation of all orderings \leq takes time $\mathcal{O}(e)$. We spend time $\mathcal{O}(k)$ in each node with k children. Thus the total time complexity of the algorithm is linear in the size of the tree, which is $\mathcal{O}(e)$.

For a pseudocode, see Algorithm 2 in Appendix A. We note that when the orderings \approx are constructed for all inner nodes, then we do not need to proceed the tree from bottom to the top. We can proceed them independently in parallel and a reordering T' of T exists if and only if we succeed in reordering of every inner node.

2.2 Extending Interval Graphs

In this section, we describe an algorithm solving Repext(INT) in time $\mathcal{O}(n+m)$ which uses Proposition 2.5 as a subroutine. Unlike in Section 2.1, the interval representations consist of closed intervals. We allow the intervals to share the endpoints and to have zero length. We first describe recognition of interval graphs. Then we show how to modify the PQ-tree approach to solve Repext(INT).

2.2.1 Recognition using PQ-trees

Recognition of interval graphs in linear time was a long-standing open problem, first solved by Booth and Lueker [8] using PQ-trees. Nowadays, there are two main approaches to recognition in linear time. The first one finds a feasible ordering of the maximal cliques which can be done using PQ-trees. The second approach uses surprising properties of the lexicographic breadth-first search, searches through the graph several times and constructs a representation if the graph is an interval graph [13].

We modify the PQ-tree approach to solve Repext(INT) in time $\mathcal{O}(n+m)$. Recall the PQ-trees from Section 2.1.

Maximal Cliques. The PQ-tree approach is based on the following characterization of interval graphs, due to Fulkerson and Gross [18]:

Lemma 2.8 (Fulkerson and Gross). A graph is an interval graph if and only if there exists an ordering of the maximal cliques such that for every vertex the cliques containing this vertex appear consecutively in this ordering.

Consider an interval representation of an interval graph. For each maximal clique, consider the intervals representing the vertices of this maximal clique and select a point in their intersection. (We know that this intersection is non-empty because intervals of the real line have the Helly property.) We call these points *clique-points*. For an illustration, see Figure 2.6. The ordering of the clique-points from left to right gives the ordering required by Lemma 2.8. Every vertex appears in consecutive maximal cliques since it is represented by an interval. For a maximal clique a, we denote the assigned clique-point by cp(a).

On the other hand, given an ordering of the maximal cliques, we place cliquepoints in this ordering on the real line. Each vertex is represented by the interval containing exactly the clique-points of the maximal cliques containing this vertex. In this way, we obtain a valid interval representation of the graph.

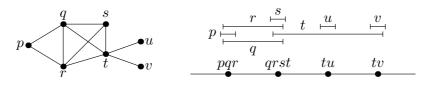


Figure 2.6: An interval graph and one of its representations with denoted clique-points.

Recognition Algorithm. Every chordal graph has at most $\mathcal{O}(n)$ maximal cliques of total size $\mathcal{O}(n+m)$ and they can be found in linear time [49]. Since every interval graph is chordal, we run this subroutine. If this subroutine fails, the input graph is not an interval graph, and the recognition algorithm outputs "no".

According to Lemma 2.8, we want to find an ordering of the maximal clique in which maximal clique containing each vertex appear consecutively. Recall the consecutive ordering problem from Section 2.1. Here, the elements E are the maximal cliques of the graph. For each vertex v, we introduce the restricting set S_v containing all the maximal cliques containing this vertex v. Using PQ-trees, we can find a feasible ordering of the maximal cliques and recognize an interval graph in time $\mathcal{O}(n+m)$.

2.2.2 Modification for REPEXT

We first sketch the algorithm. We construct a PQ-tree T for the input graph while ignoring the partial representation. The partial representation gives another restriction: an interval order \triangleleft of the maximal cliques. Using Proposition 2.5, we try to find a reordering T' of the PQ-tree T according to \triangleleft in time $\mathcal{O}(n+m)$. We are going to prove the following proposition: The partial representation is extendible if and only if the reordering algorithm succeeds.

Since our proof is constructive, we can use to build a representation \mathcal{R} extending the partial representation \mathcal{R}' . We place clique-points on the real line according to the ordering $<_{T'}$. We need to be more careful in this step. Since several intervals are pre-drawn, we cannot change their representations, so the clique-points has to be placed correctly. Using the clique-points, we construct the remaining intervals in a similar manner as in Figure 2.6.

Now, we describe everything in detail.

Restricting Clique-points. Suppose that there exists a representation \mathcal{R} extending \mathcal{R}' . Then \mathcal{R} gives some ordering < of the clique-points from left to right. We want to show that pre-drawn intervals partially specify position of some clique-points and give some necessary condition for <.

For a maximal clique a, let I(a) denote the set of all the pre-drawn intervals that are contained in a. Then I(a) restricts possible position of $\operatorname{cp}(a)$ to only those points x of the real line which are covered in \mathcal{R}' by exactly the pre-drawn intervals of I(a) and no others. We denote by (a) (resp. (a)) the leftmost (resp. the rightmost) point where the clique-point $\operatorname{cp}(a)$ can be placed, formally:

Notice that $(\curvearrowleft(a), \curvearrowright(a))$ is a subinterval of $\bigcap_{u \in I(a)} R'_u$. For an example, see Figure 2.7.

For a clique-point cp(a), the structure of all x where cp(a) can be placed is simple. The pre-drawn intervals split the line into several parts, traversed by the same intervals; denoted in Figure 2.7 by gray lines. A clique-point cp(a) can be placed only to those part which contain exactly the intervals of I(a) and no other pre-drawn intervals. So the set of all points x where cp(a) can be placed is a union of intervals.

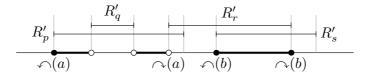


Figure 2.7: The partial representation \mathcal{R}' consisting of four pre-drawn intervals. Clique-points $\operatorname{cp}(a)$ and $\operatorname{cp}(b)$, having $I(a) = \{p\}$ and $I(b) = \{r, s\}$, can be placed to the bold parts of the real lines.

Notice also that the definition of (a) and (a) does not imply that cp(a) can be placed to all the points between (a) and (a). If a clique-point cannot be placed at all, the given partial representation is clearly not extendible.

The Interval Order \triangleleft . For two maximal cliques a and b, we define $a \triangleleft b$ if $\curvearrowright(a) \leq \curvearrowright(b)$. The definition of \triangleleft is quite natural since $a \triangleleft b$ implies that every extending representation \mathcal{R} has to place $\operatorname{cp}(a)$ to the left of $\operatorname{cp}(b)$. For example, the maximal cliques a and b in Figure 2.7 satisfy $a \triangleleft b$.

Lemma 2.9. The relation \triangleleft is an interval order.

Proof. The intervals representing \triangleleft correspond to the maximal cliques of G. To a maximal clique a, we assign an open interval $I_a = (\curvearrowleft(a), \curvearrowright(a))$. The definition of \triangleleft exactly states that $a \triangleleft b$ if and only if the intervals I_a and I_b are disjoint and I_a is on the left of I_b .

Lemma 2.10. For a sorted partial representation \mathcal{R}' , we can compute the sorted representations of \triangleleft in time $\mathcal{O}(n+m)$.

Proof. As stated above, our interval graph has $\mathcal{O}(n)$ maximal cliques containing in total $\mathcal{O}(n+m)$ vertices. We compute for every pre-drawn vertex the list of the maximal cliques containing it, and for every maximal clique a the number |I(a)| of pre-drawn vertices it contains. We initiate an empty list W and a counter i of pre-drawn intervals covering the currently sweeped point. We sweep the real line from left to right and compute the sorted representation of \triangleleft .

When sweeping there are two types of events. If we encounter a set of endpoints of pre-drawn intervals sharing a point, we first process the left endpoints, then we update \curvearrowleft and \curvearrowright for this point³, and then we process the right endpoints. If we sweep over a part, we just update \curvearrowright and \curvearrowright . We do in details the following:

- If we encounter a left endpoint ℓ_u , then we increase the counter i. For every clique a containing u, we increase its counter. If some clique a has all pre-drawn intervals placed over ℓ_u , we add a into the list W of watched cliques.
- If we encounter a right endpoint r_u , we decrease the counter of pre-drawn intervals. We ignore all maximal cliques a containing u till the end of the procedure, and naturally we also remove them from W if there are any.

³We need to update also here since it might happen that the interval (((a), (a))) is empty for some maximal clique a. This can happen only if some pre-drawn interval of I(a) is a singleton.

• Update of \curvearrowright and \curvearrowright is done for all cliques $a \in W$ such that |I(a)| = i. Notice that we currently sweep over exactly i pre-drawn intervals, and therefore we have to sweep over exactly the pre-drawn intervals of I(a). We update $\curvearrowright(a)$ to the left-most point of the current point/part if it is not yet initialized. And we update $\curvearrowright(a)$ to the right-most such point.

In the end, we output the computed \curvearrowleft and \curvearrowright naturally sorted from left to right. If for some maximal clique a, the value $\backsim(a)$ was not initiated, the clique-point $\operatorname{cp}(a)$ cannot be placed and the procedure outputs "no".

The procedure is clearly correct, and it remains to argue that this can be done in linear time. We have the cliques in W partitioned according to |I(a)|. Also notice that when we sweep over i pre-drawn intervals, then there is no $a \in W$ such that |I(a)| > i and if $a, b \in W$ such that |I(a)| = |I(b)| = i, then necessarily I(a) = I(b). But then (a) = (b) and (a) = (b), so we can ignore b for the rest of the sweep procedure and set in the end the value according to the clique a. This implementation clearly runs in O(n+m).

Using Lemma 2.10, we can construct the sorted representation of \triangleleft in time $\mathcal{O}(n+m)$. We can test Reorder (T, \triangleleft) in time $\mathcal{O}(n+m)$ using Proposition 2.5. The following proposition is key for the correctness of the algorithm.

Proposition 2.11. A partial representation \mathcal{R}' is extendible if and only if G is an interval graph and the problem REORDER (T, \triangleleft) can be solved.

Proof. The condition forced by \triangleleft are clearly necessary. So if \mathcal{R}' is extendible, then the both condition have to be satisfied. It remain to show the other implication.

Since G is an interval graph, there exists a PQ-tree T representing all feasible orderings of the maximal cliques. There exists a reordering T' of T according to \triangleleft , and we denote $\triangleleft_{T'}$ as \triangleleft . We construct a representation \mathcal{R} extending \mathcal{R}' as follows. We place the clique-points according to \triangleleft from left to right, always greedily as far to the left as possible.

Suppose we want to place a clique-point $\operatorname{cp}(a)$. Let $\operatorname{cp}(b)$ be the last placed clique-point. Consider the infimum over all the points where the clique-point $\operatorname{cp}(a)$ can be placed and that are to the right of the clique-point $\operatorname{cp}(b)$. If there is a single such point on the right of $\operatorname{cp}(b)$ (equal to the infimum), we place $\operatorname{cp}(a)$ there. Otherwise $\operatorname{c}(a) < \operatorname{c}(a)$ and we place the clique-point $\operatorname{cp}(a)$ to the right of this infimum by an appropriate epsilon, for example the length of the shortest part (see definition of $\operatorname{c}(a)$) divided by n.

We prove by contradiction that this greedy procedure cannot fail; see Figure 2.8. Let $\operatorname{cp}(a)$ be the clique-point for which the procedure fails. Since $\operatorname{cp}(a)$ cannot be placed, there are some clique-points placed on the right of \curvearrowright (a) (or possibly on \curvearrowright (a) directly). Let $\operatorname{cp}(b)$ be the leftmost one of them. If \curvearrowright (b) $\ge \curvearrowright$ (a), we obtain $a \lhd b$ which contradicts b < a since $\operatorname{cp}(b)$ was placed before $\operatorname{cp}(a)$. So, we know that \curvearrowright (b) $< \curvearrowright$ (a). To get contradiction, we question why the clique-point $\operatorname{cp}(b)$ was not placed on the left of \curvearrowright (a).

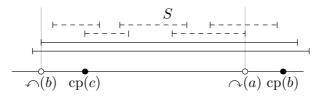


Figure 2.8: An illustration of the proof: The positions of the clique-points cp(b) and cp(c), the intervals of S are dashed.

The clique-point $\operatorname{cp}(b)$ was not placed before \curvearrowright (a) because all these positions were either blocked by some other previously placed clique-points, or they are traversed by some pre-drawn interval not in I(b). There is at least one clique-point placed to the right of \curvearrowright (b) (otherwise we could place $\operatorname{cp}(b)$ to \leadsto (b) or right next to it). Let $\operatorname{cp}(c)$ be the right-most clique-point placed between \leadsto (b) and $\operatorname{cp}(b)$. Every point between $\operatorname{cp}(c)$ and \curvearrowright (a) has to be covered by a pre-drawn interval not in I(b). Consider the set S of all the pre-drawn intervals not contained in I(b) covering any point of the real line in $[c, \curvearrowright$ (a)]; depicted dashed in Figure 2.8.

Let C be a set of all the cliques containing at least one vertex from S. Since S induces a connected subgraph, all the cliques of C appear consecutively in < because every pair of adjacent vertices from S is contained in some maximal clique of C.

Now, a and c both belong to C, but b does not. We assumed that c < b < a. Since c < b and the consecutivity of C, then a < b which contradicts b < a.

The Algorithm. We proceed in the following five steps. Only the first three steps are necessary, if we just want to answer the decision problem without constructing a representation.

- 1. Independently of the partial representation, find the maximal cliques and construct a PQ-tree T representing all feasible orderings of the maximal cliques.
- 2. Construct the sorted representation of the interval order \triangleleft by Lemma 2.10.
- 3. Using Proposition 2.5, test whether there is a reordering T' of the PQ-tree T according to \triangleleft .
- 4. Place the clique-points from left to right according to $<_{T'}$ on the real line, greedily as far to the left as possible.
- 5. Using these clique-points, construct a representation \mathcal{R} extending \mathcal{R}' .

Step 1 is the original recognition algorithm. In Step 2, we compute splitting of the real line into parts and construct a sorted representation of \triangleleft . In Step 3, we apply the algorithm of Proposition 2.5. Step 4 is the greedy procedure from the proof of Proposition 2.11. In Step 5, we construct intervals representing the vertices of $G \setminus G'$ as in Figure 2.6; we construct each such interval on top of the corresponding clique-points. See Algorithm 3 in Appendix A for a pseudocode.

Now we are ready to prove the main result of this chapter, Theorem 1.1 which states that the problem REPEXT(INT) can be solved in time $\mathcal{O}(n+m)$:

Theorem 1.1. The correcness of the algorithm is implied by Proposition 2.11. Concerning the complexity, the total size of all maximal cliques is at most $\mathcal{O}(n+m)$ and

the PQ-tree can be constructed in this time. Using Lemma 2.10, we can construct the sorted representation of \triangleleft in time $\mathcal{O}(n+m)$. According to Proposition 2.5, the PQ-tree can be reordered according to \triangleleft in time $\mathcal{O}(n+m)$. Finally, a representation \mathcal{R} extending \mathcal{R}' can be constructed, if necessary, in time $\mathcal{O}(n+m)$.

Extending Proper and Unit Interval Graphs

In this chapter, we extend representations of proper and unit interval graphs. As was already described in the introduction, even through these classes are known to be equal, they behave differently for the partial representation extension problem.

In Section 3.1, we deal with proper interval graphs and the extension algorithm is based on a simple characterization of extendible instances. In Section 3.2, we study a more general problem of bounded representations of unit interval graphs and we show that this problem is NP-complete. In Section 3.3, we deal with specific instances of the bounded representation problem for which the left-to-right order of the components is prescribed, and give an almost quadratic-time combinatorial algorithm for this problem. In Section 3.4, we solve the Repext(UNIT INT) problem by showing that it can be expressed as an instance of the bounded representation problem with prescribed ordering.

3.1 Extending Proper Interval Representations

In this section, we describe how to extend partial representations of proper interval graphs in time $\mathcal{O}(m+n)$. We also give a simple characterization of all extendible instances.

Indistinguishable Vertices. Vertices u and v are called indistinguishable if N[u] = N[v]. The vertices of G can be partitioned into groups of (pairwise) indistinguishable vertices. Note that indistinguishable vertices may be represented by the same intervals (and this is actually true for general intersection representations). Since indistinguishable vertices are not very interesting from the structural point of view, if the structure of the pre-drawn vertices allows it, we want to prune the graph to keep only one vertex per group.

Suppose that we are given an instance of Repext(PROPER INT). We compute the groups of indistinguishable vertices in time $\mathcal{O}(n+m)$ using the algorithm of Rose et. al [49]. Let u and v be two indistinguishable vertices. If u is not pre-drawn, or both vertices are pre-drawn with $R'_u = R'_v$, then we remove u from the graph, and in the final constructed representation (if it exists) we put $R_u = R_v$. For the rest of the section, we shall assume that the input graph and partial representation are pruned.

Figure 3.1: Two proper interval representations \mathcal{R}_1 and \mathcal{R}_2 with the left-to-right orderings $v_1 \triangleleft v_2 \triangleleft v_3 \triangleleft v_4 \triangleleft v_5 \triangleleft v_6 \triangleleft v_7 \triangleleft v_8$ and $v_2 \triangleleft v_1 \triangleleft v_3 \triangleleft v_4 \triangleleft v_5 \triangleleft v_7 \triangleleft v_6 \triangleleft v_8$.

An important property is that for any representation of a pruned graph, it holds that all intervals are pairwise distinct.

Left-to-right ordering. Roberts [46] gave the following characterization of proper interval graphs:

Lemma 3.1 (Roberts). A graph is a proper interval graph if and only if there exists a linear ordering $v_1 \triangleleft v_2 \triangleleft \cdots \triangleleft v_n$ of its vertices such that the closed neighborhood of every vertex is consecutive.

This linear order \triangleleft corresponds to the left-to-right order of the intervals on the real line in some valid representation of the graph. On the other hand, in each representation the order of the left endpoints is exactly the same as the order of the right endpoints, and this order satisfies the condition of Lemma 3.1. This makes recognition much simpler. For an example of \triangleleft , see Figure 3.1.

How many different orderings \triangleleft may a proper interval graph admit? In the case of a general unpruned graph possibly many, but all of them have a very simple structure. In Figure 3.1, the graph contains two groups $\{v_1, v_2, v_3\}$ and $\{v_6, v_7\}$. The vertices of each group have to appear consecutively in the ordering \triangleleft and may be reordered arbitrarily. Deng et al. [14] proved the following:

Lemma 3.2 (Deng et al.). For a connected (unpruned) proper interval graph, the ordering \triangleleft satisfying the condition of Lemma 3.1 is uniquely determined up to local reordering of groups and complete reversal.

This lemma is key for partial representation extension of proper interval graphs. Essentially, we just have to deal with a unique ordering (and its reversal) and match the partial representation on it. Notice that in the pruned graph, if two vertices are indistinguishable, then their order is prescribed by the partial representation.

Ignore the partial representation for a second. We want to construct a partial ordering < which is a simple representation of all orderings < from Lemma 3.1. There exists a proper interval representation with an ordering < if and only if < extends either < or its reversal. According to Lemma 3.2, < can be constructed by taking an arbitrary ordering < and making indistinguishable vertices incomparable. For the graph in Figure 3.1, we get

$$(v_1, v_2, v_3) < v_4 < v_5 < (v_6, v_7) < v_8,$$

where groups of indistinguishable vertices are put in brackets. This ordering is unique up to reversal and can be constructed in time $\mathcal{O}(n+m)$ [12].

Characterization of Extendible Instances. We give a simple characterization of the partial representation instances that are extendible. We start with connected instances. Let G be a pruned proper interval graph and \mathcal{R}' be a partial representation of its induced subgraph G'. Then intervals in \mathcal{R}' are in some left-to-right ordering $<^{G'}$. (Recall that the pre-drawn intervals are pairwise distinct.)

Lemma 3.3. The partial representation \mathcal{R}' of a connected graph G is extendible if and only if there exists a linear ordering \triangleleft of V(G) such that:

- (1) The ordering \triangleleft extends \triangleleft extends \triangleleft and either \triangleleft or its reversal.
- (2) Let R'_u and R'_v be two pre-drawn touching intervals, i.e., $r_u = \ell_v$, and let w be any vertex distinct from u and v. If $uw \in E(G)$, then $w \triangleleft v$, and if $vw \in E(G)$, then $u \triangleleft w$.

Proof. If there exists a representation \mathcal{R} extending \mathcal{R}' , then it is in some left-to-right ordering \triangleleft . Clearly, the pre-drawn intervals are placed the same, so \triangleleft has to extend \triangleleft According to Lemma 3.2, \triangleleft extends \triangleleft or its reversal. As for (2), clearly v has to be the right-most neighbor of u in \mathcal{R} : if R_w was on the right of R_v , it would not intersect R_u . Similarly. u is the left-most vertex of v.

Conversely, let $v_1 \triangleleft \cdots \triangleleft v_n$ be an ordering from the statement of the lemma. We construct a representation \mathcal{R} extending \mathcal{R}' as follows. We compute a common linear ordering \triangleleft of left and right endpoints from left-to-right. We start with the ordering $\ell_1 \triangleleft \cdots \triangleleft \ell_n$, into which we insert the right endpoints r_1, \ldots, r_n one-by-one. For vertex v_i , let v_j be its right-most neighbor in the ordering \triangleleft . Then, we place r_i right before ℓ_{j+1} (if it exists, otherwise we append it to the end of the ordering).

This left-to-right common order < is uniquely determined by <. Since < extends $<^{G'}$, it is compatible with the partial representation (the pre-drawn endpoints are ordered as in <). To construct the representation, we just place the non-pre-drawn endpoints equidistantly into the gaps between neighboring pre-drawn endpoints (or to the left or right of \mathcal{R}'). It is important that, if two pre-drawn endpoints ℓ_i and r_j share their position, then according to condition (2) there is no endpoint placed in between of ℓ_i and r_j in < (otherwise one of the two implications would not hold, depending whether a left endpoint is intersected in between, or a right one). See Figure 3.2 for an example.

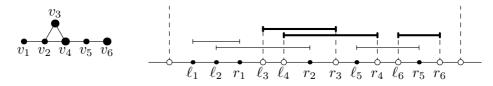


Figure 3.2: Representation of a component with order $v_1 \triangleleft v_2 \triangleleft v_3 \triangleleft v_4 \triangleleft v_5 \triangleleft v_6$. First, we compute the common order of the left and right endpoints: $\ell_1 \lessdot \ell_2 \lessdot r_1 \lessdot \ell_3 \lessdot \ell_4 \lessdot r_2 \lessdot r_3 \lessdot \ell_5 \lessdot r_4 \lessdot \ell_6 \lessdot r_5 \lessdot r_6$. The endpoints of the pre-drawn intervals split the segment into several subsegments. We place the remaining endpoints in this order and, within every subsegment, distributed equidistantly.

¹Notice that, in the partial representation, some intervals may share position. But if two endpoints ℓ_i and r_j share the position, then $v_i v_j \in E(G)$ and we break the tie by setting $\ell_i \lessdot r_j$.

We argue correctness of the constructed representation \mathcal{R} . First, it extends \mathcal{R}' , since the pre-drawn intervals are not modified. Second, it is a correct interval representation: Let v_i and v_j be two vertices with $v_i \triangleleft v_j$, and let v_k be the right-most neighbor of v_i in \triangleleft . If $v_i v_j \in E(G)$, then $\ell_i \lessdot \ell_k \lessdot r_i$ and, by consecutivity of N[u] in \triangleleft , we have $\ell_j \lessdot \ell_k$. Therefore, R_{v_i} and R_{v_j} intersect. If $v_i v_j \notin E(G)$ and $v_j \neq v_{k+1}$, then $r_i \lessdot \ell_{k+1} \lessdot \ell_j$, so R_{v_i} and R_{v_j} do not intersect. If $v_i v_j \notin E(G)$ and $v_j = v_{k+1}$, then $r_i \lessdot \ell_{k+1}$ and R_{v_i} and R_{v_j} do not intersect. Finally, we argue that \mathcal{R} is a proper interval representation. In \lessdot the order of the left endpoints is the same as the order of the right-endpoints, since r_{i+1} is always placed on the right of r_i in \lessdot .

We conclude that the representation \mathcal{R} can be made small enough to fit into any open segment of the real line that contains all pre-drawn intervals.

Now, we are ready to characterize general solvable instances.

Lemma 3.4. A partial representation \mathcal{R}' of a graph G is extendible if and only if

- (1) for each component C the partial representation \mathcal{R}'_C consisting of pre-drawn intervals in C is extendible, and
- (2) pre-drawn vertices of each component are consecutive in $<^{G'}$.

Proof. The necessity of (1) is due to Lemma 3.3 applied on each component C. For (2), if some component C would not have its pre-drawn vertices consecutive in $<^{G'}$, then $\bigcup_{u \in C} R_u$ would not be a connected segment of the real line (contradicting existence of \blacktriangleleft from Section 1.5).

Now, if the instance satisfies the both conditions we can construct a correct representation \mathcal{R} extending \mathcal{R}' as follows. Using (2), the located components are ordered from left to right, and we assign pairwise disjoint *open segments* containing all their pre-drawn intervals (there is a non-empty gap between located components we can use). To unlocated components, we assign pairwise disjoint open segments to the right of the right-most located component. See Figure 3.3. For each component, we construct a representation in its open segment, using the construction in the proof of Lemma 3.3.

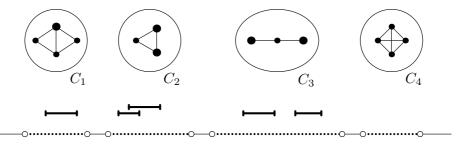


Figure 3.3: An example of a graph with four components C_1, \ldots, C_4 . The pre-drawn intervals give the order of the located components $C_1 \blacktriangleleft C_2 \blacktriangleleft C_3$. The non-located component C_4 is placed to the right. For each component, we reserve some segment in which we construct the representation.

It remains to prove that the problem RepExt(PROPER INT) can be solved in time $\mathcal{O}(n+m)$:

Proof of Theorem 1.2. We just use the characterization by Lemma 3.4, of which the conditions (1) and (2) can be easily checked in time $\mathcal{O}(n+m)$. For Lemma 3.3, we check for each component the both constraints (1) and (2). To check (2), we compute for < and its reversal the unique orderings \triangleleft . We test for each of them whether each touching pair of pre-drawn intervals is placed in \triangleleft according to (2).

If necessary, a representation \mathcal{R} can be constructed in the same running time since the proofs of Lemmas 3.3 and 3.4 are constructive.

For a pseudocode, see Algorithm 4 in Appendix A.

3.2 Bounded Representations of Unit Interval Graphs

In this section we deal with bounded representations. An input of BOUNDREP consists of a graph G and, for each vertex v_i , a lower bound lbound(v_i) and an upper bound ubound(v_i). (We allow lbound(v_i) = $-\infty$ and ubound(v_i) = $+\infty$.) The problem asks whether there exists a unit interval representation \mathcal{R} of G such that lbound(v_i) $\leq \ell_i \leq \text{ubound}(v_i)$ for each interval v_i . Such a representation is called a bounded representation.

Since unit interval representations are proper interval representations, all properties of proper interval representations described in Section 3.1 hold, in particular properties of orderings \triangleleft and \triangleleft .

3.2.1 Representations in ε -grids

Endpoints of intervals can be positioned at arbitrary real numbers. For the purpose of the algorithm, we want to work with drawings of limited resolution. For a given instance of the bounded representation problem, we want to find a lower bound for the required resolution such that this instance is solvable if and only if it is solvable in this limited resolution.

More precisely, we want to represent all intervals so that their endpoints correspond to points on some grid. For a value $\varepsilon = \frac{1}{K} > 0$, where K is an integer, the ε -grid is the set of points $\{k\varepsilon : k \in \mathbb{Z}\}$. For a given instance of BOUNDREP, we ask which value of ε ensures that we can construct a representation having all endpoints on the ε -grid. So the value of ε is the resolution of the drawing.

For the standard unit interval graph representation problem a grid of size $\frac{1}{n}$ is sufficient [12]. In the case of BOUNDREP, the size of the grid has to depend on the values of the bounds. Consider all values lbound(v_i) and ubound(v_i) distinct from

²If ε was not of the form $\frac{1}{K}$, then the grid could not contain both left and right endpoints of the intervals. We reserve K for the value $\frac{1}{\varepsilon}$ in this chapter.

 $-\infty, +\infty$, and express them as irreducible fractions $\frac{p_1}{q_1}, \frac{p_2}{q_2}, \cdots, \frac{p_b}{q_b}$. Then we define:

$$\varepsilon' := \frac{1}{\operatorname{lcm}(q_1, q_2, \dots, q_b)}, \quad \text{and} \quad \varepsilon := \frac{\varepsilon'}{n}.$$
 (3.1)

It is important that the size of this ε written in binary is $\mathcal{O}(r)$. We show that the ε -grid is sufficient to construct the bounded representation:

Lemma 3.5. If there exists a valid representation \mathcal{R}' for an input of the problem BOUNDREP, there exists a valid representation \mathcal{R} in which all intervals have endpoints on the ε -grid, where ε is defined by (3.1).

Proof. We construct an ε -grid representation \mathcal{R} from \mathcal{R}' in two steps. First, we shift intervals to the left, and then we shift intervals slightly back to the right. For every interval v_i , the sizes of the left and right shifts are denoted by $LS(v_i)$ and $RS(v_i)$ respectively. The shifting process is shown in Figure 3.4.

In the first step, we consider the ε' -grid and shift all the intervals to the left to the closest grid-point (we do not shift an interval if its endpoints are already on the grid). Original intersections are kept by this shifting, since if x and y are two endpoints satisfying $x \leq y$ before the left-shift, then $x \leq y$ also holds after the left-shift. So if $v_i v_j \in E$ and $\ell_i \leq \ell_j \leq r_i$ before the shift, then these inequalities are preserved by the shifting. On the other hand, we may introduce additional intersections by shifting two non-intersecting intervals to each other. In this case, after the shift the intervals only touch; for an example, see vertices v_2 and v_4 in Figure 3.4.

The second step shifts the intervals to the right in the refined ε -grid to remove the additional intersections created by the first step. The right-shift is a mapping

RS:
$$\{v_1, \ldots, v_n\} \to \{0, \varepsilon, 2\varepsilon, \ldots, (n-1)\varepsilon\}$$

having the right-shift property: For all pairs (v_i, v_j) with $r_i = \ell_j$, $RS(v_i) \ge RS(v_j)$ if and only if $v_i v_j \in E$.

To construct such a mapping RS, notice that if we relaxed the image of RS to $[0, \varepsilon')$, the reversal of LS would have the right-shift property, since it produces the original correct representation \mathcal{R}' . But the right-shift property depends only on the relative order of the shifts and not on the precise values. Therefore, we can construct RS from the reversal of LS by keeping the shifts in the same relative order. If $LS(v_i)$ is one of the kth smallest shifts, we set $RS(v_i) = (k-1)\varepsilon$.³ See Figure 3.4.

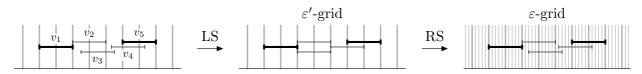


Figure 3.4: In the first step, we shift intervals to the left to the ε' -grid. The left shifts of v_1, \ldots, v_5 are $(0, 0, \frac{1}{2}\varepsilon', \frac{1}{3}\varepsilon', 0)$. In the second step, we shift to the right in the refined ε -grid. Right shifts have the same relative order as left shifts: $(0, 0, 2\varepsilon, \varepsilon, 0)$.

³In other words, for the smallest shifts we assign the right-shift 0; for the second smallest shifts, we assign ε ; for the third smallest shifts, 2ε ; and so on.

We finally argue that these shifts produce a correct ε -grid representation. The right-shift does not create additional intersections: After LS non-intersecting pairs are at distance at least $\varepsilon' = n\varepsilon$, and by RS they can get closer by at most $(n-1)\varepsilon$. Also, if after LS two intervals overlap by at least ε' , their intersection is not removed by RS. The only intersections which are modified by RS are touching pairs of intervals (v_i, v_j) having $r_i = \ell_j$ after LS. The mapping RS shifts these pairs correctly according to the edges of the graph.

Next we look at the bound constraints. If, before the shifting, v_i was satisfying $\ell_i \geq \text{lbound}(v_i)$, then this is also satisfied after $LS(v_i)$ since the ε' -grid contains the value lbound (v_i) . Obviously, the inequality is not broken after $RS(v_i)$. As for the upper bound, if $LS(v_i) = 0$ and $RS(v_i) = 0$, then the bound is trivially satisfied. Otherwise, after $LS(v_i)$ we have $\ell_i \leq \text{ubound}(v_i) - \varepsilon'$, so the upper bound still holds after $RS(v_i)$.

Additionally, Lemma 3.5 shows that it is always possible to construct an ε -grid representation having the same topology as the original representation, in the sense that overlapping pairs of intervals keep overlapping, and touching pairs of intervals keep touching. Also notice that both representations \mathcal{R} and \mathcal{R}' have the same order of the intervals.

In the standard unit interval graph representation problem, no bounds on the positions of the intervals are given, and we get $\varepsilon' = 1$ and $\varepsilon = \frac{1}{n}$. Lemma 3.5 proves in a particularly clean way that the grid of size $\frac{1}{n}$ is sufficient to construct unrestricted representations of unit interval graphs. Corneil et al. [12] show how to construct this representation directly from the ordering <, whereas we use some given representation to construct an ε -grid representation.

3.2.2 The Hardness of BOUNDREP

In this subsection we focus on hardness of bounded representations of unit interval graphs. We prove Theorem 1.3 stating that BOUNDREP is NP-complete.

We reduce the problem from 3-Partition. An input of 3-Partition consists of natural numbers k, M, and A_1, \ldots, A_{3k} such that $\frac{M}{4} < A_i < \frac{M}{2}$ for all i, and $\sum A_i = kM$. The question is whether it is possible to partition the numbers A_i into k triples such that each triple sums to exactly M. This problem is known to be strongly NP-complete (even if all numbers have polynomial sizes) [20].

Proof of Theorem 1.3. According to Lemma 3.5, if there exists a representation satisfying the bound constraints, there also exists an ε -grid representation with this property. Since the length of ε given by (3.1), written in binary, is polynomial in the size of the input, all endpoints can be placed in polynomially-long positions. Thus we can guess the bounded representation and the problem belongs to NP.

Next, we prove that the problem is NP-hard. For a given input of 3-Partition, we construct the following unit interval graph G. For each number A_i , we add a path P_{2A_i} (of length $2A_i - 1$) into G as a separate component. For all vertices x in these

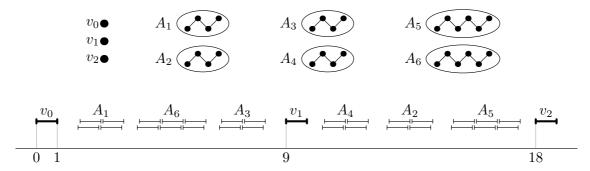


Figure 3.5: We consider the following input for 3-Partition: $k=2, M=7, A_1=A_2=A_3=A_4=2$ and $A_5=A_6=3$. The associated unit interval graph is depicted on top, and at the bottom we find one of its correct bounded representations, giving 3-partitioning $\{A_1, A_3, A_6\}$ and $\{A_2, A_4, A_5\}$.

paths, we set bounds

$$lbound(x) = 1$$
 and $ubound(x) = k \cdot (M+2)$.

In addition, we add k+1 independent vertices v_0, v_1, \ldots, v_k and make their positions in the representation fixed:

$$lbound(v_i) = ubound(v_i) = i \cdot (M+2).$$

See Figure 3.5 for an illustration of the reduction. Clearly, the reduction is polynomial.

We now argue that the bounded representation problem is solvable if and only if the given input of 3-Partition is solvable. Suppose first that the bounded representation problem admits a solution. There are k gaps between the fixed intervals v_0, \ldots, v_k each of which has space less than M+1. (The length of the gap is M+1 but the endpoints are taken by v_i and v_{i+1} .) The bounds of the paths force their representations to be inside these gaps, and each path lives in exactly one gap. Hence the representation induces a partition of the paths.

Now, the path P_{2A_i} needs space at least A_i in every representation. The representations of the paths may not overlap and the space in each gap is less than M+1, hence the sum of all A_i 's in each part is at most M. Since the total sum of A_i 's is exactly kM, the sum in each part has to be M. Thus the obtained partition solves the 3-Partition problem.

Conversely, every solution of 3-Partition can be realized in this way. \Box

3.3 Bounded Representations with Prescribed Ordering

In this section, we deal with the BOUNDREP problem when a fixed ordering ◀ of the components is prescribed. First we solve the problem using linear programming. Then we describe additional structure of bounded representations, and using this structure we construct an almost quadratic-time algorithm that solves the linear programs.

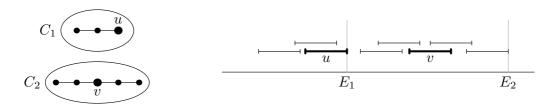


Figure 3.6: The positions of the vertices u and v are fixed by the bound constraints. The component C_1 can only be represented with u being the right-most interval, since otherwise it would block space for the component C_2 .

3.3.1 LP Approach for BOUNDREP

According to Lemma 3.2, each component of G can be represented in at most two different ways, up to local reordering of groups of indistinguishable vertices. Unlike the case of proper interval graphs, we cannot arbitrarily choose one of the orderings, since neighboring components restrict each other's space. For example, only one of the two orderings for the component C_1 in Figure 3.6 makes a representation of C_2 possible.

In the algorithm, we process components $C_1 \triangleleft C_2 \triangleleft \cdots \triangleleft C_c$ from left to right and construct representations for them. When we process a component C_t , we want to represent in on the right of the previous component C_{t-1} , and we want to push the representation of C_t as far to the left as possible, leaving as much space for C_{t+1}, \ldots, C_c as possible.

Now, we describe in details, how we process a component C_t . We calculate by the algorithm of Corneil et al. the partial ordering < described in Section 3.1 and its reversal. The elements that are incomparable by these partial orderings are vertices of the same group of indistinguishable vertices. For these vertices, the following holds:

Lemma 3.6. Suppose there exists some bounded representation \mathcal{R} . Then there exists a bounded representation \mathcal{R}' such that, for every indistinguishable pair v_i and v_j satisfying lbound $(v_i) \leq \text{lbound}(v_j)$, it holds that $\ell'_i \leq \ell'_j$.

Proof. Given a representation \mathcal{R} , we call a pair (v_i, v_j) bad if v_i and v_j are indistinguishable, lbound $(v_i) \leq \text{lbound}(v_j)$ and $\ell_i > \ell_j$. We describe a process which iteratively constructs \mathcal{R}' from \mathcal{R} , by constructing a sequence of representations $\mathcal{R} = \mathcal{R}_0, \mathcal{R}_1, \ldots, \mathcal{R}_k = \mathcal{R}'$, where the positions in a representation \mathcal{R}_s are denoted by ℓ_i^s 's.

In each step s, we create \mathcal{R}_s from \mathcal{R}_{s-1} by fixing one bad pair (v_i, v_j) : we set $\ell_i^s = \ell_j^{s-1}$ and the rest of the representation remains the same. Since v_i and v_j are indistinguishable and \mathcal{R}_{s-1} is correct, the obtained \mathcal{R}_s is a representation. Regarding bound constraints,

$$lbound(v_i) \le lbound(v_j) \le \ell_j^{s-1} = \ell_i^s < \ell_i^{s-1} \le ubound(v_i),$$

so the bounds of v_i are satisfied.

Now, in each \mathcal{R}_s the set of all left endpoints is a subset of the set of all left endpoints of \mathcal{R} . In each step, we move one left-endpoint to the left, so each endpoint is moved at most n-1 times. Hence the process terminates after $\mathcal{O}(n^2)$ iterations and produces a representation \mathcal{R}' without bad pairs as requested.

For < and its reversal, we use Lemma 3.6 to construct linear orderings \lhd : If v_i and v_j belong to the same group of indistinguishable vertices and lbound $(v_i) <$ lbound (v_j) , then $v_i \lhd v_j$. If lbound $(v_i) =$ lbound (v_i) , we choose any order \lhd between v_i and v_j .

We obtain two total orderings \triangleleft , and we solve a linear program for each of them. Let $v_1 \triangleleft v_2 \triangleleft \cdots \triangleleft v_k$ be one of these orderings. We denote the right-most endpoint of a representation of a component C_t by E_t . Additionally, we define $E_0 = -\infty$. Also, we modify all lower bounds by putting $\text{lbound}(v_i) = \max\{\text{lbound}(v_i), E_{t-1} + \varepsilon\}$ for every interval v_i , which forces the representation of C_t to be on the right of the previously constructed representation of C_{t-1} . The linear program has variables ℓ_1, \ldots, ℓ_k , and it minimizes the value of E_t . Let ε be defined as in (3.1). We solve:

Minimize:
$$E_t := \ell_k + 1$$
,
subject to: $\ell_i \leq \ell_{i+1}$, $\forall i = 1, \dots, k - 1$, (3.2)
 $\ell_i \geq \text{lbound}(v_i)$, $\forall i = 1, \dots, k$, (3.3)
 $\ell_i \leq \text{ubound}(v_i)$, $\forall i = 1, \dots, k$, (3.4)
 $\ell_i \geq \ell_j - 1$, $\forall v_i v_j \in E, v_i \lhd v_j$, (3.5)
 $\ell_i + \varepsilon \leq \ell_j - 1$, $\forall v_i v_j \notin E, v_i \lhd v_j$. (3.6)

We solve the same linear program for the other ordering of the vertices of C_t . If none of the two programs is feasible, we report that no bounded representation exists. If exactly one of them is feasible, we keep the values obtained for ℓ_1, \ldots, ℓ_k and E_t , and process the next component C_{t+1} . If the two problems are feasible, we keep the one such that the value of E_t in the solution is smaller, and process C_{t+1} .

Lemma 3.7. Let the representation of C_{t-1} be fixed. Every bounded ε -grid representation of the component C_t with the left-to-right order $v_1 \triangleleft \cdots \triangleleft v_k$ and which is on the right of the representation of C_{t-1} satisfies constraints (3.3)–(3.6).

Proof. Constraints of types (3.3) and (3.4) are satisfied, since the representation is bounded and on the right of C_{t-1} . Constraints of type (3.5) correspond to correct representation of intersecting pairs of intervals. The non-intersecting pairs of an ε -grid representation are at distance at least ε , which makes constraints of type (3.6) satisfied.

Now, we are ready to show:

Proposition 3.8. The BOUNDREP problem with prescribed \triangleleft can be solved in polynomial time.

Proof. Concerning the running time, it depends polynomially on the sizes of n and ε , which are polynomial in the size of the input r. It remains to show correctness.

Suppose that the algorithm returns a candidate for a bounded representation. The formulation of the linear program ensures that it is a correct representation: Constraints of type (3.2) make the representation respect \triangleleft . Constraints of type (3.3) and (3.4) enforce that the given lower and upper bounds for the positions of the intervals are satisfied, force the prescribed ordering \blacktriangleleft on the representation of G, and

force the drawings of the distinct components to be disjoint. Finally, constraints of type (3.2), (3.5) and (3.6) make the drawing of the vertices of a particular component C_t be a correct representation.

Suppose next that a bounded representation exists. According to Lemma 3.5 and Lemma 3.6, there also exists an ε -grid bounded representation \mathcal{R}' having the order in the indistinguishable groups as defined above. So for each component C_t , one of the two orderings \triangleleft constructed for the linear programs agrees with the left-to-right order of C_t in \mathcal{R}' .

We want to show that the representation of each component C_t in \mathcal{R}' gives a solution to one of the two linear programs associated to C_t . We denote by E'_t the value of E_t in the representation \mathcal{R}' , and by E_t^{\min} the value of E_t obtained by the algorithm after solving the two linear programming problems associated to C_t . We show by induction on t that $E_t^{\min} \leq E'_t$, which specifically implies that E_t^{\min} exists and at least one of the linear programs for C_t is solvable.

We start with C_1 . As argued above, the left-to-right order in \mathcal{R}' agrees with one of the orderings \triangleleft , so the representation of C_1 satisfies the constraints (3.2). Since $E_0 = -\infty$, the lower bounds are not modified. By Lemma 3.7, the rest of the constraints are also satisfied. Thus the representation of C_1 gives a feasible solution for the program and gives $E_1^{\min} \leq E_1'$.

Assume now that, for some C_t with $t \geq 1$, at least one of the two linear programming problems associated to C_t admits a solution, and from induction hypothesis we have $E_t^{\min} \leq E_t'$. In \mathcal{R}' , two neighboring components are represented at distance at least ε . Therefore for every vertex v_i of C_{t+1} , it holds $\ell_i \geq E_t' + \varepsilon \geq E_t^{\min} + \varepsilon$, so the modification of the lower bound constraints is satisfied by \mathcal{R}' . Similarly as above using Lemma 3.7, the representation of C_{t+1} in \mathcal{R}' satisfies the remaining constraints. It gives some solution to one of the programs and we get $E_{t+1}^{\min} \leq E_{t+1}'$.

In summary, if there exists a bounded representation, for each component C_t at least one of the two linear programming problems associated to C_t admits a solution. Therefore, the algorithm returns a correct bounded representation \mathcal{R} (as discussed in the beginning of the proof). We note that \mathcal{R} does not have to be an ε -grid representation since the linear program just states that non-intersecting intervals are at distance at least ε . To construct an ε -grid representation if necessary, we can proceed as in the proof of Lemma 3.5.

For a pseudocode, see Algorithm 5 in Appendix A.

We note that it is possible to reduce the number of constraints of the linear program from $\mathcal{O}(k^2)$ to $\mathcal{O}(k)$. Using the ordering constraints (3.2), we can replace the groups of constraints (3.5) and (3.6) by a linear number of constraints as follows. For each v_j , there are two cases. If v_j is adjacent to all vertices v_i such that $v_i \triangleleft v_j$, then we only state the constraint (3.5) for v_1 and v_j . Otherwise, let v_i be the rightmost vertex such that $v_i \triangleleft v_j$ and $v_i v_j \notin E$. Then we only state the constraint (3.5) for v_{i+1} and v_j , and the constraint (3.6) for v_i and v_j . This is equivalent to the original formulation of the problem.

In general, any linear program can be solved in $\mathcal{O}(n^{3.5}r^2\log r\log\log r)$ time by

using Karmarkar's algorithm [31]. However, our linear program is special which allows to use faster techniques:

Proposition 3.9. The BOUNDREP problem with prescribed \triangleleft can be solved in time $\mathcal{O}(n^2r + nD(r))$.

Proof. Without loss of generality, we assume that the upper and lower bounds restrict the final representation (if it exists) to lie in the interval [1, n+3]. For a given i, let j_i be the index such that v_{j_i} is the leftmost neighbor of v_i in \triangleleft . Let h_i be the index such that v_{h_i} is the rightmost vertex such that $v_{h_i} \triangleleft v_i$ and $v_{h_i}v_i \notin E$. (Notice that h_i might not be defined, in which case we ignore inequalities containing it.)

We replace the variables ℓ_1, \ldots, ℓ_k by x_0, \ldots, x_k such that $\ell_i = x_i - x_0$. We want to solve the following linear system:

Minimize:
$$E_t := x_k - x_0 + 1,$$
subject to:
$$x_i - x_{i+1} \le 0, \qquad \forall i = 1, \dots, k - 1,$$

$$x_0 - x_i \le -\text{lbound}(v_i), \qquad \forall i = 1, \dots, k,$$

$$x_i - x_0 \le \text{ubound}(v_i), \qquad \forall i = 1, \dots, k,$$

$$x_i - x_{j_i} \le 1, \qquad \forall i = 1, \dots, k,$$

$$x_{h_i} - x_i \le -1 - \varepsilon, \qquad \forall i = 1, \dots, k.$$

The obtained linear program is a system of difference constraints, since each inequality has the form $x_i - x_j \leq b_{i,j}$.

Following [10, Chapter 24.4], if the system is feasible, a solution, which is not necessarily optimal, can be found as follows. We define a weighted digraph D as follows. As the vertices, we have $V(D) = \{s, u_0, u_1, \ldots, u_k\}$ where u_i corresponds to x_i and s is a special vertex. For the edges $\vec{E}(D)$, we first have an edge (s, u_i) of the weight zero for every u_i . Then for every constraint $x_i - x_j \leq b_{i,j}$, we add the edge (u_i, u_i) of the weight $b_{i,j}$. See Figure 3.7.

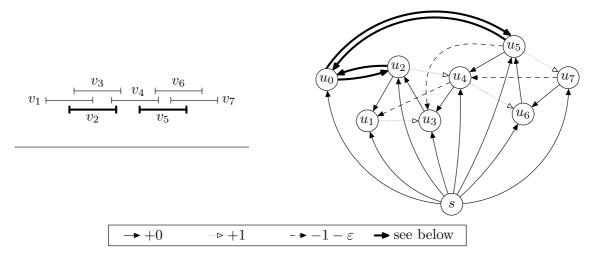


Figure 3.7: On the left, a unit interval graph with two pre-drawn intervals. On the right, the corresponding digraph D with the weight encoded as in the box. The weights of the bold edges are as follows: $w(u_0, u_2) = -\text{lbound}(v_2)$, $w(u_0, u_5) = -\text{lbound}(v_5)$, $w(u_2, u_0) = \text{ubound}(v_2)$, and $w(u_5, u_0) = \text{ubound}(v_5)$.

As proved in [10, Chapter 24.4], there are two possible cases. If G contains a negative-weight cycle, then there is no feasible solution for the system. If G does not contain negative-weight cycles, then we define $\delta(s, u_i)$ as the weight of the minimum-weight path connecting s to u_i in G. Then we put $x_i = \delta(s, u_i)$ for each i which defines a feasible solution of the system. Moreover, this solution minimizes the objective function $\max\{x_i\} - \min\{x_i\}$. We next show that this function is equivalent to the objective function in our linear program.

Suppose that we have a solution of our system, satisfying the constraints but not necessarily optimizing the objective function. Because of our assumption that the representation lies in the interval [1, n + 3], we know that $\ell_i > 0$ for all i. Therefore, $x_i > x_0$. So $\min\{x_i\}$ is always attained by x_0 , while $\max\{x_i\}$ is always attained by x_k . So minimization of the objective function $\max\{x_i\} - \min\{x_i\}$ is equivalent to the original minimization of $E_t = x_k - x_0 + 1$.

In order to find a negative-weight cycle in D or, alternatively, compute the weight of the minimum-weight paths from s to all the other vertices of D, we use the Bellman-Ford algorithm. Notice that Dijkstra's algorithm cannot be used in this case, since some edges of D have negative weight. We next analyze the running time of the whole procedure.

We assume that the cost of arithmetic operations with large numbers is not constant. The algorithm computes the value ε in the beginning which can be clearly done in time $\mathcal{O}(nD(r))$. (And instead of the least common multiple we can simply consider product of q_i 's.)

Afterwards, we compute the weights of the edges of D as multiples of ε , which takes time $\mathcal{O}(kD(r))$. Then each step of the Bellman-Ford algorithm requires time $\mathcal{O}(r)$, and the algorithm runs $\mathcal{O}(k^2)$ steps in total. The total time to solve each linear program is therefore $\mathcal{O}(k^2r + kD(r))$. Finally, the total time of the algorithm is $\mathcal{O}(n^2r + nD(r))$.

In the next subsections, we improve the time complexity of the BOUNDREP problem with prescribed \triangleleft to $\mathcal{O}(n^2 + nD(r))$. Our algorithm makes use of several structural properties of the set of all representations. We note that structural properties of the polyhedron of our linear program, in the case where all lower bounds equal zero and there are no upper bounds, have been considered in several papers in the context of semiorders [45, 4].

3.3.2 The partially ordered set $\Re \mathfrak{p}$

Let the graph G in consideration be a connected unit interval graph. We study structural properties of its representations. Suppose that we fix one of the two partial left-to-right orders < of the intervals from Section 3.1, so that only indistinguishable vertices are incomparable. We also fix some positive $\varepsilon = \frac{1}{K}$. For most of this section, we work just with lower bounds and completely ignore upper bounds.

We define \mathfrak{Rep} as the set of all ε -grid representations satisfying the lower bounds and in some the left-to-right ordering that extends <. We define a very natural partial

ordering \leq on \mathfrak{Rep} : We say that $\mathcal{R} \leq \mathcal{R}'$ if and only if $\ell_i \leq \ell'_i$ for every $v_i \in V(G)$. In this section, we study structural properties of the poset (\mathfrak{Rep}, \leq) .

If $\varepsilon \leq \frac{1}{n}$, it holds that $\mathfrak{Rep} \neq \emptyset$, since the graph G is a unit interval graph and there always exists an ε -grid representation \mathcal{R} far to the right satisfying the lower bound contraints.

The Semilattice Structure. Let us assume that $lbound(v_i) > -\infty$ for some $v_i \in V(G)$. Let S be a subset of \mathfrak{Rep} . The infimum $\inf(S)$ is the greatest representation $\mathcal{R} \in \mathfrak{Rep}$ such that $\mathcal{R} \leq \mathcal{R}'$ for every $\mathcal{R}' \in S$. In a general poset, infimums may not exist, but if they exist, they are always unique. For \mathfrak{Rep} , we show:

Lemma 3.10. Every non-empty $S \subseteq \Re \mathfrak{sp}$ has an infimum $\inf(S)$.

Proof. We construct the requested infimum \mathcal{R} as follows:

$$\ell_i = \min\{\ell'_i : \mathcal{R}' \in S\}, \quad \forall v_i \in V(G).$$

Notice that the positions in \mathcal{R} are well-defined, since the position of each interval in each \mathcal{R}' is bounded and always on the ε -grid. Clearly, if \mathcal{R} is a correct representation, it is the infimum $\inf(S)$. It remains to show that $\mathcal{R} \in \mathfrak{Rep}$.

Clearly, all positions in \mathcal{R} belong to the ε -grid and satisfy the lower bound constraints. Let v_i and v_j be two vertices. The values ℓ_i and ℓ_j in \mathcal{R} are given by two representations $\mathcal{R}_1, \mathcal{R}_2 \in S$, that is, $\ell_i = \ell_i^1$ and $\ell_j = \ell_j^2$. Notice that the left-to-right order in \mathcal{R} has to extend <: If $v_i < v_j$, then $\ell_i = \ell_i^1 \le \ell_i^2 < \ell_j^2 = \ell_j$, since \mathcal{R}_1 minimizes the position of v_i and the left-to-right order in \mathcal{R}_2 extends <. Concerning correctness of the representation of the pair v_i and v_j , we suppose that $\ell_i = \ell_i^1 \le \ell_j^2 = \ell_j$; otherwise we swap v_i and v_j .

- First we suppose that $v_i v_j \in E(G)$. Then $\ell_j^2 \leq \ell_j^1$, since \mathcal{R}_2 minimizes the position of v_j . Since \mathcal{R}_1 is a correct representation, $\ell_j^1 1 \leq \ell_i^1$. So $\ell_j 1 \leq \ell_i \leq \ell_j$, and the intervals v_1 and v_2 intersect.
- The other case is when $v_i v_j \notin E(G)$. Then $\ell_i^1 \leq \ell_i^2 \leq \ell_j^2 1 \varepsilon$, since \mathcal{R}_1 minimizes the position of v_i , \mathcal{R}_2 is a correct representation and $v_i < v_j$ in both representations. So v_i and v_j do not intersect in \mathcal{R} as requested.

Consequently, \mathcal{R} represents correctly each pair v_i and v_j , and hence $\mathcal{R} \in \mathfrak{Rep}$.

A poset is a *(meet)-semilattice* if every pair of elements a, b has an infimum $\inf(\{a,b\})$. Lemma 3.10 shows that the poset (\mathfrak{Rep}, \leq) forms a (meet)-semilattice. Similarly as \mathfrak{Rep} , we could consider the poset set of all $(\varepsilon$ -grid) representations satisfying both the lower and the upper bounds. The structure of this poset is a *complete lattice*, since all subsets have infimums and supremums. Lattices and semilattices are frequently studied, and posets that are lattices satisfy very strong algebraic properties.

The Left-most Representation. We are interested in a specific representation in \mathfrak{Rep} , called the *left-most representation*. An ε -grid representation $\mathcal{R} \in \mathfrak{Rep}$ is the left-most representation if $\mathcal{R} \leq \mathcal{R}'$ for every $\mathcal{R}' \in \mathfrak{Rep}$; so the left-most representation is left-most in each interval at the same time. We note that the notion of the left-most

representation does not make sense if we consider general representations (not on the ε -grid). The left-most representation is the infimum inf(\mathfrak{Rep}), and thus by Lemma 3.10 we get:

Corollary 3.11. The left-most representation always exists and it is unique.

There are two algorithmic motivations for studying left-most representations. First, in the linear program of Section 3.3.1 we need to find a representation minimizing E_t . Clearly, the left-most representation is minimizing E_t and in addition it is minimizing the rest of the endpoints as well. The second motivation is that we want to construct a representation satisfying the upper bounds as well, so it seems reasonable to try to place every interval as far to the left as possible. The left-most representation is indeed a good candidate for a bounded representation:

Lemma 3.12. There exists a representation \mathcal{R}' satisfying both lower and upper bound constraints if and only if the left-most representation \mathcal{R} satisfies the upper bound constraints.

Proof. Since $\mathcal{R} \in \mathfrak{Rep}$, it satisfies the lower bounds. If \mathcal{R} satisfies the upper bound constraints, it is a bounded representation. On the other hand, let \mathcal{R}' be a bounded representation. Then

$$lbound(v_i) \le \ell_i \le \ell'_i \le ubound(v_i), \quad \forall v_i \in V(G),$$

and the left-most representation is also a bounded representation.

3.3.3 Left-Shifting of Intervals

Suppose that we construct some initial ε -grid representation that is not the left-most representation. We want to transform this initial representation in $\Re \mathfrak{ep}$ into the left-most representation of $\Re \mathfrak{ep}$ by applying the following simple operation called *left-shifting*. The left-shifting operation shifts one interval of the representations by ε to the left such that this shift maintains the correctness of the representation; for an example see Figure 3.8a. The main goal of this section is to prove that by left-shifting we can always produce the left-most representation.

Proposition 3.13. For $\varepsilon = \frac{1}{K}$ and $K \geq \frac{n}{2}$, an ε -grid representation $\mathcal{R} \in \mathfrak{Rep}$ is the left-most representation if and only if it is not possible to shift any single interval to the left by ε while maintaining correctness of the representation.

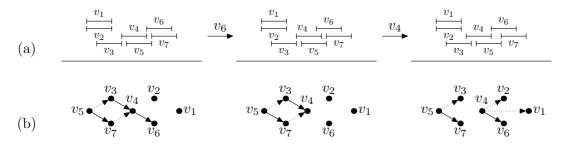


Figure 3.8: (a) A representation modified by left-shifting of v_6 and v_4 . (b) The corresponding obstruction digraph H for each of the representations. Only sinks of the obstruction digraph can be left-shifted.

Before proving the proposition, we describe some additional combinatorial structure of left-shifting. An interval v_i is called *fixed* if it is in the left-most position and cannot ever be shifted more to the left, i.e., $\ell_i = \min\{\ell'_i : \mathcal{R}' \in \mathfrak{Rep}\}$. For example, an interval v_i is fixed if $\ell_i = \text{lbound}(v_i)$. A representation is the left-most representation if and only if every interval is fixed.

Obstruction Digraph. An interval v_i , having $\ell_i > \text{lbound}(v_i)$, can be left-shifted if it does not make the representation incorrect, and the incorrectness can be obtained in two ways. First, there could be some interval v_j , $v_j \triangleleft v_i$ such that $v_i v_j \notin E(G)$ and $\ell_j + 1 + \varepsilon = \ell_i$; we call v_j a left obstruction of v_i . Second, there could be some interval v_j , $v_i \triangleleft v_j$ such that $v_i v_j \in E(G)$ and $\ell_i + 1 = \ell_j$ (so v_i and v_j are touching); then we call v_j a right obstruction of v_i . In both cases, we first need to move v_j before moving v_i .

For the current representation \mathcal{R} , we define the obstruction digraph H on the vertices of G as follows. We put V(H) = V(G) and $(v_i, v_j) \in E(H)$ if and only if v_j is an obstruction of v_i . For an edge (v_i, v_j) , if $v_j \triangleleft v_i$, we call it a left edge; if $v_i \triangleleft v_j$, we call it a right edge. As we apply left-shifting, the structure of H changes; see Figure 3.8b.

Lemma 3.14. An interval v_i is fixed if and only if there exists an oriented path in H from v_i to v_j such that $\ell_j = \text{lbound}(v_j)$.

Proof. Suppose that v_i is connected to v_j by a path in H. By definition of H, $v_x v_y \in E(H)$ implies that v_y has to be shifted before v_x . Thus v_j has to be shifted before moving v_i which is not possible since $\ell_j = \text{lbound}(v_j)$.

On the other hand, suppose that v_i is fixed, i.e., $\ell_i = \inf\{\ell'_i : \forall \mathcal{R}'\}$. Let H' be the induced subgraph of H of the vertices v_j such that there exists an oriented path from v_i to v_j . If for all $v_j \in H'$, $\ell_j > \text{lbound}(v_j)$, we can shift all vertices of H' by ε to the left which constructs a correct representation and contradicts that v_i is fixed. Therefore, there exists $v_j \in H'$ having $\ell_j = \text{lbound}(v_j)$ as requested.

For the example in Figure 3.8 on the left, if $\ell_4 = \text{lbound}(v_4)$, then the intervals v_3 , v_4 , v_5 and v_7 are fixed. Also, we can easily prove:

Lemma 3.15. If $\varepsilon = \frac{1}{K}$ and $K \geq \frac{n}{2}$, the obstruction digraph H is acyclic.

Proof. Suppose for contradiction that H contains some cycle u_1, \ldots, u_c . This cycle contains a left edges and b right edges. Recall that if (u_i, u_{i+1}) is a left edge, then $\ell_{u_{i+1}} = \ell_{u_i} - 1 - \varepsilon$, and if it is a right edge, $\ell_{u_{i+1}} = \ell_{u_i} + 1$ (and similarly for (u_c, u_1)). If we go along the cycle from u_1 to u_1 , the initial and the final positions have to be the same. Therefore $a(1 + \varepsilon) = b$.

Now if this equation holds, then a has to be a multiple of K. Therefore $a \ge K$ and $b \ge K + 1$, and thus $n \ge c = a + b \ge 2K + 1$ which is not possible.

We note that the assumption $K \geq \frac{n}{2}$ is necessary and tight. For every $\varepsilon = \frac{1}{K}$, there exists a representation of a graph with 2K+1 vertices having a cycle in H. The graph contains two cliques v_0, \ldots, v_{K-1} and w_0, \ldots, w_K such that v_i is also adjacent to

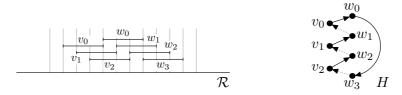


Figure 3.9: An ε -grid representation for $\varepsilon = \frac{1}{3}$ on the left and the obstruction digraph H containing a cycle on the right.

 w_0, \ldots, w_i . Then the assignment $\ell_{v_0} = 0$, $\ell_{v_i} = \ell_{v_0} + i\varepsilon$ and $\ell_{w_i} = \ell_{v_0} + 1 + i\varepsilon$ is a correct representation. Observe that H contains a cycle $w_k v_{k-1} w_{k-1} v_{k-2} w_{k-2} \ldots v_1 w_1 v_0 w_0 w_k$. See Figure 3.9 for K = 3.

Predecessors of Poset \mathfrak{Rep} . A representation $\mathcal{R}' \in \mathfrak{Rep}$ is a *predecessor* of $\mathcal{R} \in \mathfrak{Rep}$ if $\mathcal{R}' < \mathcal{R}$ and there is no representation $\bar{\mathcal{R}} \in \mathfrak{Rep}$ such that $\mathcal{R}' < \bar{\mathcal{R}} < \mathcal{R}$. We denote the predecessor relation by \prec . In a general poset, predecessors may not exist. But they always exist for a poset of a discrete structure like (\mathfrak{Rep}, \leq) : Indeed, there are finitely many representations $\bar{\mathcal{R}}$ between any $\mathcal{R}' < \mathcal{R}$, and thus the predecessors always exist. Also, for any two representations $\mathcal{R}' < \mathcal{R}$, there exists a finite *chain* of predecessors $\mathcal{R}' = \mathcal{R}_0 \prec \mathcal{R}_1 \prec \cdots \prec \mathcal{R}_k = \mathcal{R}$.

For the poset (\mathfrak{Rep}, \leq) , we are able to fully describe the predecessor structure:

Lemma 3.16. For $\varepsilon = \frac{1}{K}$ and $K \geq \frac{n}{2}$, the representation \mathcal{R}' is a predecessor of \mathcal{R} if and only if \mathcal{R}' is obtained from \mathcal{R} by applying one left-shifting operation.

Proof. Clearly, if \mathcal{R}' is obtained from \mathcal{R} by one left-shifting, it is a predecessor of \mathcal{R} .

On the other hand, suppose we have $\mathcal{R}' < \mathcal{R}$. Let H be the obstruction digraph of \mathcal{R} and \bar{H} be the subgraph of H induced by the intervals having different positions in \mathcal{R} and \mathcal{R}' . Then there are no directed edges from \bar{H} to $H \setminus \bar{H}$ (otherwise \mathcal{R}' would be an incorrect representation). According to Lemma 3.15, the digraph \bar{H} is acyclic. Therefore, it contains at least one sink v_i . By left-shifting v_i in \mathcal{R} , we create a correct representation $\bar{\mathcal{R}} \in \mathfrak{Rep}$. Clearly, $\mathcal{R}' \leq \bar{\mathcal{R}} \prec \mathcal{R}$, and so \mathcal{R}' is a predecessor of \mathcal{R} if and only if $\mathcal{R}' = \bar{\mathcal{R}}$.

Again, the assumption on the value of ε is necessary. For example in Figure 3.9, the structure of $\Re \mathfrak{p}$ is just a single chain where a predecessor of some representation is obtained by moving everything by ε to the left.

Proof of Left-shifting Proposition. The main proposition of this subsection is a simple corollary of Lemma 3.16.

Proof of Proposition 3.13. The left-most representation \mathcal{R} is $\inf(\mathfrak{Rep})$, so it has no predecessors and nothing can be left-shifted. On the other hand, if $\inf(\mathfrak{Rep}) < \mathcal{R}$, there is a chain of predecessors in between which implies using Lemma 3.16 that it is possible to left-shift some interval.

3.3.4 Preliminaries for the Shifting Algorithm

Before describing the shifting algorithm, we present several results which simplify the graph and the description of the algorithm.

Pruned Graph. The obstruction digraph H may contain many edges since each vertex v_i can have many obstructions. But if v_i has many, say, left obstructions, these obstructions have to be positioned the same. If two intervals u and v have the same position in a correct unit interval representation, then N[u] = N[v] and they are indistinguishable. Our goal is to construct a pruned graph G' which replaces each group of indistinguishable vertices of G by a single vertex. This construction is not completely straightforward since indistinguishable vertices may have different lower and upper bounds.

Let $\Gamma_1, \ldots, \Gamma_k$ be the vertices partitioned by groups of indistinguishable vertices (and the groups are ordered by < from left to right). We construct a unit interval graph G', where the vertices are $\gamma_1, \ldots, \gamma_k$ with lbound $(\gamma_i) = \max\{\text{lbound}(v_j) : v_j \in \Gamma_i\}$, and the edges E(G') correspond to the edges between the groups of G.

Suppose that we have the left-most representation \mathcal{R}' of the pruned graph G' and we want to construct the left-most representation \mathcal{R} of G. Let Γ_{ℓ} be a group. We place each interval $v_i \in \Gamma_{\ell}$ as follows. Let $\gamma_{\leftarrow}^{\ell}$ be the first non-neighbor of γ_{ℓ} on the left and $\gamma_{\rightarrow}^{\ell}$ be the right-most neighbor of γ_{ℓ} (possibly $\gamma_{\rightarrow}^{\ell} = \gamma_{\ell}$). We set

$$\ell_i = \max\{\text{lbound}(v_i), \ell_{\gamma_{\leftarrow}^{\ell}} + 1 + \varepsilon, \ell_{\gamma_{\rightarrow}^{\ell}} - 1\}, \tag{3.7}$$

and if $\gamma_{\leftarrow}^{\ell}$ does not exist, we ignore it in max. The meaning of this formula is to place each interval as far to the left as possible while maintaining the structure of \mathcal{R}' . Figure 3.10 contains an example of the construction of \mathcal{R} .

Before proving correctness of the construction of \mathcal{R} , we show two general properties of the formula (3.7). The first lemma states that each interval $v_i \in \Gamma_\ell$ is not placed in \mathcal{R} too far from the position of γ_ℓ is \mathcal{R}' .

Lemma 3.17. For each $v_i \in \Gamma_{\ell}$, it holds

$$\ell_{\gamma_{\ell}} - 1 \le \ell_i \le \ell_{\gamma_{\ell}}. \tag{3.8}$$

Proof. The first inequality is true since $\ell_{\gamma_{\ell}} - 1 \leq \ell_{\gamma_{\rightarrow}} - 1 \leq \ell_{i}$ holds according to (3.7) and the ordering < for \mathcal{R}' . The second inequality holds since since \mathcal{R}' is a correct bounded representation, and so $\ell_{\gamma_{\ell}}$ is greater than or equal to each term in (3.7). \square

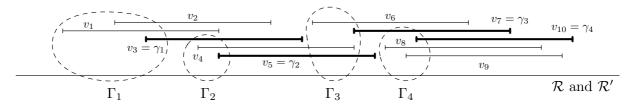


Figure 3.10: Both representations \mathcal{R} and \mathcal{R}' in one figure, with intervals of \mathcal{R}' depicted in bold. The left endpoints of intervals of each group are enclosed by dashed curves, and these curves are ordered from left to right according to <.

The second lemma states that the representations \mathcal{R} and \mathcal{R}' are intertwining each other. If \mathcal{R} is drawn on top of \mathcal{R}' , then the vertices of each group Γ_{ℓ} are in between of $\gamma_{\ell-1}$ and γ_{ℓ} ; see Figure 3.10.

Lemma 3.18. For each $v_i \in \Gamma_{\ell}$, $\ell > 1$, it holds

$$\ell_{\gamma_{\ell-1}} < \ell_i \le \ell_{\gamma_{\ell}},\tag{3.9}$$

Proof. The second inequality holds by (3.8). For the first inequality, there are two possible cases why the groups $\Gamma_{\ell-1}$ and Γ_{ℓ} are distinct:

- The first case is when $\gamma_{\leftarrow}^{\ell}$ is a neighbor of $\gamma_{\ell-1}$. Then $\ell_{\gamma_{\ell-1}} \leq \ell_{\gamma_{\leftarrow}^{\ell}} + 1 < \ell_i$; the first inequality holds since $\gamma_{\leftarrow}^{\ell} \gamma_{\ell-1} \in E(G')$ and \mathcal{R}' is a correct representation, and the second inequality is given by (3.7).
- The second case is when $\gamma_{\rightarrow}^{\ell}$ is a non-neighbor of $\gamma_{\ell-1}$. Then $\ell_{\gamma_{\ell-1}} < \ell_{\gamma_{\rightarrow}^{\ell}} 1 \le \ell_i$ by the fact that $\gamma_{\ell-1}\gamma_{\rightarrow}^{\ell} \notin E(G')$ and by (3.7).

In both cases, we get $\gamma_{\ell-1} < \ell_i$.

Now, we are ready to show correctness of the construction of \mathcal{R} .

Proposition 3.19. From the left-most representation \mathcal{R}' of the pruned graph G', we can construct the correct left-most representation \mathcal{R} of G by placing the intervals according to (3.7).

Proof. We argue the correctness of the representation \mathcal{R} . Let v_i and v_j be a pair of vertices of G. Let $v_i v_j \in E(G)$. If v_i and v_j belong to the same group Γ_ℓ , they intersect each other at position ℓ_{γ_ℓ} by (3.8). Otherwise let $v_i \in \Gamma_\ell$ and $v_j \in \Gamma_{\ell'}$, and assume that $\Gamma_\ell < \Gamma_{\ell'}$. Then $\ell_i \leq \ell_{\gamma_\ell} \leq \ell_j$ by the intertwining property (3.9). Also, $\ell_j \leq \ell_{\gamma_{\ell'}} \leq \ell_{\gamma_{\ell'}} \leq \ell_i + 1$ since $\gamma_{\ell'}$ is a right neighbor of γ_ℓ and (3.8). Therefore, $\ell_i \leq \ell_j \leq \ell_i + 1$ and v_i intersects v_j in \mathcal{R} . Now, let $v_i v_j \notin E(G)$, $v_i \in \Gamma_\ell$, $v_j \in \Gamma_{\ell'}$ and $v_i < v_j$. Then $\ell_i \leq \ell_{\gamma_\ell} \leq \ell_{\gamma_\ell'} \leq \ell_j - 1 - \varepsilon$ by (3.7) and (3.8), so v_i and v_j do not intersect. So the assignment \mathcal{R} is a correct representation of G.

It remains to show that \mathcal{R} is the left-most representation of G. We can identify each γ_{ℓ} with one interval $v_i \in \Gamma_{\ell}$ having $\operatorname{lbound}(v_i) = \operatorname{lbound}(\gamma_{\ell})$; for an example see Figure 3.10. So G' can be viewed as an induced subgraph of G. We want to show that the intervals of G' are represented in \mathcal{R} exactly the same as in \mathcal{R}' . Since $\mathcal{R}|_{G'}$ (\mathcal{R} restricted to G') is some representation of G' and \mathcal{R}' is the left-most representation of G', $\ell'_{\gamma_{\ell}} \leq \ell_{\gamma_{\ell}}$ for every γ_{ℓ} . By (3.8), we get $\ell'_{\gamma_{\ell}} = \ell_{\gamma_{\ell}}$.

We know that $\mathcal{R}|_{G'}$ is the left-most representation, or in other words each interval of G' is fixed in \mathcal{R} . The rest of the intervals are placed so that they are either trivially fixed by $\ell_i = \text{lbound}(v_i)$, or they have as obstructions some fixed intervals from G', in which case they are fixed by Lemma 3.14. Therefore, every interval of G is fixed and \mathcal{R} is the left-most representation.

For the pruned graph G', the obstruction digraph H has in and out-degree at most two. Each interval has at most one left obstruction and at most one right

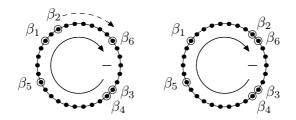


Figure 3.11: Examples of position cycles. In the cycle on the left, we can shift β_2 in clockwise direction towards β_6 , which gives a new representation whose position cycle is depicted on the right. We note that after left-shifting, v_6 is not necessarily an obstruction of v_2 .

obstruction, and these obstructions are always the same. More precisely, if v_j is a left obstruction of v_i , then $v_j = v_{\leftarrow}^i$, whereas if v_j is a right obstruction of v_i , then $v_j = v_{\leftarrow}^i$.

The pruning operation can be done in time $\mathcal{O}(n+m)$, so we may assume that our graph G is already pruned and contains no indistinguishable vertices. And the structure of obstructions in G can be computed in time $\mathcal{O}(n+m)$ as well.

Position Cycle. For each interval in some ε -grid representation, we can write its position in this form:

$$\ell_i = \alpha_i + \beta_i \varepsilon, \qquad \alpha_i \in \mathbb{Z}, \ \beta_i \in \mathbb{Z}_K,$$
 (3.10)

where $\varepsilon = \frac{1}{K}$. In other words, α_i is an integer position of v_i in the grid and β_i describes how far is this interval from this integer position.

Concerning left-shifting, the values β_i are more important. We can depict $\mathbb{Z}_K = \{0, \ldots, K-1\}$ as a cycle with K vertices where the value decreases clockwise. The value β_i assigns to each interval v_i one vertex of the cycle. The cycle \mathbb{Z}_K together with marked positions of β_i 's is called the position cycle. A vertex of the position cycle is called taken if some β_i is assigned to it, and empty otherwise. The position cycle allows us to visualize and work with left-shifting very intuitively. When an interval v_i is left-shifted, β_i cyclically decreases by one, so β_i moves clockwise along the cycle. For an illustration, see Figure 3.11.

If (v_i, v_j) is a left edge of H, then $\beta_j = \beta_i - 1$, and if (v_i, v_j) is a right edge, then $\beta_i = \beta_j$. So if v_j is an obstruction of v_i , β_j has to be very close to β_i (either at the same position or at the next clockwise position). If there is a big empty space in the clockwise direction from β_i , the interval v_i can be left-shifted many times (or till it becomes fixed by $\ell_i = \text{lbound}(v_i)$). Notice that if β_i is very close to β_j , it does not mean that ℓ_i is very close to ℓ_j because the values α_i and α_j are ignored in the position cycle.

3.3.5 The Shifting Algorithm for BOUNDREP

We want to solve an instance of BOUNDREP with a prescribed ordering \triangleleft . We work with an ε -grid which is different from the one in Section 3.2.1. The new value of ε is the value given by (3.1) refined n times, so

$$\varepsilon = \frac{1}{n^2} \cdot \varepsilon'.$$

Lemma 3.5 applies for this value of ε as well, so if the instance is solvable, there exists a solution which is on this ε -grid.

The algorithm works exactly as the algorithm of Subsection 3.3.1. The only difference is that for a component with k vertices we can solve the linear program in time $\mathcal{O}(k^2 + kD(r))$, and now we describe how to do it. We assume that the input component is already pruned, otherwise we prune it and use Proposition 3.19 to complete the representation. We expect that the left-to-right order \triangleleft of the vertices is given. The algorithm requires time $\mathcal{O}(kD(r))$ since the bounds are given in form $\frac{p_i}{q_i}$ and we need to perform arithmetic operations with these bounds. Therefore the total complexity of the algorithm for the BOUNDREP problem is $\mathcal{O}(n^2 + nD(r))$.

Overview. The algorithm for solving one linear program works in three basic steps:

- 1. We construct an initial ε -grid representation in the ordering \triangleleft which satisfies $\ell_i \geq \text{lbound}(v_i)$ for all intervals, using the algorithm of Corneil et al. [12].
- 2. We shift the intervals to the left while maintaining correctness of the representation until the left-most representation is constructed, using Proposition 3.13.
- 3. We check whether the left-most representation satisfies the upper bounds. If so, we have the left-most representation satisfying all bound constraints. This representation solves the linear program of Subsection 3.3.1 and minimizes E_t . Otherwise, the left-most representation does not satisfy the upper bound constraints, and thus by Lemma 3.12 no representation satisfying the upper bound constraints exists, and the linear program has no solution.

Input Size. Let r be the size of the input describing bound constraints. A standard complexity assumption is that we can operate with polynomially large numbers (having $\mathcal{O}(\log r)$ bits in binary) in constant time, to avoid extra factor $\mathcal{O}(\log r)$ in the complexity of most of the algorithms. However, the value of ε given by (3.1) might require $\mathcal{O}(r)$ digits when written in binary. The assumption that we can computate with numbers having $\mathcal{O}(r)$ digits in contant time would break most of the computational models. Therefore, our computational model requires a larger time for arithmetic operations with numbers having $\mathcal{O}(r)$ digits in binary. For example, the best known algorithm for multiplication/division on a Turing machine requires time $\mathcal{O}(D(r))$.

The problem is that a straightforward implementation of our algorithm working with the ε -grid would require time $\mathcal{O}(k^2r^c)$ for some c instead of $\mathcal{O}(k^2+kD(r))$. There is an easy way out. Instead of computing with long numbers having $\mathcal{O}(r)$ digits, we mostly compute with short numbers having just $\mathcal{O}(\log r)$ digits. Instead of the ε -grid, we mostly work in a larger Δ -grid where $\Delta = \frac{1}{n^2}$. The algorithm computes with the long numbers only in two places. First, some initial computations concerning the input are performed. Second, when the shifting makes some interval fixed, the algorithm estimes the final ε -grid position of the interval. All these computations can be done in total time $\mathcal{O}(kD(r))$ and we describe everything in detail later.

Left-Shifting. The basic operation of the algorithm is the LEFTSHIFT procedure which we describe here. We deal separately with fixed and unfixed intervals (and some intervals might be fixed initially). Unfixed intervals are on the Δ -grid and fixed intervals have precise positions calculated on the ε -grid. We place only unfixed intervals on the

position cycle for the Δ -grid. At any moment of the algorithm, each vertex of the position cycle is taken by at most one β_i ; this is true for the initial representation and the shifting keeps this property.

We define the procedure LeftShift(v_i) which shifts v_i from the position ℓ_i into a new position ℓ_i' such that the representation remains correct. The procedure LeftShift(v_i) consists of two steps:

- 1. Since v_i is unfixed, it has some β_i placed on the position cycle. Let k be such that the vertices $\beta_i + 1, \ldots, \beta_i + k$ of the position cycle are empty and the vertex $\beta_i + k + 1$ is taken by some β_b . Then a candidate for the new position of v_i is $\bar{\ell}_i = \ell_i k\Delta$.
- 2. We need to ensure that this shift from ℓ_i to $\bar{\ell}_i$ is valid with respect to lbound(v_i) and the positions of the fixed intervals. Concerning the lower bound, we cannot shift further than lbound(v_i). Concerning the fixed intervals, the shift is limited by positions of fixed obstructions of v_i . If v_j is a fixed left obstruction, we cannot shift further than $\ell_j + 1 + \varepsilon$, and if $v_{j'}$ a fixed right obstruction, we cannot shift further than $\ell_{j'} 1$.

The resulting position after applying LeftShift (v_i) is

$$\ell_i' = \max\{\bar{\ell}_i, \text{lbound}(v_i), \ell_j + 1 + \varepsilon, \ell_{j'} - 1\}. \tag{3.11}$$

Lemma 3.20. If the original representation \mathcal{R} is correct, than the LeftShift (v_i) procedure produces a correct representation \mathcal{R}' .

Proof. Clearly, the lower bound for v_i is satisfied in \mathcal{R}' . The shift of v_i from ℓ_i to ℓ'_i can be viewed as a repeated application of the left-shifting operation from Section 3.3.3. We just need to argue that each left-shifting operation can be applied till the position ℓ'_i is reached.

If at some point, the left-shifting operation could not be applied, there would have to be some obstruction v_j of v_i . There is no unfixed obstruction since all vertices of the position cycle $\beta_i + 1, \ldots, \beta_i + k$ are empty. And v_j cannot be fixed as well since we check positions of both possible obstructions. So there is no obstruction v_j . Therefore, by repeated applying the left-shifting operation, the interval v_i gets at a position ℓ'_i and the resulting representation is correct.

After LeftShift(v_i), if $\bar{\ell}_i$ is not a strict maximum of the four terms in (3.11), the interval v_i becomes fixed; either trivially since $\ell'_i = \text{lbound}(v_i)$, or by Lemma 3.14 since v_i becomes obstructed by some fixed interval. In such a case, we remove β_i from the position cycle.

Fast Implementation of Left-Shifting. Since we apply the LEFTSHIFT procedure repeatedly, we want to implement it in time $\mathcal{O}(1)$. Considering the terms in (3.11), the first term $\bar{\ell}_i$ is a short number (on the Δ -grid) and the remaining terms are long numbers (on the ε -grid). We first compare $\bar{\ell}_i$ to the remaining terms which are three comparisons of short and long numbers and we are going to show how to compare them in $\mathcal{O}(1)$. If $\bar{\ell}_i$ is a strict maximum, we use it for ℓ'_i . Otherwise, we need to

compute the maximum of the remaining three terms which takes time $\mathcal{O}(D(r))$. But then the interval v_i becomes fixed, and so this costly step is done exactly k times, and takes the total time $\mathcal{O}(kD(r))$.

Lemma 3.21. With the total precomputation time $\mathcal{O}(kD(r))$, it is possible to compare $\bar{\ell}_i$ to the remaining terms in (3.11) in time $\mathcal{O}(1)$ per LeftShift procedure.

Proof. Initially, we do the following precomputation for the lower bounds. By the input, we have b lower bounds given in the form $\frac{p_1}{q_1}, \ldots, \frac{p_b}{q_b}$ as irreducible fractions. For each bound, we first compute its position (α_i, β_i) on the ε -grid; see (3.10).

If $lbound(v_i) \ll lbound(v_j)$ for some vertices v_i and v_j , then $lbound(v_i)$ is never achieved since the graph is connected and every representation takes space at most k. Therefore we can increase $lbound(v_i)$ without any change in the solution of the instance. More precisely, let $\alpha = \max \alpha_i$. Then we modify each bound by setting $\alpha_i := \max\{\alpha - n - 1, \alpha_i\}$. In addition, we shift all the bounds by substructing a constant C such that each $\alpha_i - C \in [0, k + 1]$. Concerning β_i , we round the position (α_i, β_i) down to a position $(\alpha_i, \bar{\beta}_i)$ of the Δ -grid. These precomputations can be done for all lower bounds in time $\mathcal{O}(kD(r))$.

Suppose that we want to find out whether $\bar{\ell}_i \leq \text{lbound}(v_j) = \alpha_j + \beta_j \cdot \varepsilon$ where $\bar{\ell}_i$ is in the Δ -grid. Then it is sufficient to check whether $\bar{\ell}_i \leq \alpha_j + \bar{\beta}_j \Delta$ which can be done in constant-time since both α_j and $\bar{\beta}_j$ are short numbers.

When v_j becomes fixed, its precise position is computed using (3.11). Then we compute the values $\ell_j - 1$ and $\ell_j + 1 + \varepsilon$ used in (3.11) and round them down to the Δ -grid. Using these precomputed values, $\bar{\ell}_i$ can be compared with the remaining terms in (3.11) in time $\mathcal{O}(1)$. When an interval becomes fixed, time $\mathcal{O}(D(r))$ is used. Since each interval becomes fixed exactly once, this rounding takes the total time $\mathcal{O}(kD(r))$.

Notice that the representation is constructed in a position shifted by C. Later, before checking the upper bound, we shift the whole representation back.

Initial Representation. Recall that $\Delta = \frac{1}{n^2}$ and the position cycle has n^2 vertices. The algorithm of Corneil et al. [12] gives a representation in the $\frac{1}{k}$ -grid. Using the proof of Lemma 3.5, we construct from it the initial Δ -grid representation. Then we shift it such that $\ell_i \geq \text{lbound}(v_i)$ for each v_i and $\ell_i \leq \text{lbound}(v_i) + \Delta$ for some v_i . For this initial representation, each interval can be shifted to the left in total by at most $\mathcal{O}(k)$.

The initial representation obtained from the representation of the algorithm of Corneil et al. [12] places all intervals in such a way that β_i 's are almost positioned equidistantly in the position cycle; refer to the left-most position cycle in Figure 3.12. As we say in the description of the Left-Shift procedure, we only require that all β_i 's are placed to pairwise different vertices of the position cycle.

Shifting Phases. All shifting of the algorithm is done by repeated application of the LEFTSHIFT procedure. Using Lemma 3.20, we know that the representation created in each step is correct. We apply the procedure in such a way that each interval is almost always shifted by almost one. The shifting of unfixed intervals proceeds in two phases:

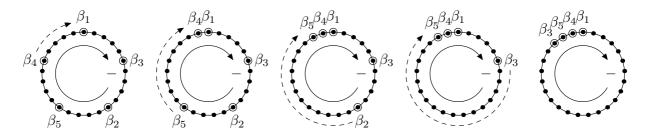


Figure 3.12: The position cycle during the first phase, changing from left to right. The first phase clusters the β_i 's by moving β_4 , β_5 , β_2 and β_3 towards β_1 . When LEFTSHIFT (v_2) is applied, v_2 becomes fixed and β_2 disappears from the position cycle.

- The first phase creates one big gap by clustering all β_i 's in one part of the cycle. To do so, we apply the LeftShift procedure to each interval, in the order given by the position cycle. Of course, some intervals might become fixed and disappear from the position cycle. We obtain one big gap of size at least n(n-1). Again, refer to Figure 3.12.
- In the second phase, we use this big gap to shift intervals one by one, which also moves the cluster along the position cycle. Again, if some interval becomes fixed, it is removed from the position cycle. The second phase finishes when each interval becomes fixed and the left-most representation is constructed. For an example, see Figure 3.13.

Putting Together. For a pseudocode, see Algorithm 7 in Appendix A. First, we show correctness of the shifting algorithm and its complexity:

Lemma 3.22. For a component having k vertices, the shifting algorithm constructs a correct left-most representation in time $\mathcal{O}(k^2 + kD(r))$.

Proof. First, we argue correctness of the algorithm. The algorithm starts with an initial representation which is correct and satisfies the lower bounds. By Lemma 3.20, after applying each Leftshift procedure, the resulting representation is still correct. The algorithm keeps a correct list of fixed intervals which is increased by shifting. So after finitely many applications of the Leftshift procedure, every interval becomes fixed, and we obtain the left-most representation.

Concerning complexity, all precomputations take total time $\mathcal{O}(kD(r))$. Using Lemma 3.21, each LeftShift(v_i) procedure can be applied in time $\mathcal{O}(1)$ unless v_i becomes fixed. The first phase is applying the LeftShift procedure k-1 times. In

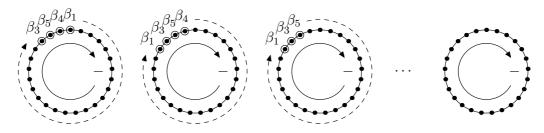


Figure 3.13: The position cycle during the second phase, changing from left to right. We shift β_i 's across the big gap till all β_i 's disappear.

the second phase, each interval is shifted by at least $\frac{n-1}{n}$ (unless it becomes fixed). Since each interval can be shifted by at most $\mathcal{O}(k)$ from its initial position, the second phase applies the LeftShift procedure $\mathcal{O}(k^2)$ times. So the total running time of the algorithm is $\mathcal{O}(k^2 + kD(r))$.

We are ready to prove that BOUNDREP with a prescribe ordering \triangleleft can be solved in time $\mathcal{O}(n^2 + nD(r))$:

Proof of Theorem 1.4. We proceed exactly as in the algorithm of Section 3.3.1, so we process the components $C_1 \blacktriangleleft \cdots \blacktriangleleft C_c$ from left to right, and for each of them we solve two linear programs. For each linear program, we find the left-most representation using Lemma 3.22, and we test for this representation (shifted back by C) whether the upper bounds are satisfied. According to Lemma 3.12, the linear program is solvable if and only if the left-most representation satisfy the upper bounds, and clearly the left-most representation minimizes E_t . The time complexity of the algorithm is $\mathcal{O}(n^2 + nD(r))$ and the proof of correctness is exactly the same as in Proposition 3.13.

We finally present an FPT algorithm for BOUNDREP with respect to the number of components c. The algorithm is based on Theorem 1.4.

Proof of Corollary 1.5. There are c! possible left-to-right orderings of the components of G. For each of them, we can decide in time $\mathcal{O}(n^2 + nD(r))$ whether there exists a bounded representation in the order, using Theorem 1.4. So the total time necessary is $\mathcal{O}((n^2 + nD(r))c!)$.

3.4 Extending Unit Interval Graphs

The Repext(UNIT INT) problem can be solved using Theorem 1.4. We just need to show that it is a particular instance of BoundRep with the known ordering ◀ of the components:

Proof of Corollary 1.6. The graph G contains unlocated components and located components. Similarly to Section 3.1, unlocated components can be placed far to the right and we can deal with them using standard recognition algorithm.

Concerning located components C_1, \ldots, C_c , they have to be ordered from left to right, which gives the required ordering \blacktriangleleft . We straightforwardly construct the instance of BOUNDREP with this \blacktriangleleft as follows. For each pre-drawn interval v_i at position ℓ_i , we put $\text{lbound}(v_i) = \text{ubound}(v_i) = \ell_i$. For the rest of the intervals, we set no bounds. Clearly, this instance of BOUNDREP is equivalent with the original REPEXT(UNIT INT) problem. And we can solve it in time $\mathcal{O}(n^2 + nD(r))$ using Theorem 1.4.

Extending Chordal Graphs and Their Subclasses

The four types FIXED, SUB, ADD, and BOTH form the lattice depicted in Figure 4.1. For a type \mathfrak{T} , we denote by $\operatorname{Gen}(\mathfrak{T}, T')$ the set of all trees T which we can generate from T' using the modifications of the type \mathfrak{T} . In addition, if T' contains pre-drawn subtrees, the trees in $\operatorname{Gen}(\mathfrak{T}, T')$ contain these (possibly subdivided) pre-drawn subtrees as well. The ordering of the types given by the lattice has this property: If $\mathfrak{T} \leq \mathfrak{T}'$, then $\operatorname{Gen}(\mathfrak{T}, T') \subseteq \operatorname{Gen}(\mathfrak{T}', T')$.

Whether a given instance is solvable depends on the set $Gen(\mathfrak{T}, T')$; so if this set contains more trees, it only helps in solving the problem. Let $\mathfrak{T} \leq \mathfrak{T}'$. If an instance of Recog* or Repext is solvable for the type \mathfrak{T} , then it is solvable for the type \mathfrak{T}' as well. Equivalently, if it is not solvable for \mathfrak{T}' , it is also not solvable for \mathfrak{T} .

For the types ADD and BOTH (and SUB for PROPER P-in-P and P-in-P), the set $Gen(\mathfrak{T}, T')$ contains a tree having an arbitrary tree T as a subtree. Therefore, the RECOG* problem for these types is equivalent to the standard RECOG problem, and we can use the known polynomial-time algorithms.

For the classes PROPER P-in-P and P-in-P, the types ADD and SUB behave differently. The type ADD allows to extend the ends of the paths. The type SUB allows to expand the middle of the path but if an endpoint of the path is contained in some pre-drawn subpath, it remains to be contained after the subdivision. The type BOTH makes the problems equivalent to the RECOG and REPEXT problems for the real line.

Indistinguishable Vertices. Let u and v be two vertices of G such that N[u] = N[v]. These two vertices are called *indistinguishable* since they can be represented exactly the same, i.e., $R_u = R_v$. (This is a common property of indistinguishable vertices for all intersection representations). From the structural point of view, groups of indistinguishable vertices are not very interesting. The goal is to construct a pruned

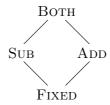


Figure 4.1: The lattice formed by four types FIXED, SUB, ADD, and BOTH.

graph where each group is represented by a single vertex. For that, we need to be little careful since we cannot prune pre-drawn vertices.

For an arbitrary graph, its groups of indistinguishable vertices can be located in time $\mathcal{O}(n+m)$ [49]. We prune the graph in the following way. If u and v are indistinguishable and u is not pre-drawn, we eliminate u from the graph (and for the representation, we can put $R_u = R_v$). In addition, if two pre-drawn subtrees are the same, we eliminate one of them. The resulting pruned graph has the following property: If two vertices u and v indistinguishable, they are both pre-drawn and represented by distinct subtrees. For the rest of the chapter, we assume that all input graphs are pruned.

Maximal Cliques. It is well known that subtrees of a tree possess the Helly property, i.e., every pairwise intersecting collection of subtrees has a non-empty intersection (which is again a subtree). Hence the following holds true for all classes of graphs considered. If K is a maximal clique of G, the common intersection $R_K = \bigcap_{u \in K} R_u$ is a subtree of T. This subtree R_K is not intersected by any other R_v for $v \notin K$ (otherwise K would not be a maximal clique). Thus the subtrees R_K corresponding to different maximal cliques are pairwise disjoint. For example, if |T| is smaller than the number of maximal cliques of G, the graph is clearly not representable in T.

4.1 Subpaths-in-Path Graphs

In this section, we deal with the classes PROPER P-in-P and P-in-P. The results obtained here are used as tools for P-in-T and T-in-T graphs in Section 4.2.

Let p_1, \ldots, p_t be the vertices of the path T'. For a located component C, we say that a vertex p_i is taken by C if there exists a pre-drawn subpath of C containing p_i .

4.1.1 Structural Results

We describe two types of orderings: Endpoint orderings for proper interval graphs and clique orderings for interval graphs. Also, we introduce an important concept called the minimum span of a component.

Endpoint Orderings of PROPER P-in-P. As stated in Section 3.1, each proper interval representation gives some ordering \triangleleft of the intervals from left to right. This is the ordering of the left endpoints from left to right, and at the same time the ordering of the the right endpoints. The following lemma of [14] states that \triangleleft is well determined:

Lemma 4.1 (Deng et al.). For a component of a proper interval graph, the ordering \triangleleft is uniquely determined up to a local reordering of the groups of indistinguishable vertices and the complete reversal.

So for a connected graph, we have a partial ordering < in which exactly the indistinguishable vertices are incomparable, and each \lhd is a linear extension of either <, or its reversal. Corneil et al. [12] describes how this ordering can be constructed in time $\mathcal{O}(n+m)$. Since the graphs we consider are pruned, all incomparable vertices in < are ordered by their positions in the partial representation. Thus we have at

most two possibilities for \triangleleft for each component C. (And two possibilities only if all pre-drawn vertices are indistinguishable.)

Minimum Spans of PROPER P-in-P. For the types FIXED and ADD, the space on the path T' is limited. So it is important to minimize the space taken by each component C. We call the minimum space required by C the minimum span of C, denoted by minspan(C). Let \mathcal{R} be a proper interval representation of C extending \mathcal{R}' , and let p_i be the left-most vertex of T' taken by C and p_i the right-most one. Then

$$\operatorname{minspan}(C) = \begin{cases} \min_{\forall \mathcal{R}} \{j - i + 1\} & \text{if some representation of } C \text{ exists,} \\ +\infty & \text{otherwise.} \end{cases}$$

A representation of C is called *smallest* if it realizes the minimum span of C.

Lemma 4.2. For every component C, the value minspan(C) can be computed in time $\mathcal{O}(n+m)$, together with a smallest representation of C.

Proof. First, we deal with unlocated components, and later modify the approach for the located ones.

Case 1: An Unlocated Component. Since there are no indistinguishable vertices, we compute in time $\mathcal{O}(n+m)$ using the algorithm of [12] any ordering \triangleleft for which we want to produce a representation as small as possible.

Let ℓ_i denote the left endpoint and r_i the right endpoint of the interval v_i . From the ordering $v_1 \lhd \cdots \lhd v_n$, we want to compute the common ordering \lessdot of both the left and the right endpoints from left to right. The starting point is the ordering of just the left endpoints $\ell_1 \lessdot \cdots \lessdot \ell_n$. Into this ordering, we insert the right endpoints r_1, \ldots, r_n one-by-one. A right endpoint r_i is inserted right before ℓ_j where v_j is the left-most non-neighbor of v_i on the right in \lhd ; if such v_j does not exist, we append r_i to the end. For an example of \lessdot , see Figure 4.2.

We build a smallest representation using \leq as follows. Let p_1, \ldots, p_k be the vertices of the tree T. We construct an assignment f which maps the endpoints of the intervals of C into T. Then for a vertex v_i we put

$$R_{v_i} = \{ p_i \mid f(\ell_i) \le p_i \le f(r_i) \}.$$

The mapping f is constructed for the endpoints one-by-one, according to \leq . Suppose that the previous endpoint in \leq has assigned a vertex p_i . If the current endpoint is a right endpoint and the previous endpoint is a left endpoint, we assign p_i to the current endpoint. Otherwise we assign p_{i+1} to it. For an example, see Figure 4.2.

In total, the component needs $2n - \ell$ vertices of T where ℓ denotes the number of changes from a left endpoint to a right endpoint in the ordering \leq ; in other words,

Figure 4.2: The ordering \lessdot is $\ell_1 \lessdot \ell_2 \lessdot r_1 \lessdot \ell_3 \lessdot \ell_4 \lessdot r_2 \lessdot r_3 \lessdot \ell_5 \lessdot r_4 \lessdot \ell_6 \lessdot r_5 \lessdot r_6$ for the component C on the left. The constructed smallest possible representation of the component C on the right, with minspan(C) = 8.

 $2n - \ell$ is the value of minspan(C). The total complexity of the algorithm is clearly $\mathcal{O}(n+m)$.

To conclude the proof, we need to show that we construct a correct smallest representation of C. A property of \lhd is that the closed neighborhood N[v] of every vertex $v \in V(G)$ is consecutive in \lhd . If $v_i v_j \in E(G)$ and $v_i \lhd v_j$, then $\ell_i \lessdot \ell_j \lessdot r_i$, and so R_{v_i} intersects R_{v_j} (between $f(\ell_j)$ and $f(r_i)$). If $v_i v_j \notin E(G)$, then $r_i \lessdot \ell_j$. Thus r_i is placed on the left of ℓ_j in \lessdot , and $R_{v_i} \cap R_{v_j} = \emptyset$ as required.

Concerning the minimality notice that in a pruned graph, $\ell_i \neq \ell_j$ and $r_i \neq r_j$ hold for every $i \neq j$. We argue that we use gaps as small as possible. Only a right endpoint r_i following a left endpoint ℓ_j can be placed at the same position. The other case of a right endpoint r_i followed by a left endpoint ℓ_j requires a gap of size one; otherwise R_{v_i} would intersect R_{v_j} but $v_i v_j \notin E(G)$. So the gaps are minimal, we construct a smallest representation, and give the value minspan(C) correctly.

Case 2: A Located Component. We modify the above approach slightly to deal with located components. We already argued that there are at most two possible orderings \triangleleft (since the indistinguishable vertices are ordered by the partial representation), and we just test both of them. Both orderings can be used if and only if all pre-drawn vertices belong to one group of indistinguishable vertices. Then these two orderings give the same minspan(C) but the minimum representations might be differently shifted, and we are able to construct both of them. If the pre-drawn intervals do not belong to one group, the ordering \triangleleft is uniquely determined. (If it is compatible with the ordering of the pre-drawn intervals at all.)

We compute the common ordering \lessdot exactly as before and place the endpoints in this ordering. The only difference is that the endpoints of the pre-drawn intervals are prescribed. So we start at the position of the left-most pre-drawn endpoint ℓ_i . We place the endpoints smaller in \lessdot than ℓ_i on the left of ℓ_i as far to the right as possible. (We approach them in the reverse order exactly as above.) Then we proceed with the remaining endpoints in the order given by \lessdot . If the current endpoint is pre-drawn, we keep it as it is. Otherwise, we place it in the same way as above. The constructed representation is smallest and gives minspan(C).

Clique Orderings of P-in-P. Recall the properties of maximal cliques from Section 1.5. For a component C, we denote by cl(C) the number of maximal cliques of C. Let \mathcal{R} be a representation of C. Since the subtrees R_K corresponding to the maximal cliques are pairwise disjoint, they have to be ordered from left to right. This ordering has the following well-known property [18]:

Lemma 4.3 (Fulkerson and Gross). A graph is an interval graph if and only if there exists an ordering of the maximal cliques $K_1 < \cdots < K_{cl(C)}$ such that for each vertex v the cliques containing v appear consecutively in this ordering.

We quickly argue about the correctness of the lemma. Clearly, in an interval representation, all maximal cliques containing one vertex v appear consecutively. (Otherwise the clique in between would be intersected by R_v in addition.) On the other hand, having an ordering < of the maximal cliques from the statement, we can

construct a representation as follows. Assign a vertex p_i of T to each clique K_i , respecting the ordering <. For each vertex v, we assign $R_v = \{p_i \mid v \in K_i\}$. Since the maximal cliques containing v appear consecutively, each R_v is a subpath.

Minimum Spans of P-in-P. We again consider the minimum span defined exactly as for proper interval graphs above. Clearly, minspan $(C) \ge \operatorname{cl}(C)$. We show:

Lemma 4.4. For an unlocated component C of an interval graph, minspan $(C) = \operatorname{cl}(C)$. We can find a smallest representation in time $\mathcal{O}(n+m)$.

Proof. We start by identifying maximal cliques in time $\mathcal{O}(n+m)$, using the algorithm of Rose et al. [49]. To construct a smallest representation, we find an ordering from Lemma 4.3, using the PQ-tree algorithm [8] in time $\mathcal{O}(n+m)$. If such an ordering does not exist, the graph G is not an interval graph and no representation exists. If the ordering exists, we can construct a representation using exactly cl(C) vertices of the path as described above, by putting $R_v = \{p_i \mid v \in K_i\}$.

We note that this approach does not translate to located components, as in Lemma 4.2 for proper interval graphs. We prove in Corollary 4.11 that finding the minimum span for a located component is an NP-complete problem. (We prove this in the setting that the problem Repext(P-in-P, Add) is NP-complete. In the reduction, we ask whether a connected interval graph has the minimum span at most (M+1)k+1 for some integers k and M.)

4.1.2 The Polynomial Cases

First we deal with all polynomial cases.

Fixed Type Recognition. We just need to use the values of minimum spans we already know how to compute.

Proposition 4.5. Both RECOG*(PROPER P-in-P, FIXED) and RECOG*(P-in-P, FIXED) can be solved in time $\mathcal{O}(n+m)$.

Proof. We process the components C_1, \ldots, C_c one-by-one and place them on T' from left to right. If $\sum_{i=1}^c \operatorname{minspan}(C_i) \leq |T'|$, we can place the components using smallest representations from Lemma 4.2 for PROPER P-in-P, resp. Lemma 4.4 for P-in-P. Otherwise, the path is too small and a representation cannot be constructed.

Add Type Extension, PROPER P-in-P. Again, we approach this problem using minimum spans and Lemma 4.2.

Proposition 4.6. The problem REPEXT(PROPER P-in-P, ADD) can be solved in time $\mathcal{O}(n+m)$.

Proof. Since the path can be expanded to the left and to the right as much as necessary, we can place unlocated components far to the left. So we only need to deal with located components, ordered $C_1 \blacktriangleleft \cdots \blacktriangleleft C_c$ from left to right. We process the components from left to right. When we place C_i , it has to be placed on the right of C_{i-1} . We

have (at most) two possible smallest representations corresponding to two different orderings of C_i . We test whether at least one of them can be placed on the right of C_{i-1} , and pick the one minimizing the right-most vertex of T taken by C_i (leaving the maximum possible space for C_{i+1}, \ldots, C_c). If neither of the smallest representations can be placed, the extension algorithm outputs "no".

If the algorithm finishes, it constructs a correct representation. On the other hand, we place each component as far to the left as possible (while restricted by the previous components on the left). So if C_i cannot be placed, there exists no representation extending the partial representation.

Non-fixed Type Recognition. The only limitation for recognition of interval graphs inside a given path is the length of the path. In the three types Sub, ADD and BOTH, we can produce a path as long as necessary. (With the trivial exception $T' = P_0$ for Sub for which the instance is solvable if and only if $G = K_n$.) For a subpaths-in-path representation, the order of the endpoints of the subpaths from left to right is the only thing that matters, not the exact positions. In a tree T with at least 2n vertices, every possible ordering is realizable.

Thus the problems are equivalent to the standard recognition of interval graphs on the real line. The recognition can be solved in time $\mathcal{O}(n+m)$; see [40, 12] for PROPER P-in-P, and [8, 13] for P-in-P.

Both Type Extension. This extension type is equivalent with the partial representation extension problems of interval graphs on the real line. Again only the ordering of the endpoints is important. The only change here is that some of the endpoints are already placed. By subdividing, we can place any amount of the endpoints between any two endpoints (not sharing the same position). Also, the path can be extended to the left and to the right which allows to place any amount of endpoints to the left of the left-most pre-drawn endpoint and to the right of the right-most pre-drawn endpoint. So any extending ordering can be realized in the BOTH type.

The partial representation extension problem for interval graphs on the real line was first considered in [36]. The paper gives algorithms for both classes P-in-P and PROPER P-in-P, and does not explicitly deal with representations sharing endpoints but the algorithms are easy to modify. The results [7, 35, 33] show that both extension problems are solvable in time $\mathcal{O}(n+m)$.

Sub Type Extension. It is possible to modify the above algorithms for partial representation extension of P-in-P and PROPER P-in-P. Instead of describing details of these algorithms, we simply reduce the problems to the type BOTH which we can solve in time $\mathcal{O}(n+m)$ (as discussed above):

Theorem 4.7. Both RepExt(PROPER P-in-P, Sub) and RepExt(P-in-P, Sub) can be solved in time $\mathcal{O}(n+m)$.

The general idea is as follows. The difference between between Sub and Both is that for the Sub type, we cannot extend the path T' at the ends. Suppose that some pre-drawn subpath R'_v contains say the left endpoint of T'. Then R'_v contains

this endpoint also in T. So we are going to modify the graph G in such a way, that every representation of the BOTH type has to place everything on the right of R'_v .

Suppose first that the graph contains some unlocated components, and we show how to deal with them. We want to find one edge $p_i p_{i+1}$ of T' which we can subdivide many times and place all unlocated components in between of p_i and p_{i+1} in T. We call an edge $p_i p_{i+1}$ expandable if no located component C takes p_j and p_k such that $j \leq i < i + 1 \leq k$.

Lemma 4.8. Let G have at least one unlocated component, and let \widetilde{G} be the graph constructed from G by removing all unlocated components. Then \mathcal{R}' is extendible to \mathcal{R} if and only if T' contains at least one expandable edge $p_i p_{i+1}$ and \mathcal{R}' is extendible to $\widetilde{\mathcal{R}}$ of \widetilde{G} .

Proof. Let \mathcal{R}' be extendible to \mathcal{R} and let C be one unlocated component placed in T such that it takes a vertex in between of p_i and p_{i+1} of T'. Clearly \mathcal{R}' is extendible to $\widetilde{\mathcal{R}}$. And $p_i p_{i+1}$ is expandable since if there would be a located component \widetilde{C} taking p_j and p_k , then C would split \widetilde{C} , contradicting existence of \blacktriangleleft in \mathcal{R} ; recall the definition of \blacktriangleleft in Section 1.5.

For the other implication, we subdivide the expandable edge $p_i p_{i+1}$ many times such that we can place all unlocated components in this area. For located components, some of them have to be placed on the left of the unlocated components, and some on the right. We can subdivide all edges of T' enough to place the endpoints in the same order as in $\widetilde{\mathcal{R}}$. Thus we get \mathcal{R} extending \mathcal{R}' .

Theorem 4.7. We describe the reduction for P-in-P, and then we slightly modify it in the last paragraph for PROPER P-in-P. We deal with unlocated components using Lemma 4.8. We just need to check existence of an expandable edge for which we first compute the ordering \blacktriangleleft of the located components (if it doesn't exist, the partial representation is clearly not extendible). If there is exactly one located component C, then at least one of p_1 and p_t is not taken by C, and say for p_1 we obtain an expandable edge p_1p_2 . And if there are at least two located components $C_1 \blacktriangleleft \cdots \blacktriangleleft C_c$, let p_i be the right-most vertex taken by C_1 . Then p_ip_{i+1} is clearly expandable. It remains to deal with located components.

Let us consider the endpoint p_1 of T'. In BOTH, we can attach in T a path P of any length on the left of p_1 . If p_1 is not taken by C_1 , we can create in T the same path P by subdividing p_1p_2 . But if p_1 is taken by C_1 , we have to forbid P to be used in the construction of \mathcal{R} . We modify both the path T' and the graphs G, and we show that any representation \mathcal{R} extending \mathcal{R}' is realized in T in between of p_1 and p_t .

The modification is as follows. Let $v_1, \ldots, v_k \in C_1$ be all pre-drawn subpaths such that $p_1 \in R'_{v_1}, \ldots, R'_{v_k}$. First, we extend the path by one by adding p_0 attached to p_1 . We introduce an additional pre-drawn vertex v_{\leftarrow} adjacent exactly to v_1, \ldots, v_k in G. We put $R'_{v_{\leftarrow}} = \{p_0\}$ and we modify $R'_{v_i} = R'_{v_i} \cup \{p_0\}$. See Figure 4.3. Indeed, we proceed exactly the same on the other side of T'; if p_t is taken by C_c , we introduce p_{t+1} and v_{\rightarrow} .



Figure 4.3: The three pre-drawn subpaths containing p_1 are $R_{v_1} = \{p_1\}$, $R_{v_2} = \{p_1, p_2\}$ and $R_{v_3} = \{p_1, p_2, p_3\}$. We add p_0 to T' and to R_{v_1}, \ldots, R_{v_3} , and we introduce additional pre-drawn subpaths $R_{v_+} = \{p_0\}$.

We use the described algorithm for Repext(P-in-P, Both) for the modified graph and the modified path, which runs in time $\mathcal{O}(n+m)$. We obtain a representation \mathcal{R} extending \mathcal{R}' if it exists. If \mathcal{R} does not exist, then the original problem is clearly not solvable. It remains to argue that if \mathcal{R} exists, then we can either construct a solution for the original Sub type problem, or we can prove that it is not solvable.

We deal only with the left side of T; for the right side the argument is symmetrical. If T' is not modified on the left side, then the edge p_1p_2 can be subdivided as necessary and we are equivalent with the BOTH type. Suppose that p_0 is added. There are no unlocated components, and so everything with the exception of v_1, \ldots, v_k has to be represented on the right of v_{\leftarrow} which is placed on p_0 .

We need to argue the issue that the newly added edge p_0p_1 can be subdivided in T. There are the following two cases:

- Case 1. If $|R'_{v_i}| \geq 3$ for each i, i.e, p_1 and p_2 belong to each R'_{v_i} , the subdivision of p_0p_1 is equivalent to the subdivision of p_1p_2 which is correct in the original SUB type problem. So nothing needs to be done.
- Case 2. Let $|R'_{v_i}| = 2$ for some i, so $R'_{v_i} = \{p_0, p_1\}$. Then $N(v_i) \setminus v_{\leftarrow}$ has to form a complete subgraph of G, otherwise the starting partial representation having $R'_{v_i} = \{p_1\}$ would not be extendible. We revert the subdivision of p_0p_1 by modifying \mathcal{R} as follows. Let p'_1, \ldots, p'_s be the new vertices of T created by the subdivision of p_0p_1 . For each $v \in N(v_i) \setminus v_{\leftarrow}$, we set $R_v = R_v \setminus \{p'_1, \ldots, p'_s\} \cup \{p_1\}$, and we remove p'_1, \ldots, p'_s by contractions. Clearly, the resulting representation is correct and still extends \mathcal{R}' .

By removing p_0 and the vertices attached to it on the left, p_{t+1} and the vertices attached to it on the right, v_{\leftarrow} and v_{\rightarrow} (of course, only if they are added), we obtain a correct representation of G inside a subdivision of T' extending the partial representation \mathcal{R}' .

Concerning PROPER P-in-P, we use almost the same approach. The only difference is that we append two vertices p_0 and \bar{p}_0 (resp. p_{t+1} and \bar{p}_{t+1}) to the end of T', and we put $R'_{v_{\leftarrow}} = \{\bar{p}_0, p_0\}$ (resp. $R'_{v_{\rightarrow}} = \{p_{t+1}, \bar{p}_{t+1}\}$), so the modified partial representation is proper.

4.1.3 The NP-complete Cases

The basic gadgets of the reductions are paths. They have the following minimum spans.

Lemma 4.9. For the class P-in-P, minspan $(P_n) = n$. For the class PROPER P-in-P and $n \ge 2$, minspan $(P_n) = n + 2$.

Proof. For P-in-P, the number of the maximal cliques of P_n is n. For PROPER P-in-P, the ordering \leq is

$$\ell_0 \lessdot \ell_1 \lessdot r_0 \lessdot \ell_2 \lessdot r_1 \lessdot \cdots \lessdot \ell_i \lessdot r_{i-1} \lessdot \cdots \lessdot \ell_n \lessdot r_{n-1} \lessdot r_n.$$

There are n changes from ℓ_i to r_{i-1} and P_n has n+1 vertices. So the minimum span equals 2(n+1)-n=n+2.

We reduce the problems from 3-Partition. An input of 3-Partition consists of positive integers k, M and A_1, \ldots, A_{3k} such that $\frac{M}{4} < A_i < \frac{M}{2}$ for each A_i and $\sum A_i = kM$. It asks whether it is possible to partition A_i 's into k triples such that the sets A_i of each triple sum to exactly M. This problem is strongly NP-complete [20] which means that it is NP-complete even when the input is coded in unary, i.e., all integers are of polynomial sizes.

Theorem 4.10. Both REPEXT(PROPER P-in-P, FIXED) and REPEXT(P-in-P, FIXED) are NP-complete.

Proof. We use almost the same reductions for both PROPER P-in-P and P-in-P. For a given input of 3-Partition (with $M \ge 4$), we construct a graph G and its partial representation as follows.

As the fixed tree we choose $T' = P_{(M+1)k}$, with the vertices $p_0, \ldots, p_{(M+1)k}$. The graph G contains two types of gadgets as separate components. First, it contains k+1 split gadgets S_0, \ldots, S_k which split the path into k gaps of the size M. Then it contains 3k take gadgets T_1, \ldots, T_{3k} . A take gadget T_i takes in each representation at least A_i vertices of one of the k gaps.

For these reductions, the gadgets are particularly simple. The split gadget S_i is just a single pre-drawn vertex v_i with $R_{v_i} = \{p_{(M+1)i}\}$. The split gadgets clearly split the path into the k gaps of the size M. The take gadget T_i is P_{A_i} for P-in-P, resp. $P_{A_{i-2}}$ for PROPER P-in-P. According to Lemma 4.9, minspan $(T_i) = A_i$. The representation is extendible if and only if it is possible to place the take gadgets into the k gaps. For an example, see Figure 4.4. The reduction is clearly polynomial.

To conclude the proof, we show that the partial representation is extendible if and only if the corresponding 3-Partition input has a solution. If the partial

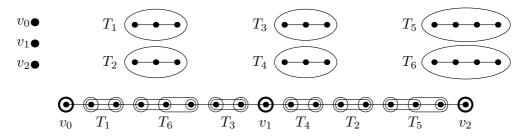


Figure 4.4: An example of the reduction for the following input of 3-Partition: k = 2, M = 7, $A_1 = A_2 = A_3 = A_4 = 2$ and $A_5 = A_6 = 3$. On top, the constructed interval graph is depicted. On bottom, the partial representation (depicted in bold) is extended.

¹Notice that if a subset of A_i 's sums to exactly M it has to be a triple due to the size constraints.

representation is extendible, the take gadgets T_i are divided into the k gaps on the path which gives a partition. Based on the constraints for the sizes of A_i 's, each gap contains exactly three take gadgets of the total minimum span M; thus the partition solves the 3-Partition problem. On the other hand, a solution of 3-Partition describes how to place the take gadgets into the k gaps and construct an extending representation.

Corollary 4.11. The problem Repext(P-in-P, Add) is NP-complete.

Proof. We use the above reduction for P-in-P with one additional pre-drawn interval v attached to everything in G. We put $R_v = \{p_0, \ldots, p_{(M+1)k}\}$, so it contains the whole tree T'. Since a representation of each take gadget T_i has to intersect R_v , it has to be placed inside of the k gaps as before.

We note that the above modification does not work for proper interval graphs. Indeed, this is not very surprising since Proposition 4.6 states that the problem REPEXT(PROPER P-in-P, ADD) can be solved in time $\mathcal{O}(n+m)$.

4.1.4 The Parameterized Complexity

In this subsection, we study the parameterized complexity. The parameters are the number c of components, the number k of pre-drawn intervals and the size t of the path T'.

By the Number of Components. In the reduction of Theorem 4.10, one might ask whether it is possible to make the reduction graph G connected. For P-in-P, it is indeed possible to add a universal vertex adjacent to everything in G, and thus make G connected as in the proof of Corollary 4.11. The following result answers this question for PROPER P-in-P negatively (unless P = NP):

Proposition 4.12. The problem Repext(PROPER P-in-P, Fixed) is fixed-parameter tractable in the number c of components, solvable in time $\mathcal{O}((n+m)c!)$.

Proof. There are c! possible orderings \blacktriangleleft of the components from left to right, and we test each of them. (The located components force some partial ordering \blacktriangleleft so we need to test less then c! orderings; see below the proof for details.) We show that for a prescribed ordering \blacktriangleleft of the components, we can solve the problem in time $\mathcal{O}(n+m)$; thus gaining the total time $\mathcal{O}((n+m)c!)$. We solve the problem almost the same as in the proof of Proposition 4.6. The only difference is that we deal with all components instead of only the located ones.

We process the components from left to right. When we process C_i , we place it on the right of C_{i-1} as far to the left as possible. For the unlocated C_i , we can take any smallest representation. For the located C_i , we test both smallest representations and take the one placing the right-most endpoint of C_i further to the left. We construct the representation in time $\mathcal{O}(n+m)$. For the correctness of the algorithm see the proof of Proposition 4.6 for more details.

We note that for NP-hardness of the problem Repext(PROPER P-in-P, Fixed) it is necessary to have some pre-drawn subpaths. On the other hand, also some unlocated components are necessary. If all the components were located, there would be a unique ordering \triangleleft and we could test it in time $\mathcal{O}(n+m)$ as described above. In general, for c components and c' located components, we need to test only $\frac{c!}{c'!}$ different orderings.

By the Number of Pre-drawn Intervals. In the reduction in Theorem 4.10, we need to have k pre-drawn intervals. One could ask, whether the problems become simpler with a small number of pre-drawn intervals. We answer this negatively. For PROPER P-in-P, the problem is in XP and W[1]-hard with respect to k. For P-in-P, we only show that it is W[1]-hard.

There are two closely related problems BINPACKING and GENBINPACKING. In both problems, we have k bins and n items of positive integer sizes. The question is whether we can pack (partition) these items into the k bins when the volumes of the bins are limited. For BINPACKING, all the bins have the same volume. For GENBINPACKING, the bins have different volumes. Formally:

Problem: BINPACKING

Input: Positive integers k, ℓ, V , and A_1, \ldots, A_{ℓ} .

Question: Does there exist a k-partition $\mathcal{P}_1, \ldots, \mathcal{P}_k$ of A_1, \ldots, A_ℓ such

that $\sum_{A_i \in \mathcal{P}_i} A_i \leq V$ for every \mathcal{P}_j .

Problem: GENBINPACKING

Input: Positive integers $k, \ell, V_1, \ldots, V_k$, and A_1, \ldots, A_{ℓ} .

Question: Does there exist a k-partition $\mathcal{P}_1, \ldots, \mathcal{P}_k$ of A_1, \ldots, A_ℓ such

that $\sum_{A_i \in \mathcal{P}_i} A_i \leq V_j$ for every \mathcal{P}_j .

Lemma 4.13. The problems BINPACKING and GENBINPACKING are polynomially equivalent.

Proof. Obviously BINPACKING is a special case of GENBINPACKING. On the other hand, let $k, \ell, V_1, \ldots, V_k$, and A_1, \ldots, A_ℓ be an instance of GENBINPACKING. We construct an instance k', ℓ', V' , and $A'_1, \ldots, A'_{\ell'}$ of BINPACKING as follows. We put $k' = k, \ell' = \ell + k$ and $V' = 2 \cdot \max V_i + 1$. The sizes of the first ℓ items are the same, i.e, $A'_i = A_i$ for $i = 1, \ldots, \ell$. The additional items $A'_{\ell+1}, \ldots, A'_{\ell+k}$ are called *large* and we put $A'_{\ell+i} = V' - V_i$ for $i = 1, \ldots, k$.

Each bin has to contain exactly one large item since two large items take more space than V'. After placing large items into the bins, we obtain the bins of the remaining volumes V_1, \ldots, V_k in which we have to place the remaining items. This corresponds exactly to the original GenbinPacking instance.

If the sizes of items are encoded in binary, the problem is NP-complete even for k=2. The more interesting version which we use here is that the sizes are encoded in unary so all sizes are polynomial. In such a case, the BinPacking problem is known to be solvable in time $t^{\mathcal{O}(k)}$ using dynamic programming where t is the total size of all

items. And it is W[1]-hard with respect to the parameter k [30]. The similar holds for REPEXT(PROPER P-in-P, FIXED):

Theorem 1.7. For a given instance of the BINPACKING problem, we can solve it by REPEXT(PROPER P-in-P, FIXED) in a similar manner as in the reduction in Theorem 4.10. As T', take a path $P_{(V+1)k}$. As G, take P_{A_i-2} for each A_i and the pre-drawn vertices v_0, \ldots, v_k such that $R_{v_i} = \{p_{(V+1)i}\}$. The rest of the argument is exactly as in the proof of Theorem 4.10.

Now, we want to solve Repext(Proper P-in-P, Fixed) using 2^k instances of GenbinPacking (which is polynomially equivalent to BinPacking), where k is the number of pre-drawn intervals.

First we deal with located components $C_1 \blacktriangleleft \cdots \blacktriangleleft C_c$. For each component, we have two possible orderings \triangleleft and using Lemma 4.2 we get (at most) two possible smallest representations which might be differently shifted. In total, we have at most $2^c \le 2^k$ possible representations keeping C_1, \ldots, C_c as small as possible leaving maximal gaps for unlocated components. We test each of these 2^c representations.

Let $C'_1, \ldots, C'_{c'}$ be the unlocated components. For each unlocated component C'_i , we compute minspan (C'_i) using Lemma 4.2. The goal is to place the unlocated components into the c+1 gaps between representations of the located components C_1, \ldots, C_c . We can solve this problem using GENBINPACKING as follows. We have k+1 bins of the volumes equal to the sizes of the gaps between the representations of C_1, \ldots, C_c . We have c' items of the sizes $A_i = \text{minspan}(C'_i)$.

A solution of GenbinPacking tells how to place the unlocated components into the k gaps. If there exists no solution, this specific representation of the located components cannot be used. We can test all 2^c possible representations of the located components. Thus we get the required weak truth-table reduction.

Corollary 4.14. The problem REPEXT(PROPER P-in-P, FIXED) is W[1]-hard and belongs to XP, solvable in time $n^{\mathcal{O}(k)}$ where k is the number of pre-drawn intervals.

Proof. Both claims follow from Theorem 1.7.

Proposition 4.15. The problems REPEXT(P-in-P, FIXED) and REPEXT(P-in-P, ADD) are W[1]-hard when parameterized by the number k of pre-drawn intervals.

Proof. We modify the reductions of Theorem 4.10 and Corollary 4.11 exactly as in the proof of Theorem 1.7. \Box

By the Size of the Path. We show that the FIXED type problems are fixed-parameter tractable with respect to the size of the path t. It is easy to find a solution by a brute-force algorithm:

Proposition 4.16. Let t be the size of the tree and let $f(t) = t^{2t^2}$. Then the problems REPEXT(PROPER P-in-P, FIXED) and REPEXT(P-in-P, FIXED) are FPT with respect to the parameter t, solvable in time $\mathcal{O}(n+m+f(t))$.

Proof. In a pruned graph, the vertices have to be represented by pairwise different intervals. There are at most t^2 possible different subpaths of a path with t vertices so the pruned graph can contain at most t^2 vertices; otherwise the extension is clearly not possible. We can test every possible assignment of the non-pre-drawn vertices to the t^2 subpaths, and for each assignment we test whether we get a correct representation extending \mathcal{R}' .

4.2 Path and Chordal Graphs

We present and prove the results concerning the classes P-in-T and T-in-T.

4.2.1 The Polynomial Cases

The recognition problems for the types ADD and BOTH are equivalent to standard recognition without any specified tree T'. Indeed, we can modify T' by adding an arbitrary tree to it. If the input graph is P-in-T or T-in-T, there exists a tree T'' in which the graph can be represented. We produce T by attaching T'' to T' in any way. Then the input graph can be represented in T as well, completely ignoring the part T'.

For path graphs, the original recognition algorithm is due to Gavril [21] in time $\mathcal{O}(n^4)$. The current fastest algorithm is by Schäffer [51] in time $\mathcal{O}(nm)$. For chordal graphs, there is a beautiful simple algorithm by Rose et al. [49] in time $\mathcal{O}(n+m)$.

4.2.2 The NP-complete Cases

All the remaining cases from the table of Figure 1.7 are NP-complete. We modify the reduction for P-in-P of Theorem 4.10. We start with the simplest reduction for the FIXED type and then modify it for the other types.

Fixed Type Recognition. For the FIXED type, we can avoid pre-drawn subtrees, using an additional structure of the tree.

Proposition 4.17. The problems RECOG*(P-in-T, FIXED) and RECOG*(T-in-T, FIXED) are NP-complete.

Proof. We again reduce from 3-Partition with an input k and M. For technical purposes, let $M \geq 8$ and so $|A_i| > 2$ for each A_i . We construct a graph G and a tree T' as follows.

The tree T' is a path $P_{(M+1)k}$ (its vertices being denoted by $p_0, \ldots, p_{(M+1)k}$) with three paths of length two attached to every vertex $p_{(M+1)i}$, for each $i=0,\ldots,k$; see Figure 4.5. Each split gadget S_i is a star, depicted on the left of Figure 4.5. When the split gadgets are placed as in T', they split the tree into k gaps exactly as the pre-drawn vertices in the proof of Theorem 4.10. Each take gadget T_i is the path P_{A_i} exactly as before. The reduction is obviously polynomial.

What remains to argue is the correctness of the reduction. Observe that given a solution of 3-Partition, we can construct a subpaths-in-tree representation of G

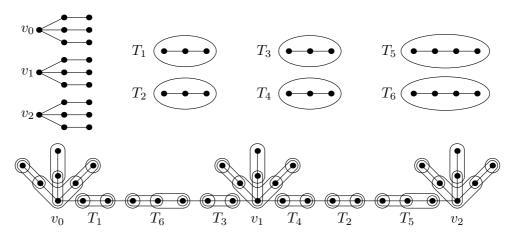


Figure 4.5: An example for the same input of 3-Partition as in Figure 4.4. On top the graph G is depicted. On bottom, a representation of G is constructed, giving the solution $\{A_1, A_3, A_6\}$ and $\{A_2, A_4, A_5\}$.

as in Figure 4.5. For the other direction, let v_0, \ldots, v_k be the central vertices of the split gadgets S_0, \ldots, S_k . We claim that each R_{v_i} contains at least one branch vertex. (Actually, exactly one since there are only n+1 branch vertices in T'.) If some R_{v_i} contained only non-branch vertices, then it would not be possible to represent three disjoint neighbors u_1, u_2 and u_3 of this v_i having each $R_{u_i} \setminus R_{v_i}$ non-empty.

Since each branch vertex is taken by one R_{v_i} , the path $P_{(M+1)k}$ is split into k gaps as before. Since $|A_i| > 2$, each T_i can be represented only inside of these gaps. Notice that the total number of the vertices in the gaps has to be equal kM, and therefore the split gadgets have to be represented entirely in the attached stars as in Figure 4.5. The rest of the reduction works exactly as in Theorem 4.10.

Sub Type Recognition. By modifying the above reduction, we get:

Theorem 4.18. The problems Recog*(P-in-T, Sub) and Recog*(T-in-T, Sub) are NP-complete.

Proof. We need to modify the two gadgets from the reduction of Theorem 4.17 in such a way that a subdivision of the tree T' does not help in placing them. Subdivision only increases the number of non-branch vertices. Thus a take gadget T_i requires A_i branch vertices. Similarly, the split gadget S_i is more complicated. See Figure 4.6 on top.

The tree T is constructed as follows. We start with a path $P_{(M+1)k}$ with vertices $p_0, \ldots, p_{(M+1)k}$. To each vertex $p_{(M+1)i}$ we attach a subtree isomorphic to the trees in Figure 4.6 on bottom. To the remaining vertices of the path, we attach one leaf per vertex. The reduction is again polynomial.

Straightforwardly, for a given solution of 3-Partition, we can construct a correct subpaths-in-tree representation in a subdivided tree. On the other hand, we are going to show how to construct a solution of 3-Partition from a given tree representation.

Recall the properties of maximal cliques from Section 1.5. Note that each triangle $u_1u_2u_3$ in each split or take gadget is a maximal clique K. Since $N[u_i] \neq N[u_j]$ for

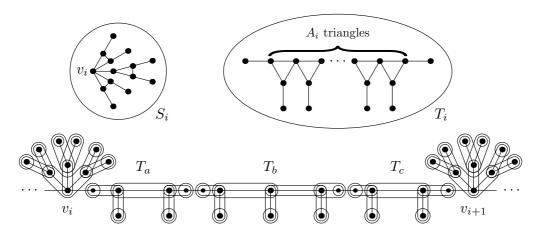


Figure 4.6: On top, the split gadget S_i is on the left and the take gadget T_i is on the right. On bottom, a part of the tree T is depicted with the small vertices added by subdivisions. The gap between two split gadgets contains three take gadgets T_a , T_b and T_c giving one triple $\{A_a, A_b, A_c\}$ with $A_a + A_b + A_c = M$.

each $i \neq j$, there has to be a branch vertex in $R_{u_i} \cap R_{u_j}$ for some i and j. The gadget S_i contains three triangles, each taking one branch vertex of T. In addition, R_{v_i} connecting them has to contain another branch vertex. So in total, S_i contains at least four branch vertices. Each gadget T_i contains A_i triangles, and so it requires at least A_i branch vertices. Since the number of branch vertices of T is limited, each S_i takes exactly four branch vertices and each T_i takes exactly A_i branch vertices.

Now, if some T_i contained a branch vertex of the subtrees attached to $p_{(M+1)j}$, at least one of its branch vertices would not be used. (Either not taken by T_i , or T_i would require at least $A_i + 1$ branch vertices.) So each S_i has to take the branch vertices of the subtrees attached to $p_{(M+1)j}$ for some j, and the take gadgets have to be placed inside the gaps exactly as before.

Proposition 4.19. Even with a single pre-drawn subtree, i.e, |G'| = 1, the problems Repext(T-in-T, Add) and Repext(T-in-T, Both) are NP-complete.

Proof. We easily modify the above reductions; for the ADD type, the reduction of Proposition 4.17, for the BOTH type, the reduction of Theorem 4.18. The modification adds into G one pre-drawn vertex v adjacent to everything such that $R_v = T'$. Since R_v spans the whole tree, it forces the entire representation \mathcal{R} into T'.

We just deal with the ADD type, for BOTH the argument is exactly the same. Let T' be the partial tree and let T be the tree in which the representation is constructed, so T' is a subtree of T. We claim that we can restrict a representation of each vertex of G into T' and thus obtain a correct representation inside the subtree T'.

Let $x \in V$. Since $xv \in E(G)$, the intersection of R_x and T' is a non-empty subtree. We put $\widetilde{R}_x = R_x \cap T'$, and we claim that $\widetilde{\mathcal{R}}$ is a representation of G in T'. To argue the correctness, let x and y be two different vertices from v (otherwise trivial). If $xy \notin E(G)$, then $R_x \cap R_y = \emptyset$, and so $\widetilde{R}_x \cap \widetilde{R}_y = \emptyset$ as well. Otherwise xyv is a triangle in G, and thus by the Helly property the subtrees R_x , R_y and $R_v = T'$ have a non-empty common intersection, giving that $\widetilde{R}_x \cap \widetilde{R}_y$ is non-empty.

For path graphs, one can use a similar technique of a pre-drawn universal vertex attached to everything. But there is the following difficulty: To do so, the input partial tree T' has to be a path. For the type BOTH, the complexity of Repext(P-in-T, Both) remains open. For the type ADD, we get the following weaker result:

Proposition 4.20. The problem Repext(P-in-T, Add) is NP-complete.

Proof. Similarly as in Proposition 4.19, add a pre-drawn universal vertex v on the path T' constructed in the reduction of Theorem 4.10 such that $R_v = T'$. The rest is exactly as above.

4.2.3 The Parameterized Complexity

We deal with parameterized complexity of the problems and we give only minor and partial results in this direction. Unlike in Section 4.1, parameterization by the number k of pre-drawn subtrees is mostly not helpful. We show that every problem with exception of Repext(P-in-T, Add) is already NP-complete for k=0 or k=1. For Repext(P-in-T, Add), we have only a weaker result that it is W[1]-hard with respect to the parameter k since Proposition 4.15 straightforwardly generalizes.

Similarly, a low number c of components does not make the problem any easier. We can easily insert a universal vertex attached to everything. So the above reductions can be modified and the problems remain $\mathsf{NP}\text{-}\mathsf{complete}$ even if the graph G is connected.

Concerning the size t of the tree, Proposition 4.16 straightforwardly generalizes:

Proposition 4.21. Let t be the size of T'. The problems Repext(P-in-T, Fixed) and Repext(T-in-T, Fixed) are fixed-parameter tractable with respect to t. They can be solved in time $\mathcal{O}(n+m+g(t))$ where

$$g(t) = 2^{t2^t}.$$

Proof. Proceed exactly as in the proof of Proposition 4.16, test all possible assignments of all vertices of a pruned graph. The only difference is that T' has at most 2^t different subtrees.

We note that a more precise bound for the number of subtrees could be use but we did not try to better estimate the function g.

5 Conclusions

In this thesis, we show recent developments in the study of the partial representation extension problem. We deal with the classes of interval graphs, proper interval graphs, unit interval graphs and chordal graphs. We describe concepts and techniques common to these algorithms, and we believe that they can be applied to other classes as well.

We conclude this thesis by describing two related problems to the partial representation extension problem. Also, we give some open problems.

5.1 Simultaneous Representations

In the introduction, we give a formal definition of the simultaneous representations problem, considered recently by Jampani and Lubiw [29, 28]. To recall the problem, the input gives graphs G_1, \ldots, G_k and the problem asks whether there exist representations $\mathcal{R}^1, \ldots, \mathcal{R}^k$ assigning the same sets to the common vertices I. For a class \mathcal{C} , we denote this problem by SIMREP(\mathcal{C}).

Here, we show that for many classes \mathcal{C} , the simultaneous representations problem is closely related to the partial representation extension problem, and many techniques useful for one problem can be applied to the other problem as well. This relation is more of a general principle than a precise mathematical statement.

Partial Representation Extension 'solves' Simultaneous Representations. Let the size of I be small. If there exists only a small number of different representations of the subgraph induced by the vertices of I, we can test all of them and solve SIMREP using Repext. So if the Repext problem is solvable in polynomial time, we can get a fixed-parameter tractable algorithm with parameter i = |I|. We show this for interval graphs:

Proposition 5.1. We can solve SIMREP(INT) in time $\mathcal{O}((n+m)\cdot(2i)!)$, where n is the total number of vertices of all graphs and m is the total number of edges of all graphs.

Proof. If a representation \mathcal{R}_I of I would be given, we could use the algorithm for REPEXT(INT) to test whether it is possible to extend it to a simultaneous representation of all the graphs. We just need to test for every graph G_j whether it is possible to extend the partial representation \mathcal{R}_I to a representation of entire G_j . This can be done in time $\mathcal{O}(n+m)$, using Theorem 1.1.

Since only the left-to-right order of endpoints is important, an interval graph with i vertices has $\mathcal{O}((2i)!)$ topologically different representations. Therefore, we can test all possible representation of I and get the running time $\mathcal{O}((n+m)\cdot(2i)!)$.

Similar relations holds for some other classes, for example proper interval graphs or circle graphs [9]. We note that this is currently the best known algorithm for SIMREP(INT) when k > 2, no polynomial-time algorithm is known. For k = 2, Jampani and Lubiw [29] give an algorithm in time $\mathcal{O}(n^2 \log n)$, and a recent paper of Bläsius and Rutter [7] improves the running time to $\mathcal{O}(n+m)$. In the case of circle graphs, Chaplick et al. [9] show that the simultaneous representations problem is NP-complete when k is a part of the input.

Simultaneous Representations "solves" Partial Representation Extension. In the opposite way, we are sometimes able to solve Repext by Simrer, which was suggested to us by Lubiw. Suppose that we have a graph G and a partial representation \mathcal{R}' of G'. We construct an instance of Simrer for k=2 as follows: I=G', $G_1=G$ and G_2 contains additional auxiliary vertices. The purpose of these auxiliary vertices is to fix a representation of I to be topologically equivalent to the partial representation \mathcal{R}' . Therefore, the solution \mathcal{R}^1 and \mathcal{R}^2 of the simultaneous representations problem gives a topologically correct representation \mathcal{R}^1 extending \mathcal{R}' .

For example in the case of interval graphs, we construct G_2 as follows. Consider the partial representation and add a long path consisting of short intervals on top of it. These intervals correspond to the auxiliary vertices added to G_2 and force a grid-like structure on the real line. This grid forces I to be represented topologically the same as in the partial representation \mathcal{R}' . Thus we can solve Repext(INT) using SIMREP(INT) which is used in [7] to obtain an algorithm solving Repext(INT) in time $\mathcal{O}(n+m)$.

We note that this relation does not always work. First, it might not be possible to construct such a graph G_2 , forcing a topologically unique representation of I. For example, the problem of simultaneous representations of chordal graphs is solvable in a polynomial time [28]. But the corresponding partial representation extension problem, considered in Chapter 4, is NP-complete; see Proposition 4.19. Second, it might happen that the topological equivalency of the representations in not sufficient, for example for partially represented unit interval graphs.

Applying Our Techniques. Concerning the classes PROPER INT and UNIT INT, the simultaneous representations problem similarly distinguishes them. One can easily construct two graphs which have simultaneous proper interval representations but not simultaneous unit interval representations. We believe that many techniques developed in Chapter 3 can be applied for the SIMREP(PROPER INT) and SIMREP(UNIT INT) problems. We already have some partial results in this direction which are not yet written, and we just sketch our ideas here.

First, one needs to construct simultaneous left-to-right orderings $<_1, \ldots, <_k$ having the same order on I. If these orderings do not exist, the simultaneous representations cannot be constructed. Using these orderings, we can easily construct simultaneous proper interval representations. For unit interval graphs, we can additionally

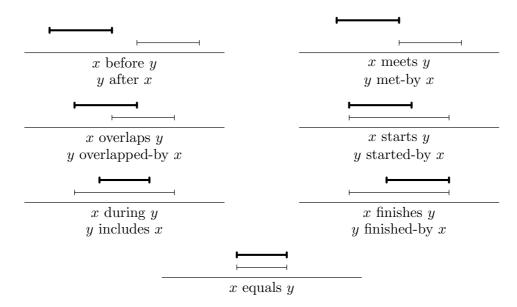


Figure 5.1: Thirteen primitive relations between the thick interval x and the thin interval y.

apply linear programming or some shifting approach.

There are two main problems for which we do not know a solution yet. We are able to construct orderings $<_1, \ldots, <_k$ only if the graphs G_1, \ldots, G_k are connected. If they are not connected, the construction of the orderings $<_1, \ldots, <_k$ is not clear. Second, for shifting algorithm, we need to choose some initial representation, and it is not clear how many shifting iterations are necessary.

5.2 Allen Algebras

The following type of problems is studied in theory of artificial intelligence and time reasoning; see Golumbic [22] for a survey. Suppose that we have several events which happened at some time. To every event, we can assign an interval of the timeline. Now, for some pairs of events we know relations. Allowing shared endpoints, Allen [1] characterized thirteen *primitive relations* between the events, see Figure 5.1. A relation is a set of several primitive relations. For some pairs of events, we specify relations in which they can occur. For example, we can specify for events x and y that either x is before y, or x is during y.

We want to find intervals representing the events such that all the specified relations are satisfied. This is called the *interval satisfiability problem*. Vilian and Kautz [56] proved that this problem is NP-complete. Golumbic and Shamir [24] gave a more simple proof, using the interval graph sandwich problem.

Notice that we can describe Repext(INT) and Repext(Proper Int) in the settings of Allen Algebras. In the case of Repext(Int), we do it in the following way. If vertices are non-adjacent, we assign them the relation {before, after}. If they are adjacent, we assign the complement of the previous relation. For pairs of pre-drawn intervals, we give singleton relations. Similarly, we can specify Repext(Proper Int).

But these more general problem are NP-complete, so they do not help in solving

REPEXT(INT) and REPEXT(PROPER INT). On the other hand, the results presented in this thesis show that some very specific problems concerning Allen Algebras are polynomially solvable.

5.3 Open Problems

We state several open problems.

Different Approach? The formulation of the first problem is not very precise. As we already described in the introduction, all the algorithm for solving partial representation extension problems we know are based on the following principle. First, we use some (known) way how to find all possible representations of an input graph. Then we derive some necessary conditions from the partial representations. Moreover, we show that these conditions are sufficient. Then, we test these conditions on all possible representations of the graph. If some representation satisfies them, we can use this representation to extend the input partial representation.

In the case of INT, we use PQ-trees. For PROPER INT and UNIT INT, we use uniqueness of the representation. The algorithms for extension of comparability, permutation and function graphs is based on properties of modular decomposition [32]. Also, the extension algorithm for planar graphs [2] uses SPQR-trees to work with all possible representations of the given planar graph.

Question 5.2. Is it possible to solve some Repext problem more "directly", without "testing" all possible representations?

Faster Algorithm for Unit Interval Graphs. Concerning Chapter 3, we give only one open problem:

Problem 5.3. Is it possible to solve Repext(UNIT INT) in time $\mathcal{O}(r)$, where r is the size of the input?

Two Problems for Chapter 4. One of the main goals of Chapter 4 is to stimulate future research in this area. Therefore, we give two open problems.

The first problems concerns the only open case in the table in Figure 1.7.

Problem 5.4. What is complexity of Repext(P-in-T, Both)?

Concerning parametrized complexity, we believe it is useful to first attack problems related to interval graphs. This allows to develop tools for more complicated chordal graphs. A generalization of Theorem 1.7 and Corollary 4.14 for P-in-P seems to be particularly interesting. The PQ-tree approach seems to be a good starting point.

Problem 5.5. Does Repext(P-in-P, Fixed) belong to XP with respect to k where k is the number of pre-drawn intervals?

Algorithms

In this appendix, we give pseudocodes of the main algorithms of this thesis. These pseudocodes are not very detail and not written formally in any programming language. They are meant to be overviews of all important steps done in each algorithm.

A.1 Reordering PQ-tree, General Partial Order

The pseudocode is in Algorithm 1. It is described in Section 2.1.1.

Algorithm 1 Reordering a PQ-tree – REORDER (T, \lhd)

Require: A PQ-tree T and a partial ordering \triangleleft .

Ensure: A reordering T' of T such that $<_{T'}$ extends < if it exists.

- 1: Construct the digraph of \triangleleft .
- 2: Process the nodes of T from bottom to the root:
- 3: for a processed node N do
- 4: Consider the subdigraph induced by the children of N.
- 5: **if** the node N is a P-node **then**
- 6: Find a topological sort of the subdigraph.
- 7: If it exists, reorder N according to it, otherwise output "no".
- 8: **else if** the node N is a Q-node then
- 9: Test whether the current ordering or its reversal is compatible with the subdigraph.
- 10: If yes, reorder the node, otherwise output "no".
- 11: **end if**
- 12: Contract the subdigraph into a single vertex.
- 13: end for
- 14: **return** A reordering T' of T.

A.2 Reordering PQ-tree, Interval Order

The pseudocode is in Algorithm 2. The description is in Section 2.1.2.

Algorithm 2 Reordering a PQ-tree, interval order – Reorder (T, \triangleleft)

Require: A PQ-tree T and an interval order \triangleleft with a sorted representation.

Ensure: A reordering T' of T such that $<_{T'}$ extends \triangleleft if it exists.

- 1: Calculate the handles for each invididual leaf of T and initiate an empty list for each endpoint.
- 2: Process the nodes of T from the bottom to the root:
- 3: for a processed node N do
- 4: Compute the handles of N using (2.3).
- 5: Add the node N to the lists of the two endpoints which are the handles of N.
- 6: end for
- 7: Iterate the sorted representation and construct the ordering $\stackrel{\sim}{\lessdot}$ for each inner node N.
- 8: Again process the nodes from bottom to the root:
- 9: for a processed node N with the ordering $\stackrel{\sim}{\lessdot}$ do
- 10: **if** the node N is a P-node **then**
- 11: Find any topological sort by removing minimal elements from \leq .
- 12: If it exists, reorder N according to it, otherwise output "no".
- 13: **else if** the node N is a Q-node **then**
- 14: Test whether the current ordering or its reversal is a topological sort.
- Process the prescribed topological sort from left to right, check for every element whether it is minimal and remove its handles from \leq .
- 16: If at least one ordering is correct, reorder the node, otherwise output "no".
- 17: **end if**
- 18: end for
- 19: **return** A reordering T' of T.

A.3 Extending Interval Graphs

The pseudocode is in Algorithm 3. It is described in Section 2.2.

Algorithm 3 Extending Interval Graphs – Repext(INT)

Require: An interval graph G and a partial representation \mathcal{R}' .

- 1: Compute maximal cliques and construct a PQ-tree.
- 2: Sweep \mathcal{R}' from left to right and construct the sorted representation of \triangleleft .
- 3: Use Algorithm 2 to reorder the PQ-tree according \triangleleft .
- 4: If any of these steps fails, no representation exists and output "no".
- 5: Place the clique-point according to the ordering $\langle T' \rangle$ from left to right:
- 6: for a clique-point cp(a) placed after cp(b) do
- 7: Compute the infimum of all points of the real line on the right of cp(b) where cp(a) can be placed.
- 8: If there is single such point, place cp(a) there.
- 9: Otherwise place cp(a) by ε on the right of the infimum, where ε is the size of the smallest part divided by n.
- 10: end for
- 11: Construct \mathcal{R} for the remaining intervals on top of the placed clique-points.
- 12: **return** A representation \mathcal{R} extending \mathcal{R}' .

A.4 Extending Proper Interval Graphs

The pseudocode is in Algorithm 4. The description is in Section 3.1.

Algorithm 4 Extending Proper Interval Graphs – Repext(PROPER INT)

Require: A proper interval graph G and a partial representation \mathcal{R}' .

- 1: Compute the pruned graph.
- 2: if pre-drawn intervals of some component do not appear consecutively then
- 3: Return "no", \mathcal{R}' is not extendible.
- 4: end if
- 5: Assign to the components pairwise disjoint open segments containing their predrawn intervals.
- 6: **for** each component **do**
- 7: Construct the partial orderings < and $<_{G'}$.
- 8: **if** there exists an ordering \triangleleft extending $\triangleleft_{G'}$, and \triangleleft or its reversal, and all touching pairs of intervals satisfy condition (2) **then**
- 9: Construct a representation of the component in its open segment as follows.
- 10: Compute a common ordering \leq , starting with $\ell_1 \leq \cdots \leq \ell_n$:
- 11: **for** each r_i **do**
- 12: Insert r_i right before ℓ_j such that v_j is the first non-neighbor of v_i on the right (or append r_i to the end, if v_j does not exist).
- 13: end for
- 14: Using *≤*, construct a representation of the component by placing non-predrawn endpoints equidistantly into the gaps.
- 15: **else**
- 16: Return "no", \mathcal{R}' is not extendible.
- 17: **end if**
- 18: end for
- 19: **return** A representation \mathcal{R} extending \mathcal{R}' .

A.5 Bounded Representation - LP Approach

The pseudocode is in Algorithm 5. The description is in Section 3.3.1.

Algorithm 5 Bounded Representation of Unit Interval Graphs – BOUNDREP

Require: A unit interval graph G, bound contraints and ordering \triangleleft .

Ensure: A representation \mathcal{R} in the ordering \triangleleft satisfying the bounds if it exists.

1: Compute the value ε according to the bounds $\frac{p_1}{q_1}, \ldots, \frac{p_b}{q_b}$:

$$\varepsilon' := \frac{1}{\operatorname{lcm}(q_1, q_2, \dots, q_b)}, \quad \text{and} \quad \varepsilon := \frac{\varepsilon'}{n}.$$

- 2: Process the components C_1, \ldots, C_c according to \blacktriangleleft .
- 3: **for** a component C_t **do**
- 4: Find the ordering <.
- 5: Construct linear orderings \triangleleft for < and its reversal.
- 6: For each ordering, solve one linear program using the previous value E_{t-1} .
- 7: **if** at least one program has a solution **then**
- 8: Use the solution minimizing the value of E_t of the linear program.
- 9: **else**
- 10: No representation exists, output "no".
- 11: end if
- 12: end for
- 13: **return** A representation \mathcal{R} in the ordering \triangleleft satisfying the bounds.

A.6 Extending Unit Interval Graphs

The pseudocode is in Algorithm 6. The description is in Section 3.4.

Algorithm 6 Extending Unit Interval Graphs – Repext(UNIT INT)

Require: A unit interval graph G and a partial representation \mathcal{R}' .

- 1: Deal with unlocated components separately, using the standard algorithm.
- 2: Construct the ordering ◀ for located components.
- 3: Set bounds exactly for each pre-drawn interval v_i at position ℓ_i , set

$$lbound(v_i) = ubound(v_i) = \ell_i.$$

- 4: Solve the located components using the bounded representation problem.
- 5: If no bounded representation exists, output "no".
- 6: **return** A representation \mathcal{R} extending \mathcal{R}' .

A.7 Bounded Representation – Shifting Algorithm

The pseudocode is in Algorithm 7. It is described in Section 3.3.5.

Algorithm 7 Bounded Representation of Unit Interval Graphs – BOUNDREP

Require: A unit interval graph G, bound contraints and ordering \triangleleft .

Ensure: A representation \mathcal{R} in the ordering \triangleleft satisfying the bounds if it exists.

- 1: Constuct a pruned graph, as described in Section 3.3.4.
- 2: Compute the value ε according to the bounds $\frac{p_1}{q_1}, \ldots, \frac{p_b}{q_b}$:

$$\varepsilon' := \frac{1}{\operatorname{lcm}(q_1, q_2, \dots, q_b)}, \quad \text{and} \quad \varepsilon := \frac{\varepsilon'}{n^2}.$$

- 3: Process the components C_1, \ldots, C_c according to \blacktriangleleft .
- 4: for a component C_t do
- 5: Find the ordering \triangleleft (which is the same as < before).
- 6: For \triangleleft and its reversal, construct the left-most representation (with additional lower bound E_{t-1}).
- 7: Precompute rounded lower bounds.
- 8: Construct an initial representation using [12].
- 9: Proceed the first phase using the LEFTSHIFTprocedure.
- 10: Proceed the second phase using the LEFTSHIFTprocedure.
- 11: **for** each LeftShift(v_i) **do**
- 12: **if** $\bar{\ell}_i$ is the strict maximum **then**
- 13: Shift v_i to position ℓ_i .
- 14: else
- 15: The interval v_i becomes fixed.
- 16: Compute precise and rounded positions of v_i .
- 17: Remove β_i from the position cycle.
- 18: end if
- 19: end for
- 20: **if** at least one left-most representation satisfies the upper bounds **then**
- 21: Use the left-most representation minimizing the value of E_t .
- 22: else
- 23: No representation exists, output "no".
- 24: end if
- 25: end for
- 26: **return** A representation \mathcal{R} in the ordering \triangleleft satisfying the bounds.

A.8 Subpath-in-Path Graphs - Sub Type Extension

The pseudocode is in Algorithm 8. The description is in Section 4.1.2.

Algorithm 8 Sub Type Extension of (Proper) Subpath-in-Path Graphs

Require: A (proper) subpath-in-path graph G and a partial representation \mathcal{R}' .

Ensure: A representation \mathcal{R} extending \mathcal{R}' if it exists.

- 1: if unlocated components can be placed then
- 2: Place them and ignore them in the rest of the algorithm.
- 3: else
- 4: No representation exists, output "no".
- 5: end if
- 6: if some subpath contains p_1 then
- 7: Add p_0 , v_{\leftarrow} (and \bar{p}_0 for PROPER P-in-P).
- 8: end if
- 9: if some subpath contains p_t then
- 10: Add p_{t+1} , v_{\rightarrow} (and \bar{p}_{t+1} for PROPER P-in-P).
- 11: end if
- 12: Apply the algorithm for type BOTH.
- 13: Deal with vertices created by subdivision of p_0p_1 and p_tp_{t+1} if necessary.
- 14: Modify the representation to construct \mathcal{R} .
- 15: **return** A representation \mathcal{R} extending \mathcal{R}' .

A.9 Proper Subpath-in-Path Graphs by Bin Packing

The pseudocode is in Algorithm 9. It is described in Section 4.1.4.

Algorithm 9 Sub Type Extension of Subpath-in-Path Graphs

Require: A proper subpath-in-path graph G and a partial representation \mathcal{R}' having k pre-drawn intervals.

- 1: Test all 2^k possible orderings of located components.
- 2: for each ordering do
- 3: Place the located components according to the ordering and minimum spans.
- 4: Distribute unlocated components into gaps using BinPacking.
- 5: end for
- 6: **if** no ordering works **then**
- 7: No representation exists, output "no".
- 8: end if
- 9: **return** A representation \mathcal{R} extending \mathcal{R}' .

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Common Notation

The following tables contain the most importation notation of this thesis, together with references to their precise definitions.

C.1 Symbols

Symbol	Description	Definition
\mathcal{R}	An intersection representation.	Chapter 1
R_u	The set representing u in \mathcal{R} .	Chapter 1
\mathcal{R}'	A partial representation.	Section 1.1
R'_u	The pre-drawn set representing u in \mathcal{R}' .	Section 1.1
N[u]	The closed neighborhood of u .	Section 1.5
◀	The left-to-right ordering of the components of an interval	Section 1.5
	representation.	
ℓ_u	The left endpoint of R_u .	Section 1.5
r_u	The right endpoint of R_u .	Section 1.5

Table C.1: The notation common for the entire thesis.

Symbol	Description	Definition
\triangleleft	The interval order of the maximal cliques.	Section 2.1.2
⋖	The left-to-right ordering of handles.	Section 2.1.2
cp(a)	A clique-point representing the clique a .	Section 2.2.1
(a), (a)	The leftmost and the rightmost position for $cp(a)$.	Section 2.2.2

Table C.2: The notation specific for Chapter 2.

Symbol	Description	Definition
\triangleleft	The left-to-right order of intervals.	Sections 3.1 and 4.1.1
<	The common left-to-right ordering of left and	Sections 3.1 and 4.1.1
	right endpoints.	

Table C.3: The notation specific for Chapter 3 and Chapter 4.

C.2 Classes and Problems

Abbreviation	Name of the class
INT	Interval graphs.
PROPER INT	Proper interval graphs.
UNIT INT	Unit interval graphs.
PROPER P-in-P	Proper subpath-in-path graphs.
P-in-P	Subpath-in-path graphs.
P-in-T	Subpath-in-tree graphs.
T-in-T	Chordal graphs, or equivalently subtree-in-tree graphs.

Table C.4: The classes of graphs used in this thesis, see Section 1.2.

Abbreviation	Name of the problem	Definition
Recog	The recognition problem.	Section 1.1
RepExt	The partial representation extension problem.	Section 1.1
$Recog^*$	The recognition problem with prescribed tree.	Section 1.2
BOUNDREP	The bounded representation problem.	Section 1.4.2
SIMREP	The simultaneous representations problem.	Section 1.3
3-Partition	The 3-partition problem.	Section 3.2.2
BinPacking	The bin packing problem.	Section 4.1.4

Table C.5: The computational problem used in this thesis.