



Master thesis

Acyclic *n*-Level Hypergraph Partitioning

Daniel Seemaier

Date: March 2, 2020

Supervisors: Prof. Dr. Peter Sanders Dr. Sebastian Schlag Priv. Doz. Dr. Christian Schulz

Institute of Theoretical Informatics, Algorithmics Department of Informatics Karlsruhe Institute of Technology

Abstract

Directed acyclic graphs are widely used to model the data flow and execution dependencies of streaming applications. Efficient parallelization of such graphs requires acyclic partitioning of the dependency graph. However, normal graphs are not always capable of precisely modelling such dependencies. Thus, we consider *directed acyclic hypergraphs* (DAHs). In this work, we present the first *n*-level hypergraph partitioning algorithm for directed acyclic hypergraphs. Moreover, we show (i) that our algorithm beats the current state of the art for directed acyclic graph partitioning in terms of solution quality on realistic instances, (ii) that *n*-level algorithms are superior to single level algorithms, and (iii) that our algorithm improves on the makespan of a parallelized image streaming application.

Zusammenfassung

Gerichtete azyklische Graphen werden häufig zur Modellierung von Datenflüssen und Ausführungsabhängigkeiten von Datenflussapplikationen genutzt. Eine effiziente automatische Parallelisierung solcher Anwendung erfordert eine azyklische Partitionierung der Abhängigkeitsgraphen. Allerdings ist eine präzise Modellierung der Abhängigkeiten mit herkömmlichen Graphen nicht immer möglich. Deswegen betrachten wir in dieser Arbeit gerichtete azyklische Hypergraphen (DAHs) und präsentieren den ersten n-Stufen Algorithmus zur azyklischen Partitionierung solcher Hypergraphen. Unsere Ergebnisse bestehen aus drei Beiträgen: wir zeigen, dass (i) unser Algorithmus häufig bessere Partitionen als der aktuell beste Algorithmus auf praxisnahen gerichteten azyklischen Graphen findet, (ii) unser n-Stufen Algorithmus auf gerichteten azyklischen Hypergraphen besser abschneidet als ein einstufiger Algorithmus und (iii) unser Algorithmus eine effizientere Parallelisierung für eine echte Datenflussapplikation ermöglicht.

Acknowledgments

I would like to thank my supervisors Dr. Sebastian Schlag and Priv. Doz. Christian Schulz, who have introduced me to research in Algorithmics and mentored me since my bachelor thesis. In many meetings over the last three years, they gave me more guidance than I could have ever asked for and supported me not only tremendously with this master thesis, but also made my time while pursuing my master's degree much more exciting. I also want thank Merten Popp, who found time during his busy workweeks to test our algorithm with a real world application, thereby strengthening the result of this research. Lastly, I thank the Steinbruch Centre for Computing for granting me access to the HPC cluster bwUniCluster. Without sufficient computational power, the experimental evaluation of this thesis would not have been possible.

Hiermit versichere ich, dass ich diese Arbeit selbständig verfasst und keine anderen, als die angegebenen Quellen und Hilfsmittel benutzt, die wörtlich oder inhaltlich übernommenen Stellen als solche kenntlich gemacht und die Satzung des Karlsruher Instituts für Technologie zur Sicherung guter wissenschaftlicher Praxis in der jeweils gültigen Fassung beachtet habe.

Karlsruhe, den 2. März 2020

Contents

Ał	Abstract iii Zusammenfassung iii					
Ζι						
1	Intro 1.1 1.2 1.3	oduction Problem Statement Contribution Structure of Thesis	1 2 2 3			
2	Fun 2.1 2.2	damentals General Definitions	5 5 8			
3	Rela 3.1 3.2	Multilevel Graph Partitioning	11 11 14 14 15			
	3.3 3.4	Acyclic Graph Partitioning	15 15 16 16 18 19			
4	Асу	clic Hypergraph Partitioning	21			
	4.14.24.3	<i>n</i> -Level Acyclic Hypergraph Partitioning 4.2.1 Initial Partitioning 4.2.2 Coarsening 4.2.3 Refinement 4.2.3 Refinement Larger Imbalances on Coarse Levels 4.3.1 Soft Rebalance 4.3.2 Hard Rebalance	 21 22 22 26 29 31 31 32 32 			
	4.4	Memetic Acyclic Hypergraph Partitioning	33			

5	Exp	erimental Evaluation	37
	5.1	DAG Model	38
	5.2	DAH Model	42
		5.2.1 Influence of Larger Imbalances on Coarse Levels	44
		5.2.2 Influence of Acyclic Coarsening	45
	5.3	Impact on Streaming Application	46
6	Con 6.1	clusion Future Work	49 49
Bil	oliogi	raphy	51
Α	Deta	ailed DAG Results	57
В	Deta	ailed DAH Results	61

1 Introduction

Directed acyclic hypergraphs (DAHs) are a generalized concept of directed acyclic graphs (DAGs) where each hyperedge can contain an arbitrary number of tails and heads. The acyclic hypergraph partitioning problem is to partition the hypernodes of a DAH into a fixed number of blocks of roughly equal size such that the corresponding quotient graph is acyclic while minimizing an objective function on the partition.

The problem is motivated by a recent paper on graph partitioning with acyclicity constraint by Moreira et al. [42] who use directed acyclic graphs to model the data flow and execution dependencies of image streaming applications. The imaging applications are executed on embedded processors with limited thermal budget and memory. To cope with these constraints, the application is distributed over multiple processors that process the data one after another. Data dependencies between parts of the application are modeled as a directed dependency graph. To distribute the application, the dependency graph is partitioned into the same number of blocks as there are processors available. Since edges between blocks correspond to interprocessor communication, the goal is to find a partition that minimizes the weighted edge cut of the partition. However, this is merely an imprecise approximation of the real objective function, which is to minimize the number of blocks containing neighbors of a node: say that one could choose between two partitions. One partition places all successors of a node into one other block, whereas the other partition splits them over two other blocks. When using a directed graph to model the data dependencies and the weighted edge cut objective function, both partitions are seen as equal, although the first option requires less interprocessor communication since the image has to be transferred to only one other processor, i.e., only once. This problem is illustrated in Figure 1.1: while the partitions shown in Figure 1.1a and 1.1b have different edge cuts, they behave the same in practice. Directed acyclic hypergraphs allow for a better model, since a single hyperedge can contain an arbitrary number of hypernodes. The *connectivity metric* then counts the number of blocks connected by a hyperedge. Using this model, both partitions are rated the same, as shown in Figure 1.1c and 1.1d.

This application leads to a lot of preexisting work on the directed acyclic graph partitioning problem, but to the best of our knowledge, none concern the directed acyclic hypergraph partitioning problem.

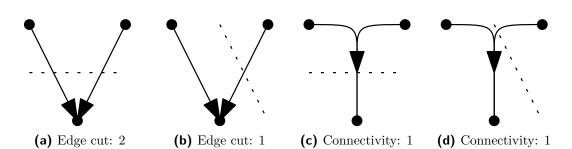


Figure 1.1: Model of an application divided into three subapplications (nodes) executed on two PEs (dashed line). Two partitions are proposed, one shown in Figures 1.1a and 1.1c, the other one shown in Figures 1.1b and 1.1d. Since in both cases, the result of one subapplication has to be sent to one other PE, both partitions perform equally good in practice and should therefore have the same cost. Figures 1.1a and 1.1b model the application using a DAG using the edge cut metric, which rates the partitions differently. In Figure 1.1c and 1.1d, the problem is modeled as a DAH using the connectivity metric, which puts the same cost to both partitions.

1.1 Problem Statement

To the best of our knowledge, there are currently no partitioning algorithms for directed acyclic hypergraphs in existing literature and therefore, it is not possible to use the DAH model in the described application domain. We close this gap by presenting the first algorithm for acyclic HGP in this thesis. Our algorithm can only handle hypergraphs where each hyperedge contains at most one head. This is sufficient to model the data flow and execution dependencies of an image streaming application, and also for a range of other applications [6,7].

1.2 Contribution

We have three main contributions: first and foremost, and to the best of our knowledge, we present the first algorithm for acyclic hypergraph partitioning by adapting techniques for DAG partitioning recently introduced by Herrmann et al. [25, 26] and Moreira et al. [42, 43]. We compare out algorithm on DAG instances with preexisting DAG partitioning algorithms to show an improvement of 10% on average. Second, we evaluate our algorithm on DAH instances. Since there are no preexisting algorithms that we could include in our benchmark set, we compare our algorithm with simpler heuristics and show that our *n*-level algorithm produces partitions with 64% lower connectivity on average than a

simple single-level k-way search with topological ordering for initial partitioning. Finally, we evaluate the impact of our improved DAH model on an image streaming application described in Ref. [42]. We show that our approach not only yields an improved model of the transfer costs, but also improves the makespan of such applications by up to 22% over the DAG model.

1.3 Structure of Thesis

The remaining content of this thesis is structured as follows. In Chapter 2, we introduce fundamental definitions from graph theory and give a precise formulation of the acyclic hypergraph partitioning problem. Chapter 3 follows with a broad overview on the current state-of-the-art graph and hypergraph partitioning techniques as well as recent approaches for solving the acyclic graph partitioning problem. This is the work that we build upon when developing our acyclic hyper-graph partitioner. The main content of this thesis is described in Chapter 4 and Chapter 5, where we present our approach to the acyclic hypergraph partitioning problem and perform an extensive experimental evaluation and comparison to previous work. Finally, we conclude our work in Chapter 6.

$1 \ Introduction$

2 Fundamentals

This chapter introduces general definitions that are used throughout this thesis. We start by defining undirected and directed graphs before generalizing them to undirected and directed hypergraphs. After defining those fundamental concepts, we give a precise formulation of the directed acyclic graph and hypergraph partitioning problems. Whenever possible, we use the notation introduced in Ref. [2,4,55].

2.1 General Definitions

An undirected weighted graph $G = (V, E, c, \omega)$ consists of finite sets V and E and weight functions c and ω . When talking about multiple graphs, we also use V(G)and E(G) to denote the node and edge set of a particular graph. The elements of V are called *nodes* and the elements of E are called *edges*. We define $n \coloneqq |V|$ and m := |E|. While the elements of V are arbitrary and of no further interest, E may only contain two-subsets of V. In other words, all edges $e \in E$ have the form $e = \{u, v\}$ with $u, v \in V$ and $u \neq v$. Figure 3.3a illustrates a simple undirected graph. We say that two nodes u and v are *adjacent* or *connected* if $\{u, v\} \in E$ and two edges e_1 and e_2 are *adjacent* if $e_1 \cap e_2 \neq \emptyset$. A node u and an edge e are *incident* if $u \in e$. The *neighborhood* $\Gamma(u)$ of a node u is the set of nodes adjacent to it. Its size $d(u) := |\Gamma(u)|$ is the degree of u. The maximum degree $\Delta(G) := \max_{u \in V} d(u)$ is the highest degree occurring in G. The node weight function $c: V \to \mathbb{R}_{>0}$ assigns non-negative weights to the nodes of G. Analogously, the edge weight function $\omega: E \to \mathbb{R}_{>0}$ assigns non-negative weights to the edges of G. The weight functions are extended to sets of nodes and edges by summing over the elements of the set, i.e., $c(V') = \sum_{v \in V'} c(v)$ for $V' \subseteq V$ and $\omega(E') = \sum_{e \in E'} \omega(e)$ for $E' \subseteq E$. If every node or every edge of a graph has the same weight, we say that the graph has unit node weights or unit edge weights and assume that c(V) = n or $\omega(E) = m$, respectively. A matching $M \subseteq E$ in a graph is a set of non-incident edges. Maximal matchings are a particular type of matchings with the property that there exists no edge $e \in E$ such that $M \cup \{e\}$ is also a matching.

The concept of *directed graphs* puts an order on the elements of each edge. More precisely, in a directed graph each edge is a pair (u, v) of some nodes $u, v \in V$ with u being the *tail* of the edge and v being its *head*. In this case, e is *directed* from u

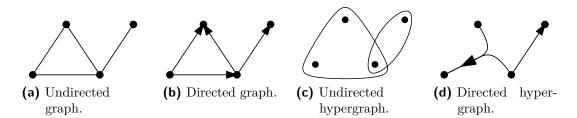


Figure 2.1: (a) An undirected graph where dots represent nodes and lines represent edges. (b) A directed graph. The Arrow of each edge points towards its head. (c) An undirected hypergraph. Hyperedges are drawn using closed polygons. (d) A directed acyclic hypergraph.

to v. In a directed graph, each node has predecessors $\Gamma^-(u) \coloneqq \{v \mid (v, u) \in E\}$ and successors $\Gamma^+(u) \coloneqq \{v \mid (u, v) \in E\}$. The degree $d(u) \coloneqq d^+(u) + d^-(u)$ of u is the sum of its out degree $d^+(u) \coloneqq |\Gamma^+(u)|$ and its in degree $d^-(u) \coloneqq |\Gamma^-(u)|$. A directed cycle $C = (v_1, \ldots, v_k, v_{k+1} = v_1)$ in a directed graph is a sequence of nodes such that $(v_i, v_{i+1}) \in E$ for $1 \leq i \leq k$ and $v_i \neq v_j$ for $1 \leq i, j \leq k, i \neq j$. A directed graph that does not contain any cycles is said to be *acyclic* and referred to as a directed acyclic graph or *DAG* for short. Figure 3.3b shows a simple directed acyclic graph. When talking about DAGs, we refer to nodes with in-degree 0 as sources and nodes with out-degree 0 as sinks of the graph. A subgraph $G' \subseteq G$ of a DAG G is a DAG that with $G'(V) \subseteq G(V)$ and $G'(E) \subseteq G'(V) \times G'(V)$. The induced subgraph $G[V'] \subseteq G$ for some $V' \subseteq G(V)$ is the subgraph of G with (G[V'])(V) = V' and $(G[V'])(E) = G(E) \cap (V' \times V')$.

A topological ordering $\tau: V \to [n]$ of the nodes of a directed graph is an order such that for every edge $(u, v) \in E$, $\tau(u) < \tau(v)$ holds. We use [n] to denote the set $\{1, \ldots, n\}$. As shown in Theorem 2.1.2, the existence of a topological ordering characterizes directed acyclic graphs.

Lemma 2.1.1. (Source: [8, Proposition 1.4.2].) Let G = (V, E) be a DAG. Then G contains a node v with in degree $d^{-}(v) = 0$.

Proof. Let $P = (v_1, \ldots, v_k)$ be a path of maximal length in G. Since G is acyclic, v_1 cannot have a predecessor in P and since P is maximal, it also cannot have a predecessor outside of P. Therefore, v_1 cannot have any predecessors, i.e., $d^-(v_1) = 0$.

Theorem 2.1.2. (Source: [8, Proposition 1.4.3].) A directed graph is acyclic if and only if there exists a topological order of the nodes of the graph.

Proof. Let G = (V, E) be a directed acyclic graph. By Lemma 2.1.1, $G_0 \coloneqq G$ has a node v_0 with in degree zero. Set $\tau(v_0) = 0$ and move on to $G_1 \coloneqq G_0 - v$. Since

removing a node from an acyclic graph keeps the graph acyclic, G_1 again contains a node v_1 with in degree 0 and we can set $\tau(v_1) = 1$ and so on. By induction, this constructs a topological order τ of G.

For the other direction, let τ be a topological ordering and assume that G contains a cycle $C = (v_1, \ldots, v_k, v_{k+1} = v_1)$. Then $\tau(v_1) < \tau(v_2) < \cdots < \tau(v_k) < \tau(v_1)$, a contradiction.

We now generalize the concept of graphs to hypergraphs by allowing edges to contain an arbitrary number of nodes. More formally, an undirected weighted hypergraph $H = (V, E, c, \omega)$ consists of a finite set of hypernodes V and a finite set of hyperedges E, where each hyperedge $e \in E$ is a non-empty subset of V, i.e., $e \subseteq V$. Hyperedges are also referred to as nets and the hypernodes contained in a net are its pins. Figure 2.1c illustrates a simple hypergraph. For a hypernode u, we define its set of incident hyperedges $I(u) \coloneqq \{e \in E \mid u \in e\}$, and its neighborhood $\Gamma(u) \coloneqq \{v \mid \{u, v\} \subseteq e \text{ for some } e \in E\}$. The size of a net is its cardinality |e|. A hypergraph where every edge has the same cardinality r is said to be r-uniform. In particular, a 2-uniform undirected hypergraph is an undirected graph. Analogously to undirected graphs, $c : V \to \mathbb{R}_{>0}$ assigns each hypernode a non-negative hypernode weight and $\omega : E \to \mathbb{R}_{>0}$ assigns each hyperedge a non-negative hyperedge weight.

The generalized version of directed graphs are *directed hypergraphs*. A directed hypergraph is an undirected hypergraph where each hyperedge $e \in E$ is divided into a set of tails $e^T \subseteq e$ and heads $e^H \subseteq e$ that fulfill $e^T \cup e^H = e$ and $e^T \cap e^H =$ \emptyset . Note that in this thesis, we only consider directed hypergraphs where each hyperedge contains at most one head pin and an arbitrary number of tail pins, i.e., hypergraphs with $|e^{H}| = 1$ for all $e \in E$. The predecessors of a hypernode u are $\Gamma^{-}(u) \coloneqq \{v \mid v \in e^{T}, u \in e^{H} \text{ for some } e \in E\}$ and its successors are $\Gamma^+(u) := \{v \mid u \in e^T, v \in e^H \text{ for some } e \in E\}.$ In a directed hypergraph, a cycle C of length k is a sequence of hypernodes, $C = (v_1, \ldots, v_k, v_{k+1} = v_1)$, such that for every i = 1, ..., k, there exists some hyperedge $e \in E$ with $v_i \in e^T$ and $v_{i+1} \in e^H$. Furthermore, we require that $v_i \neq v_j$ for $i \neq j, 1 \leq i, j \leq k$. Analogously to directed acyclic graphs, we refer to directed hypergraphs that do not contain any cycles as *directed acyclic hypergraph* or *DAH* for short. This definition of directed acyclic hypergraphs can be seen as an extension of Berge-acyclicity [10] to directed hypergraphs: Consider a bipartite graph G that contains one node for each hypernode and one node for each hyperedge of the hypergraph. For each $e \in E(H)$, add edges (u, e) for $u \in e^T$ and (e, v) for $v \in e^H$ to G. Then G is acyclic if and only if H is acyclic. An example for a directed acyclic hypergraph is shown in Figure 2.1d.

Note that given a directed hypergraph H, another way to construct a directed graph G that is equivalent to the hypergraph in regards to the acyclicity constraint is to replace each hyperedge e with a directed, bipartite graph from e^T to e^H . G is

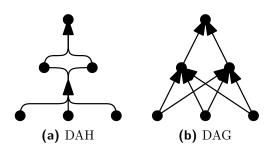


Figure 2.2: By replacing each directed hyperedge of the DAH (Figure 2.2a) with a bipartite graph from the tails of the hyperedge to its heads, we get a DAG (Figure 2.2b) that is acyclic if and only if the DAH is acyclic. Moreover, a partition of the DAH is acyclic if and only if the corresponding partition of the DAG is acyclic.

acyclic if and only if H is acyclic and an acyclic partition of G is also an acyclic partition of H and vice-versa. Figure 2.2 shows an example of this transformation.

2.2 Problem Formulation

As we did in the previous section, we start this section by defining the graph partitioning problems for undirected graphs. Then, we introduce it for directed acyclic graphs before stating both definitions for undirected and directed acyclic hypergraphs.

Given a graph G = (V, E), the k-way graph partitioning problem asks for a partition of V into k blocks $\Pi(V) \coloneqq \{V_1, \ldots, V_k\}$ such that $V = \bigcup_{1 \le i \le k} V_i$ and $V_i \cap V_j = \emptyset$ for $i \ne j$. The weight of each block is limited by the *imbalance parameter* $\epsilon \ge 0$ and might not be zero: we demand that $0 < c(V_i) \le L_{\max} \coloneqq (1+\epsilon) \lceil \frac{c(V)}{k} \rceil$ for all $1 \le i \le k$. This ensures that each block has roughly the same weight for small values of ϵ . The goal is to find a partition that minimizes an objective function. In this thesis, we only consider the edge cut objective defined by

$$\sum_{e \in \text{cut}} \omega(e),$$

where cut contains all edges with endpoints in two different blocks.

For directed acyclic graphs, we need the concept of the quotient graph. Given a directed acyclic graph G = (V, E) with partition $\Pi(V) = \{V_1, \ldots, V_k\}$, the quotient graph is a directed graph with one node v_i for each block V_i and an edge (v_i, v_j) if G contains an edge (v'_i, v'_j) with $v'_i \in V_i$ and $v'_j \in V_j$. The quotient graph can also be seen as the graph resulting from G when contracting all nodes within a

block. With this in mind, the graph partitioning problem for directed acyclic graphs is the same as before, but extended by the *acyclicity constraint*, i.e., the constraint that the partition's quotient graph must be acyclic.

Finally, we define undirected and directed acyclic hypergraphs. Given an undirected hypergraphs H = (V, E), a k-way partition $\Pi(V) = \{V_1, \ldots, V_k\}$ of H is a partition of V fulfilling the same conditions as before. For $u \in V$, we set b[u] to the block containing u, i.e., b[u] = i if and only if $u \in V_i$. Moreover, we define the *connectivity set* of a hyperedge e with $\Lambda(e) := \{V_i \mid V_i \cap e \neq \emptyset\}$ and the *connectivity* of a hyperedge e with $\lambda(e) := |\Lambda(e)|$. Two blocks $V_i, V_j \in \Pi$ are adjacent if there is a hyperedge $e \in E$ with $V_i \cap E \neq \emptyset$ and $V_j \cap E \neq \emptyset$. A hypernode u and a block V_i are adjacent if $u \notin V_i$ but there exists a hyperedge e with $u \in e$ and $e \cap V_i \neq \emptyset$. The most prominent objective functions in hypergraph partitioning are the *cut* and *connectivity* objectives. Given a partition $\Pi(V) = \{V_1, \ldots, V_k\}$, the former one is defined as

$$\sum_{e \in \mathrm{cut}} \omega(e),$$

where cut contains all hyperedges that are cut, i.e., have $\lambda(e) > 0$. The connectivity metric is defined as the sum

$$\sum_{e \in \text{cut}} (\lambda(e) - 1) \omega(e).$$

In other words, a hyperedge that only contains hypernodes from one block does not contribute to the connectivity metric whereas an edge containing hypernodes from two different blocks contributes one and so on. Note that both metrics are equal to the ordinary edge cut metric used in graph partitioning for 2-uniform hypergraphs. This allows us to compare our algorithm to the current state-of-the-art for DAG partitioning later on.

The quotient graph Q of a partitioned directed acyclic hypergraph H again contains a node v_i for each block V_i and an edge (v_i, v_j) if H contains an hyperedge e with tail pins in V_i and head pins in V_j . More formally, $V(Q) := \Pi$ and $E(Q) := \{(V_i, V_j) \mid \exists e \in E(H) : e^T \cap V_i \neq \emptyset \text{ and } e^H \cap V_j \neq \emptyset\}$. The hypergraph partitioning problem for directed acyclic hypergraphs is the same as before, but with the further restriction that the resulting quotient graph must also be acyclic.

Both graph partitioning problems are NP-complete [30, 42] and there are no constant factor approximation algorithms [5, 42]. Since the corresponding hypergraph partitioning problems are generalized versions of them, they are also NP-complete. In practice, we must therefore focus on good heuristics to find high-quality partitions of large graphs.

Fundamentals

3 Related Work

The general graph partitioning problem has been studied for a long time and subsequently, there is an enormous amount of preexisting literature. For a broad overview, we refer to existing literature [14,58]. More specialized literature focuses on hypergraph partitioning and acyclic graph partitioning. In this chapter, we give a brief introduction to each topic. We start by giving a general introduction to graph partitioning. Afterwards, we introduce hypergraph partitioning by briefly presenting the core components of KaHyPar [2, 4, 27, 54, 55] and PaToH [15]. Finally, we discuss recent approaches for acyclic graph partitioning introduced by Moreira et al. [42, 43] and Herrmann et al. [25, 26].

3.1 Multilevel Graph Partitioning

Most recent graph partitioning algorithms employ the multilevel paradigm first introduced by Hendrickson and Leland [24]. Note that this scheme isn't limited to ordinary node-based graph partitioning, but has been applied successfully to a wide range of problems, such as sequential and distributed edge partitioning [38,57], graph drawing [39] and even support-vector machines [56]. We also make heavy use of this scheme in this thesis. Therefore, we give a brief introduction to multilevel graph partitioning algorithms in this section. We start by outlining the multilevel partitioning scheme. Afterwards, we present the refinement algorithm by Fiduccia and Mattheyses (FM algorithm) [23] since we use this algorithm as a basis for our own refinement algorithm.

Outline. Multilevel graph partitioning typically consists of three phases, namely coarsening, initial partitioning and uncoarsening or refinement. The whole process is depicted in Figure 3.1. During coarsening, the algorithm constructs a hierarchy of roughly $\log(n)$ coarser graphs by contracting matchings or clusters. Both variants can be implemented efficiently. For instance, Birn et al. [12] present an efficient 2-approximative parallel matching algorithm with linear running time (on a single core) and low I/O complexity. Meyerhenke et al. [40] present a clustering algorithm using size-constrained label propagation which also has linear running time. Contracting a set $S \subseteq V$ of nodes works as follows: remove all nodes in S from

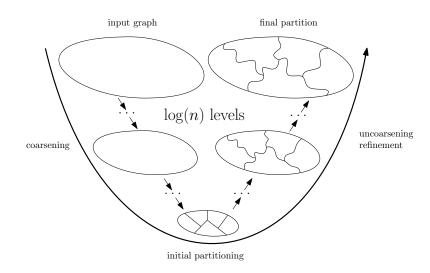


Figure 3.1: Multilevel partitioning scheme: coarsening, initial partitioning and uncoarsening with refinement.

the graph and insert a new node s with node weight c(s) = c(S) and neighborhood $\Gamma(s) = \bigcup_{u \in S} \Gamma(u)$. If this process creates parallel edges e_1, \ldots, e_m , they are merged into one edge e with accumulated edge weight $\omega(e) = \omega(\{e_1, \ldots, e_m\})$. After roughly $\log(n)$ levels of coarser graphs, the coarsening algorithm decides that the graph is small enough and terminates. We use the following terminology when referring to different graphs of the hierarchy: the *finest graph* is the input graph, i.e., the graph without any contractions. When contracting the finest graph, we get a series of *coarser graphs.* The smallest graph is referred to as the *coarsest graph.* The next phase computes a partition for the coarsest graph. This partition is referred to as *initial partition.* Since the coarsest graph is small compared to the input graph, the initial partitioning can be relatively slow without affecting the overall running time much. During uncoarsening, the partition of a coarser graph is projected onto the finer graph. This is done by assigning all nodes that were contracted into one node to that nodes partition in the coarser graph. After each projection, a *refinement* algorithm improves the the current partition. The most prominent refinement algorithms are variants of the algorithm by Fiduccia and Mattheyses (FM) [23] or the one by Kernighan and Lin (KL) [35]. Since we use the FM algorithm as basis of our own refinement algorithm, we elaborate on it in the following paragraph. Once refinement on the finest level was executed, the algorithm terminates.

The intuition of multilevel graph partitioning is based on the following observations: first, projecting a partition from a coarser level onto a finer level preserves the edge cut. Thus, the final edge cut is *at least* as low as the edge cut of the initial partition, assuming that the refinement algorithm guarantees no worsening. The other observation is that the multilevel scheme allows a more global view on the partition landscape compared to a single-level algorithm. During the coarser levels of the graph hierarchy, the refinement algorithm considers to move chunks of nodes (since a single node on the coarsest graph corresponds to many nodes on the finest graph), allowing a global view. During finer levels, refinement performs fine-grained moves.

Fiduccia–Mattheyses Algorithm. The FM algorithm [23] is a linear-time local search heuristic originally proposed as a 2-way refinement algorithm. The algorithm alternates between both blocks and moves the node with the highest gain (i.e., the change in the objective caused by the move) from the current block to the other one. To do that, the algorithm maintains two priority queues, one for each block. The queues are initialized with all nodes belonging to the corresponding block, with priority equal to their gain value. This allows an implementation with amortized linear running time when using bucket queues as priority queues; for details, see Ref. [23]. Once a node was moved, it is excluded from the rest of the current pass of the algorithm. The pass terminates as soon as one priority queue runs empty. Then, the algorithm undos moves in reverse order until the best partition observed during refinement is restored.

While Fiduccia and Mattheyses move all nodes during a pass of one iteration, Karypis and Kumar [34] reduce its running time significantly by only considering boundary nodes for movement and stopping the search after a constant number of fruitless moves, i.e., moves that did not yield an improved partition. Moreover, they extended the original 2-way FM algorithm to a k-way refinement algorithm using only a single global priority queue while maintaining the linear running time. Osipov and Sanders [47] present an adaptive stopping criterion (instead of a constant number of fruitless moves) based on a random walk model. The criterion terminates the current pass as soon as further improvements become unlikely. Moreover, Osipov et al. [52] introduce a highly localized version of the FM algorithm called multi-try FM. Instead of initializing a pass of the algorithm with all border nodes, they choose to repeatedly initialize it with only a single node. The intuition behind this improvement is that it allows the algorithm to find localized improvements before the partition gets trashed at other places.

For hypergraph partitioning, Akhremtsev et al. [2] implement a variant of the FM algorithm that maintains k priority queues, one for each block, and uses an extension of the adaptive stopping criterion introduced by Osipov and Sanders [47]. Each priority queue contains all hypernodes that can be moved to the corresponding block, i.e., that are in another block but adjacent to it, with the gain value as priority. Similar to multi-try FM [52], their FM implementation is highly localized and gets repeatedly initialized with only a couple of hypernodes. Since calculating and updating gain values on hypergraphs is expensive, they implement several novel techniques to improve the performance of their FM algorithm. First and foremost,

they cache gain values across multiple passes to prevent having to calculate the same gain value multiple times. Secondly, they exclude large hyperedges since those bottleneck the gain value calculation and are unlikely to change their connectivity.

3.2 Undirected Hypergraph Partitioning

In this section, we briefly introduce hypergraph partitioning by presenting two algorithms for undirected hypergraph partitioning, namely KaHyPar [2,4,27,28,55] and PaToH [15]. We present KaHyPar because it computes the best partitions out of all publicly available HGP algorithms on a large collection of hypergraph instances [54] and we use it as a framework to implement our own algorithms. PaToH is generally one of the fastest HGP algorithm [54] and we use it during initial partitioning in our of our experiments. For an in-depth introduction to high-quality hypergraph partitioning and KaHyPar, we refer to the doctoral thesis of Sebastian Schlag [54].

3.2.1 KaHyPar

Many recent hypergraph partitioners [15, 21, 32, 59] and graph partitioners [29, 33, 50, 52] use the multilevel paradigm first introduced by Hendrickson and Leland [24] with roughly $\log(n)$ levels to partition an undirected hypergraph or undirected graph. KaHyPar uses an extreme version of this paradigm known as the *n*-level multilevel graph partitioning paradigm that was first introduced by Osipov and Sanders with KaSPar [47].

Like multilevel algorithms for graph partitioning, KaHyPar constructs a hierarchy of coarser graphs by contracting pairs of hypernodes during coarsening. Contracting a pair (u, v) of hypernodes works as follows: remove both hypernodes from the hypergraph and replace them by a new hypernode w with hypernode weight c(w) = c(u) + c(v). Add w to every hyperedge that contained u or v. If this process yields parallel hyperedges e_1, \ldots, e_k , the parallel hyperedges are replaced by a single hyperedge e with hyperedge weight $\omega(e) = \omega(\{e_1, \ldots, e_k\})$. In contrast to the normal multilevel paradigm, KaHyPar stores the resulting hypergraph after every single pair contraction in a hierarchy of coarser hypergraphs, giving the paradigm its name. At some point, the coarsening algorithm decides that the hypergraph is small enough and terminates. KaHyPar then computes an initial partition for the coarsest hypergraph. During uncoarsening, the hierarchy is unrolled and the hypernodes get uncontracted in reverse order. After each uncontraction operation, a localized refinement algorithm tries to improve the connectivity of the partition around the uncontracted hypernodes. Once all hypernodes were uncontracted, the partitioning process is complete.

KaHyPar implements various algorithms for each phase. For coarsening, KaHy-Par rates pairs of adjacent hypernodes according to the *heavy-edge* rating function [55] and contracts the pair with the highest rating next. For initial partitioning, it implements several algorithms including random hypernode assignment, breathfirst search and greedy hypergraph growing to obtain an initial k-way partitioning using recursive bisection. Finally, KaHyPar implements several refinement algorithms, namely localized 2-way FM refinement [55], localized k-way FM refinement [2] and 2-way flow based refinement algorithms [27]. Further techniques outside the n-level paradigm include the detection of community structures to guide the coarsening algorithm [28] and a memetic algorithm [4]. For more details on the various component as well as a general introduction to hypergraph partitioning, we refer to Ref [54].

3.2.2 PaToH

PaToH [15] is a multilevel hypergraph partitioning algorithm with roughly $\log(n)$ levels by Çatalyürek and Aykanat that originated from sparse-matrix vector multiplication. It optimizes the connectivity objective and uses recursive bisection to obtain a k-way partition. During coarsening, PaToH constructs a hierarchy of coarser hypergraphs using either a matching based hierarchical clustering algorithm or a hierarchic-agglomerative clustering algorithm. Then, it obtains an initial partition on the coarsest hypergraph using greedy hypergraph growing. Refinement is done using a variant of the 2-way FM algorithm that only keeps border hypernodes in its priority queues. After every mode, the algorithm updates the gain values of its adjacent hypernodes, removes hypernodes that are no longer boundary hypernodes from its priority queue and adds those that became border nodes due to the move.

3.3 Acyclic Graph Partitioning

Research on acyclic graph partitioning dates back multiple decades, although early work only considers simple heuristics to partition the graph. Two approaches make use of a maximum-fanout-free cone clustering [16] of the graph, a technique commonly used to make circuits sparser. Kocan et al. [37] cluster the graph and greedily join the relatively small clusters into blocks to obtain a final partition. Cong et al. [19] obtain an initial partition based on a topological ordering of the graph, then cluster it and use a variant of the FM algorithm with cycle detection to improve the partition on the clustered graph. In a subsequent work, Cong et al. [17] improve on their initial research by presenting FLARE, a 2-level partitioning algorithm using edge separability-based circuit clustering [18] and scheduled 2-way FM refinement. These approaches are superseded in terms of solution quality by recent multilevel algorithms from Herrmann et al. [25, 26] and Moreira et al. [42, 43]. The former presents new coarsening algorithms to produce acyclic coarser graphs and shows how to use a pre-existing undirected graph partitioner to obtain an initial partition of a DAG. The latter introduces more sophisticated k-way refinement algorithms and an evolutionary algorithm. To the best of our knowledge, the latest work of Herrmann et al. [25] is the current state-of-the-art in terms of solution quality. Since we make use of these techniques, we present them in detail in the following sections.

3.3.1 Initial Partitioning

In contrast to the classic multilevel partitioning scheme, both Herrmann et al. and Moreira et al. compute the initial partition of the DAG on the finest level of the graph hierarchy, i.e., before computing the coarser graphs. Moreira et al. do this because their coarsening algorithm might create cycles in the graph, making it potentially impossible to obtain a balanced initial partition otherwise. Herrmann et al. tried both variants but report better results for computing the initial partition on the input graph.

Moreira et al. [42] use a simple greedy heuristic to obtain an initial k-way partition of the input DAG that fulfills the acyclic constraint. To be more precise, they compute a topological order τ of the DAG and assign nodes $\tau^{-1}(i \cdot \lceil \frac{n}{k} \rceil), \ldots, \tau^{-1}((i + 1) \cdot \lceil \frac{n}{k} \rceil - 1)$ to block $i, i = 0, \ldots, k - 1$. While the resulting partition is acyclic and balanced, the heuristic ignores its edge-cut.

Herrmann et al. [26] propose a more sophisticated heuristic for computing initial bipartitions that makes use of pre-existing undirected graph partitioners such as METIS [33]. The algorithm is outlined in Figure 3.2: first, they treat the DAG as undirected graph and use the pre-existing undirected graph partitioner to obtain an initial bipartition. The resulting bipartition is balanced and optimizes the correct objective, but does generally violate the acyclic constraint. Therefore, additional work is required to restore the constraint.

3.3.2 Coarsening

Moreira et al. [43] use sized-constrained label propagation [40] to identify clusters and contract them. In regards to the acyclic property of the directed graph, they contract arbitrary sets of nodes which can cause the coarser graph to become cyclic, although this is unproblematic when calculating the initial partition before coarsening.

Herrmann et al. [26] present a novel coarsening algorithm that computes an acyclic coarser graph. On a high level, the algorithm computes a clustering of the graph while avoiding *forbidden edges*, i.e., edges that might

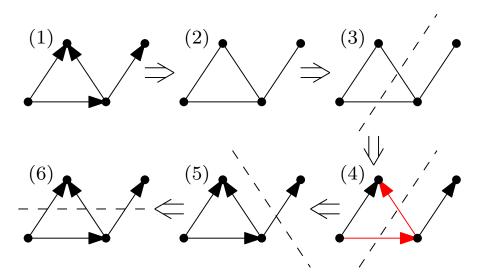


Figure 3.2: Obtaining an initial partition using a pre-existing initial partitioner: Treat the (1) directed input graph as (2) undirected graph and (3) use a pre-existing graph partitioner to partition the undirected graph. The (4) resulting partition might violate the acyclic constraint, which (5) can be restored by moving nodes. Those movements might imbalance the partition, requiring (6) a final re-balance step.

induce a cycle when contracted. Afterwards, it simultaneously contracts all clusters to obtain the coarser graph.

To identify forbidden edges, they introduce the concept of *toplevel* values as outlined in Definition 3.3.1.

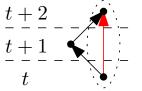
Definition 3.3.1. (Toplevel of a node. Source: [26].) Let G = (V, E) be a DAG. The toplevel of a node $v \in V$, denoted by top[v], is the length of a longest path from any source in G to v. In particular, sources s of the graph have toplevel top[s] = 0.

Observe that the contraction of an edge e = (u, v) induces a cycle in the coarser graph if and only if the graph contains an *u*-*v*-path avoiding *e*. Theorem 3.3.2 states conditions that are sufficient to identify such edges.

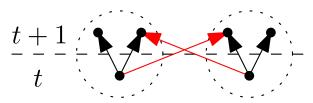
Theorem 3.3.2. (Source: [26].) Let G = (V, E) be a DAG and $C = \{C_1, \ldots, C_k\}$ be a clustering of V. If C is such that

- for any clustering C_i and for all $u, v \in C_i$, $|top[u] top[v]| \leq 1$, and
- for two different clusters C_i and C_j and for all $u \in C_i$ and $v \in C_j$, either $(u, v) \notin E$, or $top[u] \neq top[v] 1$,

then contracting all clusters in C yields an acyclic coarser graph.



 (a) Contracting the cluster induces a cycle of length 2. The cluster is forbidden by the first condition since the head of the red edge has toplevel > t + 1.



(b) Contracting both clusters induces a cycle of length 2. The clustering is forbidden by the second condition since the red edges connect nodes with toplevel difference ≤ 1 from different clusters.

Figure 3.3: Cyclic formations prevented by the (a) first and (b) second condition of Theorem 3.3.2.

The first condition prevents the formation of cycles that only contain one of the contracted clusters, while the second condition prevents the formation of cycles containing multiple contracted clusters. An example for each case is given in Figure 3.3.

Initially, all nodes are in their own cluster. The algorithm iterates over all nodes still in a singleton cluster and tries to add it to one of the neighboring clusters. A node can only be added to a cluster if it does not violate the criteria formulated in Theorem 3.3.2. The algorithm then selects the best neighboring cluster based on some rating function such as the edge weight and adds the node to the cluster. If no neighboring clusters are viable, the node stays in its singleton cluster.

3.3.3 Refinement

Moreira et al. [42] describe simple k-way refinement algorithms that move nodes with the highest gain values, where the gain value of a move is the change of the objective function on the partition caused by it. Since arbitrary node movements can cause the partition to become cyclic, they propose several restrictions on the set of potential movements and suggest the use of an online cycle detection algorithm. For the following paragraphs, let $\Pi = \{V_1, \ldots, V_k\}$ be the acyclic k-way partition with quotient graph Q and let $\tau : Q(V) \to [k]$ be a topological order of the quotient graph. They propose the following restrictions on possible target blocks of a node.

The simple move heuristic only considers to move a node in block $\tau^{-1}(i)$ to blocks $\tau^{-1}(i-1)$ (if the node does not have predecessors in its current block) and $\tau^{-1}(i+1)$ (if the node does not have successors in its current block). Observe that a simple move does not change the topological order of the quotient graph. The *advanced move* heuristics determines the highest block $\tau^{-1}(a)$ (in the topological order) containing a predecessor of a node and the lowest block $\tau^{-1}(b)$ containing a successor. Then, the viable target blocks are $\tau^{-1}(a), \tau^{-1}(a+1), \ldots, \tau^{-1}(b-1), \tau^{-1}(b)$. An advanced move might add a new edge to the quotient graph, but the topological order is preserved. *Global moves* consider all target blocks. Whenever a global move would introduce a new edge in the quotient graph, they use Kahn's algorithm [31] to determine whether the move creates a cycle in the quotient graph. Although Kahn's Algorithm has linear running time, this approach is viable since the size of the quotient graph is usually small.

Both Moreira et al. [42] and Herrmann et al. [26] suggest the use of the 2-way FM algorithm. Moreira et al. schedule the FM algorithm on pairs on blocks to refine a k-way partition, whereas Herrmann et al. use recursive bisection and only consider bipartitions. In 2-way refinement, there is no need for cycle detection since all moves are simple moves. Therefore, it is easy to identify movable nodes.

3.4 Memetic Algorithms

Memetic algorithms are inspired by the theory of evolution and were first introduced by Pablo Moscato [44]. For a broad introduction to memetic algorithms, we refer to the work by Moscato and Cotta [45]. Here, we only summarize the metaheuristic briefly. The basic building blocks of memetic algorithms are *mutation*, *recombination*, *local search*, a *fitness function*, and a population of individuals (i.e., solutions to the optimization problem). Initially, the algorithm produces individuals that form the initial population. Then, mutation and recombination operations are repeated until the population converges or the time limit is exceeded. Both operations produce offsprings that form the next generation of individuals. If the population is full, old individuals have to be evicted to make room for the new ones. To chose which individuals to evict, the fitness function is used. In general, individuals with low fitness are most likely to evicted from the population. The operations to produce offsprings work as follows. Mutation selects on individual and changes it, for example by applying local search to it. Recombination selects two individuals and combines them, forming the offspring.

KaHyPar also includes a memetic algorithm [4], which uses natural combine operations provided by the multilevel partitioning scheme. Since we use this implementation as starting point for our own memetic algorithm, we briefly describe its core components. The algorithm first produces a population of high-quality hypergraph partitions obtained using KaHyPar-C [28]. For mutation, it selects a random individual and iterates coarsening and local search using random seeds. For recombination, it selects two individuals using tournament selection [41] and combines them using the following operation: Coarsen the graph, but only contract pairs of hypernodes that are in the same blocks in both individuals. This ensures that either individual can be used as initial partition on the coarsest hypergraph. They use the better one and proceed with unrolling the hypergraph hierarchy while improving the partition using local search. When inserting a new offspring into the population, it evicts the individual most similar to the new offspring to ensure a diverse population. Furthermore, KaHyPar implements more sophisticated mutation and recombination operations to produce a more diverse population.

Besides undirected hypergraph partitioning, memetic algorithms have been applied successfully to a broad field of problems, including graph partitioning [53] and clustering [11], node separators [53], and the territory design problem [1]. Recently, Moreira et al. [43] proposed a memetic algorithm for the DAG partitioning problem. For more applications and trends, we refer to recent surveys [36,46] on the subject.

4 Acyclic Hypergraph Partitioning

Our hypergraph partitioner with acyclicity constraint is based on KaHyPar [2,4,55], a state-of-the-art hypergraph partitioner. To cope with the acyclicity constraint, we extend KaHyPar with directed hypergraphs and implement new algorithms for coarsening, initial partitioning, and the refinement of directed hypergraphs and acyclic partitions. We start this chapter with an outline that summarizes the interaction between each component of our DAH partitioner. The following subchapters then explore each component in detail.

Just like Moreira et al. [42, 43] and Herrmann et al. [25, 26], we start Outline. the partitioning process by computing an initial partition of the hypergraph on the input hypergraph. In this phase, we consider two alternatives: partitioning via topological order and partitioning via a pre-existing undirected hypergraph partitioner. Once the initial partition has been obtained, we run a 2-way refinement algorithm on the initial partition, followed by a V-cycle structured as follows. For coarsening, we use our acyclic coarsening algorithm described in Chapter 4.2.2. Since the hypergraph is already partitioned, the coarsening algorithm does not contract pairs of hypernodes in different blocks. This ensures that the initial partition can be projected onto the coarsest hypergraph of the *n*-level hypergraph hierarchy. The contracted hypernodes are then uncontracted. After each uncontraction, we run our 2-way refinement algorithm initialized with the uncontracted hypernodes. Once we obtained a k-way partition, we run a V-cycle with our k-way refinement algorithm presented in Chapter 4.2.3. The whole algorithm is also outlined in Figure 4.1. To further improve the result of our multilevel algorithm, we introduce an evolutionary algorithm in Chapter 4.4.

4.1 Data Structure for Directed Hypergraphs

In memory, KaHyPar represents hypergraphs as undirected bipartite graphs stored in an adjacency array [55]. The bipartite graph contains one node for each hypernode and one for each hyperedge. The nodes representing hypernodes have outgoing edges to the nodes representing incident hyperedges and analogously, nodes representing hyperedges are connected to nodes representing their pins. We extend this format by introducing a *head counter* h for each hypernode and each hyperedge of the bipartite graph. For hyperedges, h stores the number of head pins and for hypernodes u, it stores the number of hyperedges that contain u as a head pin. During hypergraph construction, we ensure that head pins are placed first in the adjacency array of a hyperedge and likewise, we ensure that hyperedges containing a head pin u are placed first in the adjacency array of u. This allows us to effectively iterate over only the heads or tails of a hyperedge.

4.2 *n*-Level Acyclic Hypergraph Partitioning

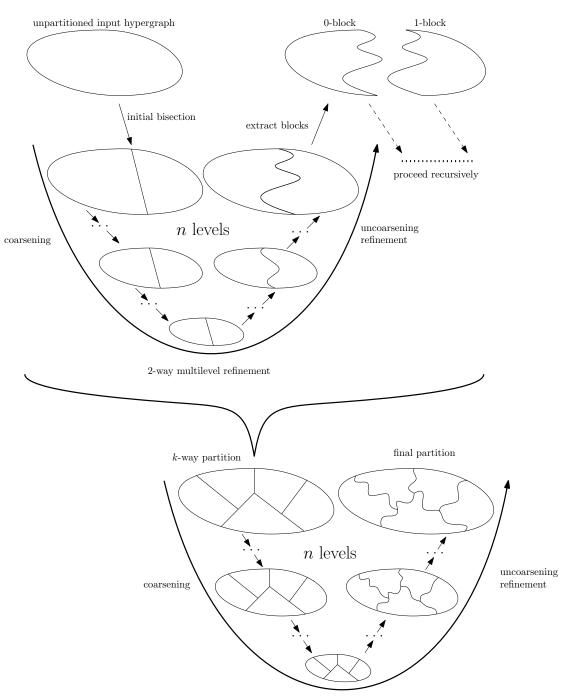
Recall that KaHyPar uses the n-level hypergraph partitioning scheme [47] described in Chapter 3.2.1 to obtain hypergraph partitions of high quality. We use this implementation and exchange the algorithms used for initial partitioning, coarsening and refinement with our own. These algorithms are presented in the following sections.

4.2.1 Initial Partitioning

This section describes our approaches for obtaining an initial partition of the directed acyclic hypergraph. Each algorithm starts with an unpartitioned directed acyclic hypergraph H = (V, E) and produces a partition of V into blocks V_1, \ldots, V_k for a fixed number of blocks k. Recall that depending on the configuration, H could refer to the input hypergraph or the coarsened hypergraph, i.e., the output of the coarsening algorithm described in Chapter 4.2.2.

Initial Partitioning via Topological Ordering. Recall that Moreira et al. [42] compute their initial partition based on a topological ordering of the graph. We implement the same approach for directed acyclic hypergraphs to obtain an initial k-way partition. First, we calculate a topological ordering of the nodes of the hypergraph using Kahn's algorithm [31] adapted for directed hypergraphs (Algorithm 1). Using the topological order, our algorithm greedily assigns nodes to blocks until they are full. More precisely, it assigns the first nodes to the first block until its weight exceeds $\lceil \frac{c(V)}{k} \rceil$ before it assigns the next nodes to block two and so on. Note that this approach always produces a balanced initial partition for hypergraphs with unit node weight. For weighted hypergraphs, it might produce an initial partition violating the balance constraint due to the greedy assignment of nodes to blocks. In this case, the refinement step must balance the partition.

Initial Partitioning via Undirected Partitioning. This algorithm is based on the initial partitioning algorithm for DAG partitioning presented by Herrmann et al. [26]



k-way refinement

Figure 4.1: High-level overview over our algorithm with all components. First, we obtain an initial partition using recursive bisection. For each bisection step, we use one of the initial partitioning algorithms. Then, we improve the partition using our 2-way refinement algorithm. Once the hypergraph is partitioned into k blocks, we run one V-cycle with our k-way refinement algorithm. Algorithm 1: Kahn's algorithm [31] for DAHs.

Data: DAH H = (V, E). Result: Topological ordering τ of H. 1 $i \leftarrow 0$ 2 while exists $v \in V$ with $d^+(v) = 0$ do 3 $\begin{bmatrix} \tau(i) = v \\ H \leftarrow H - v \\ 5 \end{bmatrix} \begin{bmatrix} \tau(i) = v \\ H \leftarrow i + 1 \end{bmatrix}$ 6 return τ

described in Chapter 3.3.1. Given a DAH, we compute an initial bipartition of the hypergraph as follows. First, we turn the DAH into an undirected hypergraph by merging the tails and heads of each hyperedge. Note that due to our data structure, this is merely a conceptual step and does not require any actual work. Next, we pass the undirected hypergraph to a preexisting hypergraph partitioner that minimizes the connectivity metric to obtain an initial bipartition. We use KaHyPar-MF [27] and PaToH [15]. To the best of our knowledge, KaHyPar-MF regularly finds partition with the lowest connectivity metric out of all hypergraph partitions while PaToH is the fastest one. This bipartition is the projected onto the original DAH. In general, the resulting bipartition violates the acyclicity constraint and therefore, we must perform further steps to make it acyclic.

The algorithm to make the bipartition acyclic is shown in Algorithm 2. On a high level, we select one edge in the quotient graph that we want to remove and move hypernodes from one block to the other one accordingly. Denote the two blocks by V_1 and V_2 and assume that we want to remove the quotient graph edge from V_1 to V_2 . We start a breadth-first search at every hypernode in V_1 that has successors in V_2 . The search only scans successors in V_2 and moves every node from V_2 to V_1 . Once the search has completed, no hypernode in V_1 has successors in V_2 and therefore, the quotient graph edge from V_1 to V_2 is removed.

The resulting acyclic partition might become imbalanced due to the movements from one block to the other one. To cope with this problem, we run an additional balancing step afterwards. This step simply moves hypernodes from the overloaded block to the underloaded block. Note that cannot move arbitrary hypernodes while keeping the bipartition acyclic. More precisely, if we have an acyclic bipartition with blocks V_1 and V_2 and a quotient graph edge from V_1 to V_2 , we can only moves hypernodes in V_1 that have no successors in V_1 . In an effort the keep the connectivity of the bipartition low, we sort the movable hypernodes in the overloaded block by their gain value using a priority queue. The whole process in depicted in Algorithm 4. Algorithm 3 puts it all together: first, it obtains a bipartition violating the acyclic constraint. Then, it runs Algorithm 2 and Algorithm 4 twice, once removing the quotient graph edge from V_1 to V_2 and once removing the reverse edge. Finally, it selects the bipartition with the lower connectivity and returns it. We run this algorithm twice; once as depicted, and then a second time using predecessors instead of successors in Algorithm 2. We then select the best partition out of all options.

We also considered to skip the balancing step by using an asymmetric number of further bisections to split the resulting two blocks: after bisecting a block B(that should ultimatly be split into k blocks) of weight c(B) into smaller blocks B_1 and B_2 , we set $k_1 := \lfloor k \cdot \frac{c(B_1)}{c(B)} \rfloor - 1$ and $k_2 := \lfloor k \cdot \frac{c(B_2)}{c(B)} \rfloor - 1$ and continue by splitting B_i into k_i blocks, i = 1, 2. Once the algorithm obtained all k blocks, we use the hard rebalancing algorithm described in Chapter 4.3.2 to ensure a balanced k-way partition. However, since initial experiments did not show an improvement over Algorithm 3, we did not pursue this approach any further.

Algorithm 2: Subroutine $FixCyclic(\cdot)$ referenced in Algorithm 3: moves nodes to make a bipartition acyclic.

Data: Cyclic bipartition (V_1, V_2) of DAH H = (V, E). Result: Acyclic bipartition. 1 S := new Stack()2 for $u \in V_1$ with $\Gamma^+(u) \cap V_2 \neq \emptyset$ do $S := S \cup \{u\}$ 3 while $S \neq \emptyset$ do 4 u := S.pop()5 for $v \in \Gamma^+(u) \cap V_2$ do 6 $S := S \cup \{v\}$ 7 $V_1 := V_1 \cup \{v\}$ 8 $u := V_2 \setminus \{v\}$ 9 return (V_1, V_2)

Algorithm 3: Initial partitioning algorithms that makes use of a preexisting hypergraph partitioner $HG(\cdot, \cdot)$ for undirected hypergraphs.

Data: DAH H = (V, E). Result: Bipartition $V = V_1 \dot{\cup} V_2$. 1 $(V_1, V_2) \coloneqq HG(H, k)$ // partition as undirected hypergraph 2 $(V'_1, V'_2) \coloneqq Balance(FixCyclic(V_1, V_2))$ // break (V_1, V_2) 3 $(V''_1, V''_2) \coloneqq Balance(FixCyclic(V_2, V_1))$ // break (V_2, V_1) 4 return min $\{(V'_1, V'_2), (V''_1, V''_2)\}$ // select bipartition with lower connectivity Algorithm 4: Subroutine $Balance(\cdot)$ referenced in Algorithm 3: moves nodes from the overloaded block to the underloaded one to balance the bipartition.

Data: Imbalanced acyclic bipartition (V_1, V_2) of DAH $H = (V, E)$. Let V_1			
be the overloaded and V_2 be the underloaded block and assume that			
the quotient graph edge goes from V_1 to V_2 .			
Result: Acyclic balanced bipartition.			
1 $Q \coloneqq \texttt{new PriorityQueue}()$			
$_2$ for $u\in V_1$ do			
3 $ $ if $\Gamma^+(u) \cap V_1 = \emptyset$ then			
$4 \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \ \$			
5 while $c(V_1) > (1+\epsilon) \lceil \frac{c(V)}{2} \rceil$ do			
$_{6} \mid u \coloneqq Q.\texttt{deleteMax}()$			
7 $V_1 \coloneqq V_1 \setminus \{u\}$			
9 UpdateGainValues (Q) // update gain values of neighbors of u in Q			
10 UpdateMovableHypernodes (Q) // remove/add hypernodes from/to Q			
11 return (V_1,V_2)			

4.2.2 Coarsening

The coarsening phase iteratively selects a set of hypernodes and contracts them, yielding a hierarchy of *n* levels of coarser hypergraphs. When contracting pairs of hypernodes, the hypergraph can become cyclic as illustrated in Figure 4.2. We explore two types of coarsening algorithms: the first one is the coarsening algorithm already implemented in KaHyPar-K. It selects pairs of hypernodes to be contracted without constraints in regards to keeping the directed hypergraph acyclic. The second approach restricts the algorithm implemented in KaHyPar-K to pairs of hypernodes that can be contracted safely while keeping the hypergraph acyclic. This approach is based on the acyclic clustering by Herrmann et al. [26] presented in Chapter 3.3.2 and described in the rest of this chapter.

When contracting a pair of hypernodes that are both pins of the same hyperedge, but one is one of the hyperedge's tails and the other one is one of the hyperedge's heads, it is not obvious whether the resulting hypernode should be a tail or a head of the hyperedge. Recall that in this thesis, we focus on directed hypergraphs where each hyperedge contains *at most* one head. For those instances, whenever one of the contraction partners is a head of a hyperedge e, the resulting hypernode is also a head of e. For the more general case, we propose two differ-

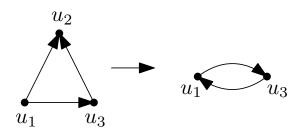


Figure 4.2: Acyclic directed hypergraph with three hypernodes u_1, u_2, u_3 on the left. Contracting the hyperedge $\{u_1, u_2\}$ yields the cyclic hypergraph on the right.

ent solutions. First, one could exclude such pairs altogether, i.e., only contract pairs of hypernodes that have the same role in all shared hyperedges. We ran some initial experiments with this setting and observed that it seemed to decrease partition quality. Hence, we did not pursue it further. As an alternative solution (that we did not implement), we propose to introduce *mixed pins* and leave it to the refinement algorithms to treat them appropriately.

Acyclic Hypergraph Coarsening. The algorithm is based on Theorem 4.2.2, which identifies pairs of hypernodes that may not be in the same cluster.

Definition 4.2.1. (Mixed-level and single-level clusters.) Let H = (V, E) be a DAH and $C = \{C_1, \ldots, C_k\}$ be a clustering of V such that for each $C_i \in C$ and for all $u, v \in C_i$, $|top[u] - v[v]| \leq 1$. We refer to clusters C_i where all $u, v \in C_i$ have top[u] = top[v] as single-level clusters and clusters C_j that contain at least one pair of nodes $u, v \in C_j$ with |top[u] - top[v]| = 1 as mixed-level clusters.

Theorem 4.2.2. (Based on Theorem 3.3.2.) Let H = (V, E) be a DAH and $C = \{C_1, \ldots, C_k\}$ be a clustering of V, such that

- for any clustering C_i and for all $u, v \in C_i$, $|top[u] top[v]| \leq 1$, and
- for two different **mixed-level clusters** C_i and C_j and for all $u \in C_i$ and $v \in C_j$, either $(u, v) \notin E$ or |top[u] top[v]| > 1.

Then the coarser hypergraph $H \circ C$ is acyclic.

Proof. First, recall that a DAH can be seen as a DAG by replacing each hyperedge e with a directed bipartite graph from e^T to e^H . This transformed graph is acyclic if and only if the DAH is acyclic. Hence, we can proof the theorem based on graphs rather than hypergraphs.

Let G be the corresponding DAG. Assume that $G \circ C$ contains a cycle. By Theorem 3.3.2, the cycle must contain at least one single-level cluster C_i . Moreover, since the nodes of C_i have the same toplevel, the cycle must have length at least 2. Let C_{i-1} be the predecessor and C_{i+1} be the successor of C_i in the cycle. Let t be the lowest toplevel of nodes in C_{i-1} . Then the toplevel of nodes in C_i and C_{i+1} is at least t + 1, which forbids a path from C_{i+1} to C_{i-1} , a contradiction.

Note that the difference between Theorem 3.3.2 and Theorem 4.2.2 lies in the distinction between single-level and mixed-level clusters: the second condition must only hold for pairs of mixed-level clusters. Since adjacent nodes in DAGs always have different toplevels, the clustering algorithm by Herrmann et al. [25] only produces mixed-level clusters. In contrast, DAHs might contain adjacent hypernodes with the same toplevel, justifying this distinction.

Based on this theorem, our clustering algorithm works as follows. We start on the input hypergraph and compute all toplevels, then start the clustering process. Generally, we only consider putting hypernodes in the same cluster whose toplevel differs by at most one, i.e., we only build clusters that do not violate the first condition of Theorem 4.2.2. Since the second condition is a fairly strong restriction, we follow Herrmann et al. [25] and opt to allow clusters violating the condition. Instead, whenever we add a hypernode that a cluster that would violate the second condition of the theorem, we run a cycle detection subroutine to ensure that the operation does not induce a cycle in the coarser hypergraph. Once the clustering has been computed, we contract pairs of hypernodes placed in the same cluster pair by pair, yielding one coarser graph after each contraction. If the coarser hypergraph after the last contraction is still too large, we repeat the algorithm. We stop it once one the hypergraph is small enough (i.e., has less than $180 \cdot k$ hypernodes, the same stopping criterion used in KaHyPar [2]) or one repetition of the algorithm could not find any more non-singleton clusters.

More precisely, our algorithm works as follows. At the beginning of the algorithm, all hypernodes are in their own singleton cluster. For each hypernode u that is still in a singleton cluster, we rate each neighbor using the heavy-edge metric already implemented in KaHyPar. We select the highest rated neighbor v whose cluster can include u without violating the first condition from Theorem 4.2.2. If all hypernodes in v's cluster have the same toplevel as u, we know that we can safely add u to v's cluster without inducing a cycle in the coarser hypergraph. Otherwise, we temporarily add u to v's cluster and search the hypergraph for edges violating the second condition in Theorem 4.2.2, then check whether they induce a cycle in the contracted hypergraph. To be more precise, let the toplevel of hypernodes in v's cluster be t and t + 1. We maintain a queue of hypernodes that are to be processed. Initially, the queue contains all hypernodes in v's cluster with toplevel t. For each hypernode x in the queue, we explore its successors. Whenever we scan a successor y that is in another cluster, we add all hypernodes from y's cluster with toplevel t to the queue. If y is in v's cluster, but x is not, the search found a cycle in the coarsened hypergraph. At this point, we abort the search, remove u from

v's cluster and move on to the next hypernode. If the search does not find a cycle, we leave u in v's cluster and move on. Once the clusters are found, we contract hypernodes in the same cluster pair by pair to obtain a n-level clustering.

4.2.3 Refinement

2-way FM Refinement. On a high level, the 2-way FM refinement algorithm moves hypernodes with the highest gain values between the blocks of a bipartition while making sure to only consider movements that keep the partition acyclic. Over the course of the algorithm, it keeps track of the best bipartition. Once a stopping criterion decides that the refinement is unlikely to find a further improvement of the bipartition, it rollbacks to the best partition found.

More precisely, the algorithm uses two priority queues (one for each block) to keep track of hypernodes and their gain values. Each priority queue contains movable hypernodes in the corresponding block and their gain value, i.e., the change in the connectivity metric when moving the hypernode to the other block. A hypernode is movable if and only if it can be moved to the other block without causing the partition to become cyclic. During 2-way refinement, this is easy to decide: denote the blocks of the bipartition by V_1 and V_2 and assume that V_2 is the successor of V_1 in the quotient graph. Then a hypernode in V_1 can be moved to V_2 if and only if it does not have any successors in V_1 . Analogously, a hypernode in V_2 can be moved to V_1 if and only if it does not have any predecessors in V_2 . Therefore, it is sufficient to keep track on the number of successors or predecessors that a hypernode has in the same block. We implement this using a simple array that we compute once at the start of the uncoarsening phase and then update it appropriately after every uncontraction operation or movement. In particular, we can use this counter to decide whether new hypernodes become movable (counter becomes zero) or unmovable (counter becomes nonzero) after a movement. We then insert those hypernodes into the appropriate priority queue or remove them.

Initially, the priority queues are empty. After uncontracting a hypernode, the resulting hypernodes and their partners are inserted into the priority queues if they are movable. If none of those are movable, the refinement step is skipped and the next hypernode is uncontracted. Otherwise, the algorithm pulls the hypernode with the highest gain value from the priority queue. The hypernode is only moved to the other block if that does not violate the balance constraint. Nonetheless, the hypernode is marked and therefore excluded from the rest of this refinement round. If the node was moved, all unmarked movable neighbors are inserted into their corresponding priority queue and the gain values of all affected hypernodes are updated.

Once the stopping criteria decides that any further improvement of

the partition if unlikely or both priority queues are empty, the algorithm reverts to the best partition found.

k-way FM Refinement. The *k*-way FM refinement aims to improve a given k-way partition and is based on the *k*-way FM refinement algorithm implemented in KaHyPar [2]. The algorithm maintains k priority queues, one queue for each block. Each queue holds hypernodes that can be moved to the block, with the priority being the gain value of the respective move. Viable target blocks for a hypernode depends on the configuration of the refinement algorithm; we consider two alternatives, namely the use of *advanced moves* and *global modes*, described in the following paragraph. We limit the set of movable hypernodes to border hypernodes. The algorithm always performs the best move across all priority queues and after the stopping criterion is reached, the best partition found during the process is restored. Similar to the refinement algorithm of Moreira et al. [42], we consider the following alternatives for viable moves.

Advanced Moves. In this configuration, we only consider hypernode movements that can never create a cycle in the quotient graph. Let $\tau : Q(V) \to [k]$ be a topological order on the quotient graph Q. For a hypernode $u \in V$, let a be the maximum index such that $\tau^{-1}(a)$ contains a predecessor of u and let b be the minimum index such that $\tau^{-1}(b)$ contains a successor of u, i.e., $a = \max\{\tau(b[v]) \mid v \in \Gamma^{-}(u)\}$ and $\tau(b) = \min\{\tau(b[v]) \mid v \in \Gamma^{+}(u)\}$. Note that $a \leq \tau(b[u]) \leq b$ since τ is a topological ordering. We then only consider blocks $\tau^{-1}(a), \ldots, \tau^{-1}(b)$ as viable target blocks for u. This ensures that moving u can never create a cycle in the quotient graph. We therefore do not need any cycle detection algorithm. Since we only need to quotient graph to compute the initial topological ordering, we do not need to keep it up-to-date during refinement.

Global Moves. This configuration considers all adjacent blocks as viable target blocks for a hypernode movement. In particular, it computes gain values for movements that might create a cycle in the quotient graph. Therefore, we must use a cycle detection algorithm to prevent those movements. Initially, we considered the use of advanced online cycle detection algorithms [9,49], but eventually, we settled on Kahn's algorithm [31] since cycle detection did not prove to be a bottleneck in our final implementation. After executing a move, we use the cycle detection algorithm to scan the quotient graph for any cycles. If the move created one, we reverse it, mark the hypernode and remove it from all priority queues. Since the quotient graph might change after every movement, we must keep it up-to-date.

4.3 Larger Imbalances on Coarse Levels

In graph partitioning, temporarily allowing larger imbalances on coarse levels is a well-known technique to improve the resulting partition quality. For instance, Meyerhenke et al. [40] relax the balance constraint on the coarsest level by a constant offset, then narrow it down linearly while unrolling the graph hierarchy until the desired balance constraint is enforced on the finest level of the hierarchy. Previously, Walshaw and Cross [61] formalized this idea and adjusted the allowed imbalance on a given level more carefully, trying to prevent the balance constraint from shrinking too rapidly and therefore losing partition quality. In general, this technique is based on the intuition that a larger imbalance on the coarse levels of the graph hierarchy gives the refinement algorithm more freedom to build partitions that would otherwise be unreachable due to the strict balance constraint. In acyclic DAH partitioning, the partition landscape is even more fractioned than in ordinary graph partitioning due to the additional acyclicity constraints. Hence, we try this technique to further improve the quality of our partitioning algorithm.

Outline. We run the k-way FM refinement algorithm on the coarsest level of the hypergraph hierarchy using a looser imbalance factor $\epsilon' > \epsilon$, producing a partition that violates the ϵ -balance constraint. During the uncoarsening and refinement phase, we then gradually improve the balance of the partition until we end with an ϵ -balanced partition on the finest level. To be more precise, we linearly lower ϵ' to ϵ , i.e., use $\epsilon_i = (\epsilon' - \epsilon)/(n - n_0 + i)$ after the *i*-th uncontraction operation, where n_0 denotes the number of hierarchies. To improve the balance of the partition, we use a soft rebalancing algorithm that moves hypernodes based on their gain value to improve the balance while not increasing the connectivity metric. If this step is unable to sufficiently improve the balance of the partition, we follow up with a hard rebalancing algorithm that moves hypernodes from overloaded blocks at the cost of increasing the connectivity metric. After balance is restored, we run the k-way FM refinement algorithm from Chapter 4.2.3 initialized with all hypernodes.

4.3.1 Soft Rebalance

The soft rebalancing algorithm is a version of the FM algorithm that only performs *Advanced Moves*. It accepts a partition if it lowers the imbalance of the partition while not increasing its connectivity.

The priority queues are organized as suggested by Träff [60]. We use one priority queue for each block as well as one priority queue for each hypernode. The queue

corresponding to a hypernode contains an entry for each block that the hypernode can be moved to with its corresponding gain value as priority. The queue corresponding to a block contains all movable hypernodes in that block indexed by their highest possible gain value, i.e., the maximal key in their hypernode priority queue.

The queues are initialized with all movable hypernodes at the start of the uncoarsening and refinement phase. After each uncontraction operation, one round of the algorithm is executed. A round of the soft rebalancing algorithm works as follows. It enables all priority queues that belong to overloaded blocks and pulls the hypernode with the maximal gain across all queues. The target of the move is determined by the maximal element in the hypernode priority queue. If the move reduces the imbalance of the partition, the hypernode is moved to the target block, the priority queues and gains are updated accordingly and the hypernode is inserted into the priority queue of the target block. If the hypernode was not moved, the inviable target block is removed from the hypernode's priority queue and is re-inserted into the block's priority queue.

The round terminates if the improved balance of the partition meets the current imbalance factor or a certain number of moves was performed without producing a viable partition. The algorithm then reverts hypernode movements until the last accepted partition.

4.3.2 Hard Rebalance

Since the algorithm described in the last section is not guaranteed to sufficiently improve the balance of the partition, we also present an algorithm that always succeeds at the cost of a higher connectivity metric. The hard rebalancing algorithm selects an overloaded block and an underloaded block. It then moves nodes along subsequent blocks in the topologically ordered quotient graph from the overloaded block to the underloaded block. The blocks are selected such that the sum of the gain values of all movements is maximal.

More precisely, the algorithm keeps two priority queues for each block, for a total of 2k priority queues. One queue contains all hypernodes that can be moved to the previous block while the other one contains all hypernodes movable to the next block. The priorities are the corresponding gain values. Let $\tau : Q(V) \to [k]$ be a topological ordering of the quotient graph Q. Moving hypernodes from one block $\tau^{-1}(i)$ to block $\tau^{-1}(j)$ with i < j involves moving one hypernode from block $\tau^{-1}(i)$ to block $\tau^{-1}(i+1)$, one from $\tau^{-1}(i+1)$ to $\tau^{-1}(i+2)$ and so on. We therefore select a pair of an overloaded and an underloaded block such that the sum of the gain values of all of those moves is maximal among all possible (overloaded, underloaded) pairs of blocks. Note that this is not a precise approach to find the best pair of blocks, since moves might change the gain values of subsequent moves, although

our experiments indicate that it is a good approximation in practice. After moving a hypernode, we insert it into the priority queue of its new block (if it is still movable) and update gain values of adjacent hypernodes.

This algorithm can always restore the balance of a DAH partition. To see this, we need Lemma 4.3.1.

Lemma 4.3.1. Let H be a DAH with an acyclic k-way partition Π , and topological ordering $\tau : \Pi \to [k]$ of the quotient graph. Then every block $\tau^{-1}(i)$, $1 < i \leq k$ has at least one hypernode movable to block $\tau^{-1}(i-1)$ and every block $\tau^{-1}(j)$, $1 \leq j < k$ has at least one hypernode movable to block $\tau^{-1}(i-1)$.

Proof. Recall that a H can be seen as a DAG G by replacing each hyperedge e with a directed bipartite graph from e^T to e^H . This transformed graph is acyclic if and only if the DAH is acyclic. Hence, we can prove the statement for DAGs rather than DAHs. Let $1 < i \leq k$. Consider the block-induced subgraph $G' := G[\tau^{-1}(i)]$. Since G is a DAG, so is G' and we can use Lemma 2.1.1 to see that G' contains a node $u \in \tau^{-1}(i)$ with indegree zero in G'. Moving this node to block $\tau^{-1}(i-1)$ does not create a backward edge in the quotient graph with respect to τ , proving the first statement of the lemma. The second statement follows analogously.

Moreover, observe that the topological ordering of the k-way partition does not change during the coarse of the algorithms: Since we only move hypernodes between subsequent blocks, the moves can only create new quotient graph edges between subsequent blocks. Hence, since the topological ordering is static during the course of the algorithm, and since we always have movable hypernodes, the algorithm can always perform movements until the desired balance is reached.

However, this is only true when working on a DAH. If the hypergraph contains cycles, it might fail when a block has no movable hypernodes. While our input hypergraphs are always acyclic, one of the coarsening algorithms described in Chapter 4.2.2 might produce cyclic coarser hypergraphs. In this case, we simply stop the rebalancing step and try again after uncontracting the next hypernode. Since the finest hypergraph is acyclic, the algorithm succeeds eventually.

4.4 Memetic Acyclic Hypergraph Partitioning

In Chapter 3.4, we referenced memetic algorithms as a successful metaheuristic for high-quality graph and hypergraph partitioning. Now, we present a memetic algorithm for the DAH partitioning problem. We use the memetic algorithm already implemented in KaHyPar [4] as framework and exchange its building blocks for mutation, recombination and generating the initial partition with new algorithms. From a metaheuristic point of view, the algorithm remains unchanged. Hence, we follow the description from Ref. [4] closely.

Population. The memetic algorithm starts by generating the initial population \mathcal{P} . The population consists of individuals, which are always ϵ -balanced k-way partitions of the input hypergraphs. We generate the initial individuals using our multilevel algorithm introduced in previous chapters. The size of \mathcal{P} is choosing dynamically by measuring the time t_I it takes to generate one individuals, i.e., the running time of our multilevel algorithm: $|\mathcal{P}| \coloneqq \max(3, \min(50, \delta \cdot (t/t_I)))$, where δ is a configuration parameter that we set to 0.15.

The fitness of an individual is its connectivity, since that is the objective that we want to optimize. An individual with lower connectivity is fitter than one with higher connectivity. The initial population is evolved over several generational cycles using the *steady-state* paradigm [20]: We generate only a single offspring per generation. When inserting a new individual I_1 into the population, we evict an old one I_2 based on similarity: the difference between both individuals is defined as $d(I_1, I_2) := |D(I_1) \ominus D(I_2)|$, where D(I) is a multi-set that contains each hyperedge $e \in E$ exactly $\lambda(e) - 1$ times and \ominus is the symmetric difference. In other words, we consider two individuals to be somewhat similar if all hyperedges have roughly the same connectivity in both partitions.

Recombination Operator. For recombination, we select parents using binary tournament selection [41]: First, we select two individuals at random and choose the fitter one as first parent P_1 . Then, we repeat the process to select the second parent P_2 . We then run a modified V-cycle to combine both parents. During coarsening, we only allow the contraction of two hypernodes u and v if they are in the same block in both P_1 and P_2 , i.e., if $b_1[u] = b_1[v]$ and $b_2[u] = b_2[v]$. This allows us to use P_1 or P_2 as initial partition once coarsening has terminated. We apply the fitter one of both parents as initial partition. Finally, we unroll the graph hierarchy and improve the partition using our k-way refinement algorithm. Note that this recombination operator produces offsprings that are at least as good as the better of both parents.

Mutation Operations. We implement two mutation operations. The first one starts by selecting a random individual I, then perform a modified V-cycle that works as follows. During coarsening, it only contracts pairs of hypernodes u and v that are in the same block in I, i.e., b[u] = b[v]. On the coarsest level of the graph hierarchy, I is used as initial partition and the graph hierarchy is unrolled and improved using our k-way refinement algorithm. This operation produces an offspring that is at least as good as I. The second mutation operation also selects a random individual I_1 , then generates a new individual I_2

as described earlier. Next, both individuals get recombined, but we always use I_2 as initial partition. This operator can therefore produce offsprings that are worse than the individual selected from the population.

4 Acyclic Hypergraph Partitioning

5 Experimental Evaluation

In this chapter, we evaluate the performance of our algorithm and the influence of various components on solution quality. We start by presenting our methodology, the systems used for evaluation and the benchmark setup. We then present our main results by comparing our algorithm to previous DAG partitioning approaches in Chapter 5.1. For these experiments, we only use DAG instances, since the other algorithm cannot cope with hypergraphs. Afterwards, we compare our memetic multilevel algorithm to simpler approaches for DAH partitioning in Chapter 5.2 and evaluate the impact of DAH partitioning on an image streaming application in Chapter 5.3. Finally, we evaluate the influence of the acyclic coarsening algorithm and present partitioning approaches that did not prove to be beneficial in Chapter 5.2.2 and Chapter 5.2.1, respectively.

Methodology and Setup. We implement all algorithms described in Chapter 4 in the KaHyPar hypergraph partitioning framework [2,4,27,28,54,55]. We use mlDHGP to refer to our multilevel algorithm and memDHGP for our memetic multilevel algorithm. The code is written in C++ and compiled using g++ 9.1 using -O3-march=native as compile flags. Our implementation is based on the KaHyPar version from march 2019¹. The source code of our preliminary version is available at github.com/danielseemaier/kahypar/tree/HyperDAG, and we plan to integrate our algorithm into the next release of KaHyPar available at kahypar.org.

We ran our experiments on two different machines. Machine A is a single node from the HPC cluster bwUniCluster equipped with two Intel Xeon E5-2670 Octa-Core (Sandy Bridge) processor clocked at 2.6 GHz, 64 GB main memory, 20 MB L3-Cache and 8x256 KB L2-Cache. Machine B has two Intel Xeon E5-2650 v2 Octa-Core (Sandy Bridge) clocked at 2.6 GHz, 128 GB main memory, 20 MB L3-Cache and 8x256 KB L2-Cache.

Performance Profiles. We use performance profiles [22] to compare the solution quality of different algorithms. The plots show one curve for each algorithm included in the comparison. The x-axis shows $\tau \in [1, 100]$ and the y-axis shows the fraction

¹Commit hash 84c7e7c523701efb0e51752053656e34d206cf4c, repository github.com/ SebastianSchlag/kahypar.

of instances for which each algorithm computed a partition that is within a factor of τ of the best partition computed by any algorithm for that instance. In particular, $\tau = 1$ shows the fraction of instances for which each algorithm computed the best partition (i.e., partition with the lowest connectivity). A value of 0.8 on the y-axis for $\tau = 1$ shows that the algorithm computed the best result on 80% of all instances and a value of 1 for $\tau = 1.1$ reveals that the algorithm computes partitions within a factor of 1.1 of the partition of the best algorithm on *every* instance.

Instances. We use three sets of graphs for our experimental evaluation. The first set of graphs are deduced from the PolyBench Benchmark Suite [51]. Those graphs were kindly provided to us in DAG format by Herrmann et al. and were also used for evaluation in previous papers on DAG partitioning [25, 26, 43]. Next, we use graphs from the ISPD98 Circuit Benchmark Suite [3]. These instances contain one node for each cell and a directed edge from the source of a net to each of its sinks. In case the resulting instance does not form a DAG, i.e., contains cycles, we do the following: We gradually add directed edges and skip those that would create a cycle. We use these instances for our main experiments. Some of our initial experiments use DAGs based on the circuits from the ISCAS85 Combinational Benchmark Circuits [13]. Those DAGs contain one node for each logic gate and a directed edge from the output of a logic gate to the input of another logic gate. Basic properties of these instances are provided in Table 5.1.

To conduct experiments on hypergraphs, we transform all DAGs into DAHs using the row-net model on their adjacency matrices: A hypergraph contains one hypernode for each node in the DAG and a hyperedge for each node u with outgoing edges. The head of the hyperedge is u and the tails are the successors of u.

When evaluating the impact of DAH partitioning on an image streaming application, we use DAGs modeling the data flow of an advanced imaging algorithm [48] and transform them into DAHs. This transformation works differently from the previous description and is described in Chapter 5.3.

5.1 DAG Model

We start our experimental evaluation by comparing our algorithm to the current state-of-the art on DAG partitioning. Note that the comparison is possible even though previous algorithms optimize the edge-cut on DAGs, whereas our algorithm optimizes the connectivity metric on DAHs since both objectives are equal for 2-uniform hypergraphs. The comparisons includes the algorithm by Herrmann et al. [25] and the one by Moreira et al. [43]. We name the former algorithm HOUKC (first letters of the authors last names), the latter one Moreira (last name of the first

Graph	n	m	Ref.	Graph	n	m	Ref.
	PolyBen	ch			ISPE	98	
2mm	36500	62200	[51]	ibm01	13865	42767	[3]
$3\mathrm{mm}$	111900	214600	[51]	ibm02	19325	61756	[3]
adi	596695	1059590	[51]	ibm03	27118	96152	[3]
atax	241730	385960	[51]	ibm04	31683	108311	[3]
covariance	191600	368775	[51]	ibm05	27777	91478	[3]
doitgen	123400	237000	[51]	ibm06	34660	97180	[3]
durbin	126246	250993	[51]	ibm07	47830	146513	[3]
fdtd-2d	256479	436580	[51]	ibm08	50227	265392	[3]
gemm	1026800	1684200	[51]	ibm09	60617	206291	[3]
gemver	159480	259440	[51]	ibm10	74452	299396	[3]
gesummv	376000	500500	[51]	ibm11	81048	258875	[3]
heat-3d	308480	491520	[51]	ibm12	76603	392451	[3]
jacobi-1d	239202	398000	[51]	ibm13	99176	390710	[3]
jacobi-2d	157808	282240	[51]	ibm14	152255	480274	[3]
lu	344520	676240	[51]	ibm15	186225	724485	[3]
ludcmp	357320	701680	[51]	ibm16	189544	648331	[3]
mvt	200800	320000	[51]	ibm17	188838	660960	[3]
seidel-2d	261520	490960	[51]	ibm18	201648	597983	[3]
symm	254020	440400	[51]				
syr2k	111000	180900	[51]				
syrk	594480	975240	[51]				
trisolv	240600	320000	[51]				
trmm	294570	571200	[51]				
		IS	SCAS8	5			
c432	196	336	[13]	c2670	1 4 2 6	2075	[13]
c499	243	408	[13]	c3540	1719	2936	[13]
c880	443	729	[13]	c5315	2485	4386	[13]
c1355	587	1064	[13]	c6288	2448	4800	[13]
c1908	913	1497	[13]	c7552	3719	6144	[13]

 $\label{eq:table 5.1: Basic properties of DAG instances.$

author), our own multilevel algorithm mlDHGP+X, and our own memetic algorithm memDHGP+X, where $X \in \{KaHyPar, PaToH\}$ is the undirected hypergraph partitioner used during initial partitioning. Whenever we omit +X, we default to KaHyPar since it showed the best results. When using +KaHyPar, we invoke KaHyPar using the newest configuration available² and when using +PaToH, we invoke it using the default configuration. To run the HOUKC algorithm on our benchmark set, we use the implementation publicly available on the first author's website³. Since the author's implementation of Moreira belongs to Intel and is therefore not publicly available, we refrain from running it on our benchmark set and opt to use the numbers provided in Ref. [43]. We have therefore no data for this algorithm on the ISPD98 instances, but can still compare it to the other algorithms on the PolyBench instances.

We partition each graph in our benchmark set 5 times for different values of $k \in \{2, 4, 8, 16, 32\}$ and use $\epsilon = 3\%$ as imbalance parameter. These values were chosen because they were also used in previous work on DAG partitioning [25,43]. Each algorithm gets allocated 8 hours (except for adi, for which we give all solvers 24 hours) for each instance of our benchmark set. We run them on a single core of Machine A. While the memetic algorithms take the time limit as an input parameter, non-memetic algorithms do not. We therefore run them repeatedly using random seeds until the time limit exceeds.

The effect of the undirected hypergraph partitioner used during initial partitioning and of our memetic component is summarized in Figure 5.1. As we can see in Figure 5.1a, using a state-of-the undirected hypergraph partitioner pays off: With KaHyPar, our algorithm computes the same or a better solution on 65% out of all instances (PolyBench and ISPD98), whereas the same is only true for 45% out of all instances when using it with PaToH. Hence, we will only consider memDHGP+KaHyPar and m1DHGP+KaHyPar in the rest of the experiments. Moreover, Figure 5.1b summarizes the influence of our memetic component. We observe the memDHGP computes a strictly better partition on 65% out of all instances compared to repeated runs using random seeds (i.e., m1DHGP) with the largest improvement made on 2mm with k = 8 (10%) and ibm01 with k = 2 (8%). Based on these observations, we conclude that our memetic algorithm is more effective than repeated restarts of our multilevel algorithm.

Next, we compare our algorithm to HOUKC and Moreira. The best edge cuts found by each algorithm are compared in Figure 5.2. Detailed results with per- instance edge cuts are available in Tables A.1–A.4. We observe that memDHGP+KaHyPar outperforms the other algorithms on both the PolyBench and ISPD98 instances. Looking at Figure 5.2a, we see that it computes the best partition on over 82% of

²github.com/SebastianSchlag/kahypar/blob/master/config/km1_kahypar_mf_jea19. ini

³people.bordeaux.inria.fr/julien.herrmann

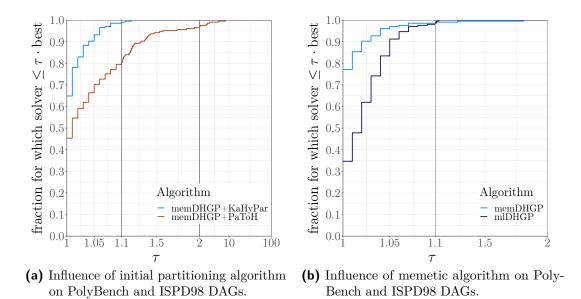


Figure 5.1: Influence of the algorithm used for undirected initial partitioning and of our memetic algorithm on PolyBench and ISPD98 DAGs.

all PolyBench instances, while HOUKC and Moreira only compute the best partition on 25% and 14%, respectively. Moreover, memDHGP+KaHyPar is within a factor of 1.1 of the best algorithm on over 95%, whereas the other algorithms are only within a factor of 1.1 of the best algorithm on 51% and 22%, respectively. The ISPD98 instances reveal a similar observation: Here, memDHGP+KaHyPar computes the best solution on over 87% instances compared to HOUKC. The average improved to the previous state-of-the art HOUKC is 11.1% (PolyBench instances) and 9.7% (ISPD98 instances). The best improved is observed on the graph covariance from the PolyBench instances with k = 2: While HOUKC computed an edge cut of 34 307, our algorithm finds a partition with an edge cut of 11 281.

We finish this chapter by taking a look at the running times of mlDHGP+KaHyPar and HOUKC in Figure 5.2c. Note that we exclude Moreira from this comparison since we did not run their code ourself and the authors did not report per-instance running times for their algorithm. As can be seen in the figure, mlDHGP+KaHyPar is slower than HOUKC by several orders of magnitude. This has several reasons: HOUKC uses METIS [33] during initial partition rather than KaHyPar, only uses 2-way refinement, uses a multilevel scheme with only log(n) levels rather than n levels, and only supports DAGs and no DAHs. Computing and updating gain values is faster when optimizing the edge cut objective on graphs than the connectivity objective on hypergraph. In particular, due to the acyclicity constraint, edge cut gain values

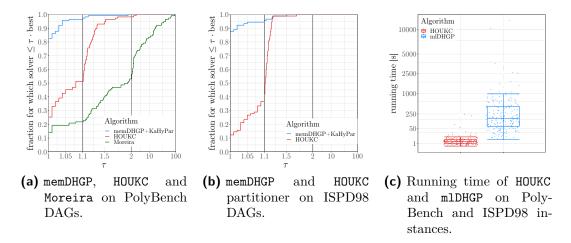


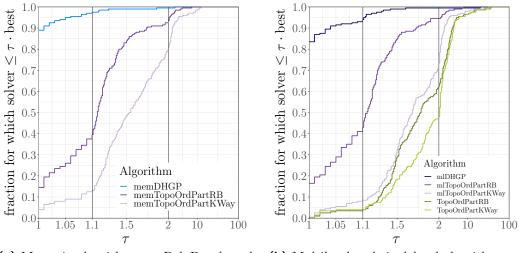
Figure 5.2: Performance and running time of DAG partitioners on PolyBench and ISPD98 DAGs.

never change when moving nodes between the blocks of a bipartition of a normal graph. Hence, HOUKC does not need to update gain values during 2-way refinement.

5.2 DAH Model

After establishing the state-of-the art for DAG partitioning, we move on the DAH instances. For these experiments, we use the PolyBench and ISPD98 instances transformed into DAHs as described above, but exclude the adi graph from the PolyBench instances since it is too large to be partitioned by our algorithm within a reasonable time frame. We use the same values for k and ϵ as before. Since we are not aware of any other algorithms for DAH partitioning, we compare the performance of our algorithm to simpler heuristics, namely TopoOrdPartRB and **TopoOrdPartKWay**. The former one partitions the graph using recursive bisection. For each bisection, it uses Kahn's algorithm to compute a topological order of the graph, splits the graph based on that, and then uses our 2-way refinement algorithm to improve the bisection. The latter one again uses Kahn's algorithm to compute a topological order, but then directly splits the graph into k blocks and improves it using our k-way refinement algorithm. We also test these heuristics with multilevel refinement and name those algorithms mlTopoOrdPartRB and mlTopoOrdPartKWay. Finally, we use those partitions as input to our memetic algorithm and name the result memTopoOrdPartRB and memTopoOrdPartKWay. As before, we give the memetic algorithm 8 hours time and use Machine A to perform these experiments.

The results are shown in Figure 5.3 and Table 5.2 with detailed per-instance



(a) Memetic algorithms on PolyBench and ISPD98 DAHs.

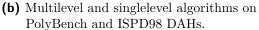


Figure 5.3: Comparison of memDHGP and mlDHGP to simpler approaches.

results available in Tables B.1–B.4. Looking at the performance profile shown in Figure 5.3a, we observe that memDHGP outperforms the simpler heuristics, computing the best result on almost 90% out of all instances, whereas the second best approach, memTopoOrdPartRB, only computes the best result on less than 15% out of all instances. On average, memDHGP computes partitions with 20% lower connectivity than memTopoOrderPartRB. We can therefore conclude that using a high quality undirected hypergraph partitioner during initial partitioning improves the overall result significantly.

Focusing on Figure 5.3b and Table 5.2, we see that using recursive bisection gives significantly better results than using direct k-way partitioning. Looking at the single-level algorithms TopoOrderPartRB and TopoOrderPartKWay, the approach using recursive bisection is 10% better on average than direct k-way. This advantage increases when looking at the multilevel algorithms: Here, recursive bisection improves partitions by 49%. We believe that this is due to the fact that the k-way search space is much more fractured than the 2-way search space due to the acyclicity constraint. Moreover, we observe that the multilevel algorithms are better than the single-level algorithms; this is expected as the multilevel component adds a more global view to the optimization landscape.

Algorithm	gmean
TopoOrderPartRB	16 571
mlTopoOrderPartRB	9128
memTopoOrderPartRB~(8h)	8 071
TopoOrderPartKWay	18 161
mlTopoOrderPartKWay	13605
$\verb memTopoOrderPartKWay (8h) $	10 643
mlDHGP	7 2 4 4
memDHGP (8h)	6526

Table 5.2: Geometric mean solution quality of the best results out of multiple repetitions for different algorithms on PolyBench and ISPD98 instances (as DAHs).

5.2.1 Influence of Larger Imbalances on Coarse Levels

Next, we evaluate the influence of allowing a larger imbalance on coarse levels. This algorithm is described in Chapter 4.3. We executed the experiment on Machine B and used the ISCAS85 and PolyBench DAHs with $k = 2, 4, 8, 16, 32, \epsilon = 3\% = 0.03$ (maximum imbalance of the resulting partition) and various values for ϵ' (imbalance on the coarsest level of the hypergraph hierarchy).

The result of this experiments is summarized in Figure 5.4. In our initial experiment, we obtained an initial k-way partition using a topological ordering of the DAH, i.e., using mlTopoOrdKWay from the previous chapter. We then ran a V-cycle with larger imbalance on coarse levels as described in Chapter 4.3. The result of this run can be seen in Figure 5.4a. We observe that using $\epsilon' = 20\%$ computed the best partition on more than on third out of all instances, whereas the configuration that does not allow a larger imbalance on coarse levels ($\epsilon' = \epsilon = 3\%$) only computes the best partition on 28% out of all instances. Moreover, when using $\epsilon' = 7\%$, the computed partition is within a factor of 1.1 of the best partition on 74% out of all instances, whereas the same can only be said for 54% out of all instances when not allowing a larger imbalance on coarse levels. Based on this, it seems that this techniques can improve a partition computed using mlTopoOrdKWay on ISCAS85 instances. However, as we have seen in the previous chapter, mlTopoOrdKWay computes partitions of low quality while mlDHGP computes much better partitions. Hence, we repeated the same experiment with mlDHGP instead of mlTopoOrdKWay. With this configuration, we see in Figure 5.4b that $\epsilon' \in \{7\%, 20\%\}$ no longer compute better partitions, but worsen the overall result. The same observation applies when partitioning the PolyBench instances, as can be seen in Figure 5.4c. Here, we see that not allowing larger imbalances on coarse

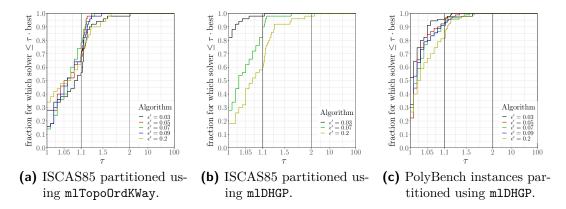


Figure 5.4: Influence of larger imbalance on coarse levels for ISCAS85 and PolyBench DAHs and various values for ϵ' .

levels computes the best partition on most instances and generally outperforms all other configurations, and partition quality decreases with increasing values of ϵ' .

In our experiments, we observed that almost all balance improvements stem from our hard rebalancing algorithm that decreases partition quality. This decrease in partition quality outweighs the improvement found during the initial k-way refinement with larger imbalance on the coarsest level and all improvements found during uncoarsening. We think that this is due to the observation that k-way refinement and our soft rebalancing algorithm are too restricted by the acyclicity constraint.

5.2.2 Influence of Acyclic Coarsening

In this experiment, we evaluate the influence of our acyclic coarsening algorithm presented in Chapter 4.2.2 by comparing it to the coarsening algorithm that is already implemented in KaHyPar [2]. We used Machine B to execute this experiment. The instances are ISPD98 DAHs that were partitioned into k = 2, 4, 8, 16, 32 blocks with maximum imbalance $\epsilon = 3\%$.

The average connectivity using the acyclic coarsening is 16 086, whereas the coarsening algorithm already implemented in KaHyPar yields an average connectivity of only 20 915. Hence, the acyclic coarsening algorithm produces partitions with 23% lower connectivity on average. This improvement can also be seen in Figure 5.5: With acyclic coarser hypergraphs, we compute strictly better partitions on almost all instances. Using the coarsening algorithm that is already implemented in KaHyPar, we only get within a factor of 1.1 of the partition computed using the acyclic coarsening algorithm on 10% out of all instances. We believe that this is due to the fact that hypernodes in a cycle cannot be moved to another block

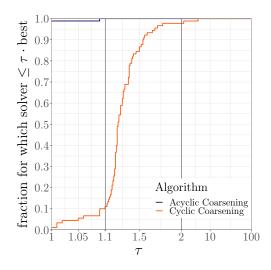


Figure 5.5: ISPD98 instances partitioned using the acyclic coarsening algorithm and the one that is already implemented in KaHyPar and does not keep the hypergraph acyclic.

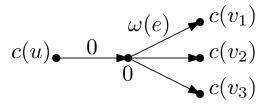
without violating the acyclicity constraint, therefore restricting the refinement algorithm. Judging from this experiment, we conclude that acyclic coarsening algorithms are necessary to obtain high-quality acyclic DAH partitions. This was already observed by Herrmann et al. [25] for the DAG case.

5.3 Impact on Streaming Application

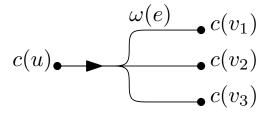
The experiment described in this chapter was conducted by Merten Popp. The author of this thesis only implemented the transformation of DAGs into DAHs.

To evaluate the impact of our DAH partitioning algorithm on image streaming applications, we integrate the algorithm into a toolchain implementing the *Local Laplacian filter* that was also used to evaluate earlier work on DAG partitioning [42, 43]. The filter is an edge-aware image processing filter using concepts of *Gaussian pyramids* and *Laplacian pyramids* as well as a point-wise remapping function to enhance image details without creating artefacts. For more details on the image processing algorithm see Ref. [48].

The toolchain implementing the algorithm invokes our algorithm on a DAG structured as depicted in Figure 5.6a. The DAG consists of *compute nodes* and *output nodes*. Compute nodes correspond to parts of the filter application and have the size of the program as node weight. Output nodes are merely conceptual and have node weight zero. Each compute node can have an arbitrary number of output nodes as successors; the edge from compute node to output node has



(a) DAG model: Nodes with nonzero weight are *compute nodes*, nodes with zero weight are *output nodes*. One compute node can have multiple output nodes. The weight of edges from output nodes to compute nodes correspond to the amount of data that has to be transferred. Note that all outgoing edges from an output node have the same weight.

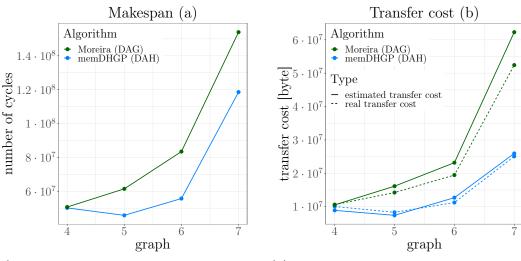


(b) DAH model: Each output node is replaced by a single hyperedge that models the data transfer.

Figure 5.6: Structure of the data dependency and execution flow graphs, once as DAG (Figure 5.6a) and once as DAH (Figure 5.6b). The DAGs are kindly provided to us by Moreira et al. and we perform the transformation to the corresponding DAHs ourself.

edge weight zero. The successors of output nodes are other compute nodes, which depend on the output of the preceding compute node. All outgoing edges from an output node have the same edge weight, namely the size of the compute node's output. To benefit from DAH partitioning, we transform the supplied DAG into a DAH as depicted in Figure 5.6b. Each output node is replaced with a single hyperedge that has the same hyperedge weight as its outgoing edges. The head of the output hyperedge is the preceding compute node and its tails are the succeeding compute nodes. This DAH is then partitioned using memDHGP with a time limit of 10 minutes and the final partition is projected back onto the original DAG by placing output nodes into the same block as its preceding compute node.

The filtering algorithm is configurable by parameter K, the number of levels that the image pyramids have. A higher value for K improves the filters result, but also increases the size of the DAG (DAH). To evaluate our algorithm, we use values $K \in \{4, 5, 6, 7\}$, resulting in a DAG with at most n = 752 and m = 862 for K = 7. We compare the impact of DAH partitioning to Moreira, the algorithm previously used in the toolchain. The result is illustrated by Figure 5.7. By looking at Figure 5.7a, we observe that DAH partitioning generally results in a lower makespan compared to DAG partitioning. The best improvement is a reduction in makespan by 22%, observed on the largest filtering algorithm, i.e., K = 7. Figure 5.7b hints that this improvement is in part due to a more accurate modeling of transfer costs. The figure compares the edge cuts reported by the respective



(a) Makespan of each filtering algorithm. The values are obtained using a cycle-true compiled simulator of the target platform. Lower is better.

(b) Edge cut as reported by the respective algorithm (i.e., estimated data transfer) versus actual data transfer. Closer gap is better.

Figure 5.7: Impact of the DAH model with improved partitioning results on image streaming applications: Lower makespan and more accurate modeling of the transfer cost.

algorithms (i.e., the estimated transfer cost) to the actual transfer costs. We observe that the DAH model indeed estimates the transfer cost much better than the DAG model, confirming our initial hypothesis. Based on this, we conclude that the DAH model is better suited than the DAG model in this application domain.

6 Conclusion

Motivated by the shortcomings of the DAG model for representing dataflow and execution dependencies in image streaming applications, we presented the first hypergraph partitioning algorithm for directed acyclic hypergraphs that can cope with the acyclicity constraint. We implemented our novel algorithm using the hypergraph partitioning framework KaHyPar and ran extensive experiments to benchmark its components and to compare it to the previous state-of-the art on DAG partitioning.

Indicated by our experimental evaluation, we observed that our *n*-level partitioning algorithm improved on the current state-of-the-art algorithm for DAG partitioning: Compared to previous algorithms, we computed the best partitions on 82% out of all instances in our benchmark set and improved the edge cut by 10%on average. Based on this, we concluded that our algorithm can be seen as the new state-of-the art for DAG partitioning in terms of partitioning quality. Since there are no previous algorithms for DAH partitioning, we compared our algorithms to more simple heuristics and showed significant improvements. Compared to direct *k*-way partitioning based on a topological ordering with single-level FM refinement, we showed that our memetic algorithm could lower the connectivity of the resulting partition by 64% on average. Getting back to the problem that motivated our work, we showed that acyclic DAH partitioning allows a significantly more efficient parallelization of image streaming applications on embedded devices, improving the makespan of an advanced image filter by 22%. We could therefore conclude that our algorithm outperforms previous DAG partitioner that were used in the application domain.

6.1 Future Work

We have left several paths for future extensions and improvements. First, our current implementation can only handle DAHs where every hyperedge has at most one head. This limitation could be lifted with changes to the coarsening or refinement step. Secondly, since k-way local search is a bottleneck of our algorithm, future work could work towards parallelized refinement algorithms. Lastly, we plan to integrate our algorithm into the next KaHyPar release.

 $6 \ Conclusion$

Bibliography

- [1] Nitin Ahuja, Matthias Bender, Peter Sanders, Christian Schulz, and Andreas Wagner. Incorporating road networks into territory design. In *Proceedings of* the 23rd SIGSPATIAL International Conference on Advances in Geographic Information Systems, SIGSPATIAL '15, New York, NY, USA, 2015. Association for Computing Machinery.
- [2] Yaroslav Akhremtsev, Tobias Heuer, Peter Sanders, and Sebastian Schlag. Engineering a direct k-way hypergraph partitioning algorithm. In 2017 Proceedings of the Ninteenth Workshop on Algorithm Engineering and Experiments (ALENEX), pages 28–42.
- Charles J. Alpert. The ispd98 circuit benchmark suite. In Proceedings of the 1998 International Symposium on Physical Design, ISPD '98, pages 80–85, New York, NY, USA, 1998. ACM.
- [4] Robin Andre, Sebastian Schlag, and Christian Schulz. Memetic multilevel hypergraph partitioning. In *Proceedings of the Genetic and Evolutionary Computation Conference*, GECCO '18, pages 347–354, New York, NY, USA, 2018. ACM.
- [5] Konstantin Andreev and Harald Räcke. Balanced graph partitioning. In Proceedings of the Sixteenth Annual ACM Symposium on Parallelism in Algorithms and Architectures, SPAA '04, page 120–124, New York, NY, USA, 2004. Association for Computing Machinery.
- [6] Giorgio Ausiello, Paolo G. Franciosa, and Daniele Frigioni. Directed hypergraphs: Problems, algorithmic results, and a novel decremental approach. In *Theoretical Computer Science*, pages 312–328, Berlin, Heidelberg, 2001. Springer Berlin Heidelberg.
- [7] Giorgio Ausiello and Luigi Laura. Directed hypergraphs: Introduction and fundamental algorithms—a survey. *Theoretical Computer Science*, 658:293 – 306, 2017. Horn formulas, directed hypergraphs, lattices and closure systems: related formalism and application.
- [8] Jørgen Bang-Jensen and Gregory Z. Gutin. Digraphs: Theory, Algorithms and Applications. Springer-Verlag London, 2nd edition, 2009.
- [9] Michael A. Bender, Jeremy T. Fineman, Seth Gilbert, and Robert E. Tarjan.

A new approach to incremental cycle detection and related problems. ACM Trans. Algorithms, 12(2), December 2015.

- [10] Claude Berge and Edward Minieka. Graphs and Hypergraphs. North-Holland mathematical library. North-Holland Publishing Company, 1976.
- [11] Sonja Biedermann, Monika Henzinger, Christian Schulz, and Bernhard Schuster. Memetic Graph Clustering. In 17th International Symposium on Experimental Algorithms (SEA 2018), volume 103 of Leibniz International Proceedings in Informatics (LIPIcs), pages 3:1–3:15, Dagstuhl, Germany, 2018. Schloss Dagstuhl-Leibniz-Zentrum fuer Informatik.
- [12] Marcel Birn, Vitaly Osipov, Peter Sanders, Christian Schulz, and Nodari Sitchinava. Efficient parallel and external matching. In *Euro-Par 2013 Parallel Processing*, pages 659–670, Berlin, Heidelberg, 2013. Springer Berlin Heidelberg.
- [13] Franc Brgles and Hideo Fujiwara. A neutral netlist of 10 combinational circuits and a target translator in fortran. *IEEE International Symposium on Circuits* and Systems, 1985.
- [14] Aydin Buluç, Henning Meyerhenke, Ilya Safro, Peter Sanders, and Christian Schulz. Recent Advances in Graph Partitioning. In Algorithm Engineering -Selected Results and Surveys, pages 117–158. Springer, 2016.
- [15] Umit V. Catalyürek and Cevdet Aykanat. Hypergraph-Partitioning-Based Decomposition for Parallel Sparse-Matrix Vector Multiplication. *IEEE Transactions on Parallel and Distributed Systems*, 10(7):673–693, Jul 1999.
- [16] Jason Cong and Yuzheng Ding. On area/depth trade-off in lut-based fpga technology mapping. In *Proceedings of the 30th International Design Automation Conference*, DAC '93, page 213–218, New York, NY, USA, 1993. Association for Computing Machinery.
- [17] Jason Cong and Sung Kyu Lim. Performance driven multiway partitioning. In Proceedings of the 2000 Asia and South Pacific Design Automation Conference, ASP-DAC '00, page 441–446, New York, NY, USA, 2000. Association for Computing Machinery.
- [18] Jason Cong and Sung Kyu Lim. Edge separability based circuit clustering with application to circuit partitioning. In *Proceedings 2000. Design Automation Conference. (IEEE Cat. No.00CH37106)*, pages 429–434, Jan 2000.
- [19] Jason Cong, Zheng Li, and Rajive Bagrodia. Acyclic multi-way partitioning of boolean networks. In 31st Design Automation Conference, pages 670–675, June 1994.
- [20] Kenneth A. De Jong. Evolutionary computation a unified approach. MIT Press, 2006.

- [21] Karen D. Devine, Erik G. Boman, Robert T. Heaphy, Rob H. Bisseling, and Umit. V. Catalyürek. Parallel Hypergraph Partitioning for Scientific Computing. In 20th International Conference on Parallel and Distributed Processing (IPDPS), pages 124–124. IEEE, 2006.
- [22] Elizabeth D. Dolan and Jorge J. Moré. Benchmarking optimization software with performance profiles. *Mathematical Programming*, 91(2):201–213, 2002.
- [23] Charles M. Fiduccia and Robert M. Mattheyses. A linear-time heuristic for improving network partitions. In *Proceedings of the 19th Design Automation Conference*, DAC '82, pages 175–181, Piscataway, NJ, USA, 1982. IEEE Press.
- [24] Bruce Hendrickson and Robert Leland. A multilevel algorithm for partitioning graphs. In *Proceedings of the 1995 ACM/IEEE Conference on Supercomputing*, Supercomputing '95, New York, NY, USA, 1995. ACM.
- [25] Julien Herrmann, Jonathan Kho, Bora Uçar, Kamer Kaya, and Umit Çatalyürek. Acyclic partitioning of large directed acyclic graphs. In 2017 17th IEEE/ACM International Symposium on Cluster, Cloud and Grid Computing (CCGRID), pages 371–380, May 2017.
- [26] Julien Herrmann, Yusuf Özkaya, Bora Uçar, Kamer Kaya, and Umit Çatalyürek. Multilevel algorithms for acyclic partitioning of directed acyclic graphs. SIAM Journal on Scientific Computing, 41(4):A2117–A2145, 2019.
- [27] Tobias Heuer, Peter Sanders, and Sebastian Schlag. Network flow-based refinement for multilevel hypergraph partitioning. ACM Journal of Experimental Algorithmics (JEA).
- [28] Tobias Heuer and Sebastian Schlag. Improving coarsening schemes for hypergraph partitioning by exploiting community structure. In 16th International Symposium on Experimental Algorithms (SEA 2017), London, UK, 21th - 23rd June 2017. Ed.: C. Iliopoulos, volume 75 of LIPIcs - Leibniz International Proceedings in Informatics, page Art. Nr. 21. Schloss Dagstuhl - Leibniz-Zentrum für Informatik, Wadern, 2017. 46.12.02; LK 01.
- [29] Manuel Holtgrewe, Peter Sanders, and Christian Schulz. Engineering a Scalable High Quality Graph Partitioner. Proceedings of the 24th IEEE International Parallal and Distributed Processing Symposium, pages 1–12, 2010.
- [30] Laurent Hyafil and Ronald Rivest. Graph partitioning and constructing optimal decision trees are polynomial complete problems. IRIA-Laboria, Rocquencourt, France, 1973.
- [31] Arthur B. Kahn. Topological sorting of large networks. Commun. ACM, 5(11):558–562, November 1962.
- [32] George Karypis, Rajat Aggarwal, Vipin Kumar, and Shashi Shekhar. Multilevel hypergraph partitioning: applications in vlsi domain. *IEEE Transactions on Very Large Scale Integration (VLSI) Systems*, 7(1):69–79, March 1999.

- [33] George Karypis and Vipin Kumar. A fast and high quality multilevel scheme for partitioning irregular graphs. SIAM J. Sci. Comput., 20(1):359–392, December 1998.
- [34] George Karypis and Vipin Kumar. Multilevel k-way partitioning scheme for irregular graphs. J. Parallel Distrib. Comput., 48(1):96–129, January 1998.
- [35] George Karypis and Vipin Kumar. Multilevelk-way partitioning scheme for irregular graphs. J. Parallel Distrib. Comput., 48(1):96–129, January 1998.
- [36] Hye-Jin Kim and Yong-Hyuk Kim. Recent progress on graph partitioning problems using evolutionary computation. CoRR, abs/1805.01623, 2018.
- [37] Fatih Kocan and Mehmet H. Gunes. Acyclic circuit partitioning for path delay fault emulation. In *The 3rd ACS/IEEE International Conference on Computer Systems and Applications, 2005.*, page 22, Jan 2005.
- [38] Lingda Li, Robel Geda, Ari B. Hayes, Yanhao Chen, Pranav Chaudhari, Eddy Z. Zhang, and Mario Szegedy. A simple yet effective balanced edge partition model for parallel computing. *SIGMETRICS Perform. Eval. Rev.*, 45(1):6, June 2017.
- [39] Henning Meyerhenke, Martin Nollenburg, and Christian Schulz. Drawing large graphs by multilevel maxent-stress optimization. *IEEE transactions on* visualization and computer graphics, 24(5):1814–1827, 2018.
- [40] Henning Meyerhenke, Peter Sanders, and Christian Schulz. Partitioning complex networks via size-constrained clustering. In *Experimental Algorithms*, pages 351–363, Cham, 2014. Springer International Publishing.
- [41] Brad L. Miller and David E. Goldberg. Genetic algorithms, tournament selection, and the effects of noise. *Complex Systems*, 9, 1995.
- [42] Orlando Moreira, Merten Popp, and Christian Schulz. Graph partitioning with acyclicity constraints. In 16th International Symposium on Experimental Algorithms, SEA 2017, June 21-23, 2017, London, UK, pages 30:1–30:15, 2017.
- [43] Orlando Moreira, Merten Popp, and Christian Schulz. Evolutionary multi-level acyclic graph partitioning. In Proceedings of the Genetic and Evolutionary Computation Conference, GECCO 2018, Kyoto, Japan, July 15-19, 2018, pages 332–339, 2018.
- [44] Pablo Moscato. On evolution, search, optimization, genetic algorithms and martial arts - towards memetic algorithms. *Caltech Concurrent Computation Program*, 10 2000.
- [45] Pablo Moscato and Carlos Cotta. A Modern Introduction to Memetic Algorithms, pages 141–183. Springer US, 2010.
- [46] Pablo Moscato and Luke Mathieson. Memetic algorithms for business analytics

and data science: A brief survey. Moscato P., de Vries N. (eds) Business and Consumer Analytics: New Ideas. Springer, Cham, pages 545–608, 05 2019.

- [47] Vitaly Osipov and Peter Sanders. n-Level Graph Partitioning. In Proceedings of the 18th European Conference on Algorithms: Part I, volume 6346 of LNCS, pages 278–289. Springer, 2010.
- [48] Sylvain Paris, Samuel W Hasinoff, and Jan Kautz. Local laplacian filters: Edge-aware image processing with a laplacian pyramid. *ACM Trans. Graph.*, 30(4):68, 2011.
- [49] David J. Pearce and Paul H. J. Kelly. A dynamic topological sort algorithm for directed acyclic graphs. ACM Journal of Experimental Algorithms, pages 1–7, 2006.
- [50] François Pellegrini and Jean Roman. Scotch: A software package for static mapping by dual recursive bipartitioning of process and architecture graphs. In *High-Performance Computing and Networking*, pages 493–498, Berlin, Heidelberg, 1996. Springer Berlin Heidelberg.
- [51] Louis-Noël Pouchet. Polybench/c: The polyhedral benchmark suite. https: //web.cse.ohio-state.edu/~pouchet.2/software/polybench/, 2012.
- [52] Peter Sanders and Christian Schulz. Engineering Multilevel Graph Partitioning Algorithms. In Proceedings of the 19th European Symposium on Algorithms, volume 6942 of LNCS, pages 469–480. Springer, 2011.
- [53] Peter Sanders and Christian Schulz. Distributed evolutionary graph partitioning. In *Proceedings of the Meeting on Algorithm Engineering and Experiments*, ALENEX '12, page 16–29, USA, 2012. Society for Industrial and Applied Mathematics.
- [54] Sebastian Schlag. High-Quality Hypergraph Partitioning. PhD thesis, 2019.
- [55] Sebastian Schlag, Vitali Henne, Tobias Heuer, Henning Meyerhenke, Peter Sanders, and Christian Schulz. k-way hypergraph partitioning via n-level recursive bisection. In 2016 Proceedings of the Eighteenth Workshop on Algorithm Engineering and Experiments (ALENEX), pages 53–67.
- [56] Sebastian Schlag, Matthias Schmitt, and Christian Schulz. Faster support vector machines. In 2019 Proceedings of the Twenty-First Workshop on Algorithm Engineering and Experiments (ALENEX), pages 199–210.
- [57] Sebastian Schlag, Christian Schulz, Daniel Seemaier, and Darren Strash. Scalable edge partitioning. In 2019 Proceedings of the Twenty-First Workshop on Algorithm Engineering and Experiments (ALENEX), pages 211–225.
- [58] Christian Schulz and Darren Strash. Graph partitioning: Formulations and applications to big data. In *Encyclopedia of Big Data Technologies*, pages 1–7. Springer, 2018.

- [59] Aleksandar Trifunovic and William J. Knottenbelt. Parallel Multilevel Algorithms for Hypergraph Partitioning. *Journal of Parallel and Distributed Computing*, 68(5):563 – 581, 2008.
- [60] Jesper Larsson Träff. Direct graph k-partitioning with a kernighan-lin like heuristic. *Operations Research Letters*, 34(6):621 629, 2006.
- [61] Chris Walshaw and Marg Cross. Mesh partitioning: A multilevel balancing and refinement algorithm. *SIAM J. Sci. Comput.*, 22(1):63–80, January 2000.

A Detailed DAG Results

Table A.1: Detailed per instance results on the ISPD98 benchmark set [3]. HOUKC refers to the algorithm developed by Herrmann et. al. [25]. mlDHGP + X refers to our multilevel algorithm and memDHGP + X refers to our memetic algorithm with X as undirected hypergraph partitioner for initial partitioning. The *Best* column reports the best edge cut found during 8 hours of individual runs. For mlDHGP + X, the *Average* column reports the average edge cut of 5 individual runs and the *Best* column reports the best edge cut found during 8 hours. For memDHGP + X, the *Best* column reports the best edge cut found during 8 hours. For memDHGP + X, the *Best* column reports the best edge cut found during 8 hours. For memDHGP + X, the *Best* column reports the best result found after running for 8 hours. The *Overall Best* column shows the best cut found by any tool with the following identifiers: H: HOUKC, N: one of the new approaches. In general, lower is better.

		HO	UKC	mlD	HGP 1	memDHGP	mlDHGP	memDHGP	Over	all
				1	with KaHy	Par	with	PaToH	Bes	t
Graph	Κ	Average	Best (8h)	Average	Best (8h)	Best (8h)	Average	Best (8h)	Result S	olver
	2	3175	2 752	3 2 3 5	2428	2 2 5 5	2 7 3 0	2 2 9 0	2255	N
	4	6 0 9 2	5099	5 4 3 4	5028	4848	5325	4841	4 8 4 1	N
ibm01	8	7449	6880	8 0 2 6	7240	6958	8 268	6 6 3 9	6 6 3 9	N
	16	10555	8 603	9 1 3 1	8 1 3 5	8 0 2 8	8 870	7627	7 6 2 7	N
	32	12652	11 1 19	10 909	10086	9572	11 107	9404	9 4 0 4	N
	2	8 5 4 0	4 708	8 772	3262	5873	8 806	8 599	3262	Ν
	4	13264	11375	12 290	11374	11497	12317	11400	11374	N
ibm02	8	17832	16591	17 557	16522	16253	17520	16387	16253	Ν
	16	24856	23002	21 708	20209	19727	22 1 28	20455	19727	N
	32	30407	29082	26 379	25263	24264	26659	25393	24264	Ν
	2	14601	13 687	15278	12584	11870	14265	12 051	11870	N
	4	21802	20077	20652	18622	17757	18 840	17835	17757	N
ibm03	8	26051	24361	25 370	21494	20579	22975	20699	20579	N
	16	30776	27238	29 885	24637	24006	28097	23837		N
	32	33439	31034	32 134	27309	27093	30035	27085	27085	N
	2	9 5 1 8	9 1 0 8	9 7 2 7	8508	8 2 3 7	9 7 2 7	8 2 3 7	8 2 3 7	Ν
	4	14226	13190	12 668	11512	10970	12358	10944	10944	N
ibm04	8	18508	16683	18 677	16983	16298	18 811	15878	15878	N
	16	25885	22874	24 363	22800	21812	24298	21373	21373	Ν
	32	30512	27107	27 882	26486	25078	28127		25078	N
	2	8 360	5882	7 4 9 4	5478	5830	7 285	6979	5478	N
	4	17040	13278	14 932	10740	10710	15035	11885	10710	N
ibm05	8	23170	19480	19618	16076	15980	19803	15934	15934	Ν
	16	29747	25590	25 512	22049	20771	24914	21604		N
	32	34495	30721	29 437	27465	27582	30155	26899		N
	2	14049	12736		11804	11341	12832	11285		Ν
	4	23206	20317	21 641	19097	18197	21 705	18374	18197	Ν
ibm06	8	30875	26980	25 402	23202	22455	25155	22263	22263	Ν
	16	34069	30848	29 421	27435	26384	29 793	27263		N
	32	38243	36197	32 781	31310	30839	32 826	30597	30597	Ν
	2	16341	15855	15 738	15356	13681	16 003	12965	12965	Ν
	4	26842	23522	22 608	21583	20499	22 273	20348		N
ibm07	8	29 702	27069	26 935	25655	24464	27 186	24586		Ν
	16	36633	33606	31 746	30788	29808	32195	29797		N
	32	43083	40205	36 959	35901	34648	37 017		34648	Ν
	2	25139	24481	24 418	22381	22079	24 384	21925	21925	Ν
	4	52118	38711	41 350	38644	38495	41 402	38330		N
ibm08	8	84 639	81587	50 063	49238	48429	50 043	47124		Ν
	16	96107	88135	88 727	87323	85996	89 513		85996	Ν
	32	109264	96746	93 556	92591	90 779	94 172	90 660		Ν
	2	19509	15084	17 233	12661	12305	16 307	12127		Ν
	4	28408	25120	26 143	23342	22557	26 184	20 892		Ν
ibm09	8	36168	31734	33 276	30411	29654	34 341		29654	Ν
	16	42373	39154	39 712	37301	35902	39 529	34707	34707	Ν
	32	50041	45987	45 226	41007	40701	45131	39887	39887	N

Table A.2: Detailed per instance results on the ISPD98 benchmark set [3]. HOUKC refers to the algorithm developed by Herrmann et. al. [25]. mlDHGP + X refers to our multilevel algorithm and memDHGP + X refers to our memetic algorithm with X as undirected hypergraph partitioner for initial partitioning. The *Best* column reports the best edge cut found during 8 hours of individual runs. For mlDHGP + X, the *Average* column reports the average edge cut of 5 individual runs and the *Best* column reports the best edge cut found during 8 hours. For memDHGP + X, the *Best* column reports the best edge cut found during 8 hours. For memDHGP + X, the *Best* column reports the best edge cut found during 8 hours. For memDHGP + X, the *Best* column reports the best result found after running for 8 hours. The *Overall Best* column shows the best cut found by any tool with the following identifiers: H: HOUKC, N: one of the new approaches. In general, lower is better.

		HOI	UKC		HGP vith KaHy			memDHGP PaToH	Overa Bes	
Graph	К	Average	Best (8h)			Best (8h)		Best (8h)	Result S	
	2	24 983	24073		21 575	21 328	22 560	21 310	21 310	
	4	38 620	35 083	39 383	33 217	36 352	39 288	32 101	32 101	
bm10		49646	44820	47827	40 423	39 202	46 082	38 238	38 2 38	
011110	16	63 960	54164	55 610	50 854	49 632	56 1 2 9	49 892	49632	
	32	69 990	65 302	64229	61 838	59 914	64 105	59 180	59180	
	2	19 224	16 926	21 879	14374	13 578	15748	13 318	13 318	
	4	36 346	26 5 39	26 9 1 9	22750	21623	24724	21 310	21310	
bm11	8	39755	32812	32816	30401	28563	33 247	28477	28477	
	16	52698	45779	40 706	38055	39 294	43 773	37 257	37257	
	32	63925	57699	50612	47999	47 331	52963	47 930	47331	
	2	29 359	27 238	30 315	27 860	27365	29 6 20	27688	27238	
	4	50457	47 922	49225	44 108	42728	49591	46 107	42728	
bm12	8	60 024	53 785	57394	52487	51425	57046	50 955	50955	
	16	72429	65979	66 486	62965	61 186	67 160	61 484	61 1 86	
	32	84 328	76 066	73872	70 503	68 7 39	73 252	68712	68712	
	2	30 698	19 008	21 700	17161	17484	22 151	17659	17161	
	4	39781	29 1 98	39 288	31 700	32 060	38 609	26 500	26500	
bm13	8	54061	39 453	55253	42 881	44535	47765	41 596	39453	
	16	71208	60 006	65263	55070	49820	65 962	49 993	49820	
	32	89 053	76762	81 831	72262	74997	81 416	70 987	70 987	
	2	33 205	31 988	51 511	48 338	48 140	52 065	49 670	31 988	
	4	55342	49 972	69 320	64 838	62 888	70 364	66 680	49972	
bm14		76 297	68 992	68 051	62718	60 929	67 598	56 972	56972	
	16	96 638	80 591	79801	74705	73 224	80 0 29	73 861	73224	
		104543	96 677	91 692	89 688	87 904	92 823	86 504	86 504	
	2	74713	71 593	66 301	63 603	63 1 36	82679	67 804	63 1 36	
	4	105577	95 911	97786	87 849	92 812	96479	88 349		
bm15	8	146984	123 993		112014	113564	124 884	108 619		
	16		153693		135061	124709	143941	133614		
	32		174057	158765	154660	149 558	160 815	148763		
	2	55 871	52 980	51 699	48 222	48 063	50 167	45 371	45 371	
	4	108576	93 874	98471	93 941	91481	99729	89 976	89976	
bm16		130302	117 375		115437	119 439	126431	114458		
	16		148626		136 916	134387	142235	132412		
		181 924	172909		158854	157879	164 966	153490		
	2	75 860	57 177	70 331	59100	59 470	61 401	56 895	56 895	
	4	100287	89849	121023	78692	77 889	121175	107211	77889	
bm17	8	151126	141679	152455	124639	126610	147 848	130 307		
	16	182272	166847	171507	153 812	155789	165498	150 026		
	32		198404	188792	167274	173762	194056	182853		
	2	37 123	34 949	35 907	33 434	33 394	36 651	33 277	33 277	
	4	63 000	53 948	64540	53190	53 237	58432	48 482	48 482	
bm18	8	92 636	78 164	86 580	76686	75 728	81 435	70 558	70 558	
	-	121219	108 744		93 018	88 959	113 181	98 976		
		144219	132 289			110 816	128 875	119 170		
Mean	_	41189	36205	37 828	33459	33 007	37 382	33 088		_

Table A.3: Detailed per instance results on the PolyBench benchmark set [51]. HOUKC refers to the algorithm developed by Herrmann et.al. [25]. Moreira refers to the algorithm developed by Moreira et al. [43]. mlDHGP + X refers to our multilevel algorithm and memDHGP + X refers to our memetic algorithm with X as undirected hypergraph partitioner for initial partitioning. For HOUKC, the Average column reports the better average from Table A.1 and Table A.2 in [25] and the Best column reports the best edge cut found during 8 hours of individual runs or the best edge cut reported in [25], if that is lower (marked with a star). For mlDHGP + X, the Average column reports the average edge cut of 5 individual runs and the Best column reports the best edge cut found during 8 hours. For memDHGP + X, the Best column reports the best result found after running for 8 hours. The Overall Best column shows the best cut found by any tool with the following identifiers: H: HOUKC, N: one of the new approaches, M: Moreira et.al. In general, lower is better.

			HOUKC	Moreira	a et. al.	mlD	HGP	memDHGP	mlDHGP	memDHGP	Ove	rall
							with KaHy			PaToH	Be	
Graph	Κ	Average	Best (8h) or [25]	Average	Best		Best (8h)	Best $(8h)$	Average	Best $(8h)$	Result	
	2	200	200	200	200	200	200	200	200	200		H,M,N
	4	2 160	946	947	930	1065	930	930	1 006	930	930	
2mm	8	5 361	2 910	7 181	6 604	2819	2576	2465	5 563	5 1 1 0	2 465	N
	16 32	11 196 15 911	8 103 12 708	13330 14583	13092 14321	7 090 11 397	5963 10 635	$5435\10398$	7881 12228	6 632 11 012	5 435 10 398	N N
	2	1000	800	14 383	14321	800	800	800	12 228	11012	800	H,N
	4	9 264	2 600	38 722	37 899	2 647	2 600	2 600	2 600	2 600	2 600	
3mm	8	24 330	7 735	58 129	49 559	8 596	6 967	6 861	14 871	9 560	6 861	N
0	16	37 041	21 903	64 384	60 127	23 513	19 625	19675	28 021	23 967		N
	32	46437	36718	62 279	58190	34 721	30908	31423	38879	34353		Ν
	2	142719	*134 675	134945	134675	138433	138057	138 279	138520	138 329	134675	H,M
	4	212938	210979	284666	283892	213255	212709	212851	213390	212564	210 979	H
adi	8	256302	229563	290823	290672	253885	252271	253206	254282		229563	Н
	16	282485	271374	326963	326923	281068	277337	280437	281751		271374	Н
	32	306075	305 091	370 876	370413	309 930	303078	299 387	309 757		302 157	N
	2	39876	32 451	47 826	47 424	39 695	24 150	23 690	45 130	43 450		N
	4	48 645	43 511	82 397	76 245	50 725	42 028	39 316	50 144	47 486		N
atax	8	51 243	48 702			54 891	48 824	47 741	52 163	49 450		N
	$\frac{16}{32}$	59 208 69 556	52 127 57 930	127687 132092	$125146 \\ 130854$	68 153 66 267	50 962 54 613	51256 56051	53256 56 773	51 191 54 536	51 191 54 536	N N
	32 2		4775									
	2 4	27 269 61 991	*34 307	66 520 84 626	66445 84213	4 775 12 362	4 775 11 724	477511281	5893 13339	5641 12344	4 775 11 281	H,N N
covariance		74 325	*50 680		102425	24 429	21460	21106	51 984	41 807		
covariance	16	119 284	99 629	125 816	102425 123276	62 011	60 1 4 3	58 875	65 302	59 153		N
	32	121 155	94 247	142 214	137 905	76 977	73758	72 090	80 464	74 770	72 090	N
	2	5 0 3 5	3 000	43 807	42 208	3 0 0 0	3 0 0 0	3 000	3 000	3 000	3 000	H,N
	4	37051	9 0 0 0	72115	71082	11029	9 000	9 0 0 0	28317	27852	9 0 0 0	H,N
doitgen	8	51283	36790	76977	75114	36 326	34912	34682	42185	38491	34 682	Ń
	16	62 296	50481	84 203	77436	51064	48992	50486	50993	48193	48 193	Ν
	32	68350	59632	94135	92739	59159	58184	57408	57208	55721	55 721	Ν
	2	12997	12997	12997	12997	12997	12997	12997	12997	12997		$_{\rm H,M,N}$
	4	21566	*21 566	21641	21641	21556	21557	21541	21556	21541	21 541	Ν
durbin	8	27519	27518	27571	27571	27511	27508	27509	27511	27509	27 509	Ν
	16	32 852	32 841	32 865	32 865	32 869	32 824	32 825	32 852	32 825	32 825	N
	32	39 738	39 732	39 726	39 725	39 753	39 717	39701	39 753	39 701	39 701	N H
	2 4	6 024 15 294	4 381 11 551	5 494 15 100	5494 15099	5233 11670	$4756 \\ 9325$	4 604 9 240	6318 11572	6 285 10 232	4 381 9 240	H N
fdtd-2d	8	23 699	19 527	33 087	32355	17 704	9325 15906	9240 15653	17 990	10232 15758		N
lata-2a	16	32 917	28 065	35 714	35 239	25 170	22866	22 041	24 582	22 003		N
	32	42 515	39 063	43 961	42507	32 658	30 872	29 868	32 682	29 772	29 772	N
	2	4 200	4 200	383 084	382 433	4 200	4 200	4 200	4 768	4 6 9 0	4 200	H,N
	4	59854	12 600	507250	500526	12 600	12 600	12 600	13 300	12 600		
gemm	8	116990	33 382	578951	575004	70 827	31413	30 912	188172	175495	30 912	Ń
-	16	263050	224173	615342	613373	185872	164235	148040	202920	194017	148 040	Ν
	32	330 937	277879	626472	623271	270 909	265771	258607	280849	275188	258607	Ν
	2	20 913	*20 913	29349	29270	22725	19485	19 390	20 317	18 930	18 930	Ν
	4	40299	35431	49361	49229	38 600	35021	33324	37632	34328	33 324	Ν
gemver	8	55266	43716	68163	67094	50440	44253	43276	47799	42548		N
	16	59072	54012	78 115	75596	53 819	48618	48182	53775	46563		N
	32	73 131	63 012	85 331	84 865	58 898	53 581	54 953	59 210	52 404		N
	2	500	500	1 666	500	500	500	500	500	500		H,M,N
	4	10 316	1 500	98 542	94 493	5 096	1 500	1 500	1 548	1 500		N
gesummv	8 16	9618	4 0 2 1	101 533	98 982	25 535	3 500	3 500	3 640	3 500	3 500	N N
	16 32	$35686 \\ 45050$	11 388 28 295	112064 117752	104866 114812	30 215 31 740	7 500 15 620	7 500 16 339	7883 16 144	7500 15 500	7 500 15 500	IN N
	32 2	9378	8 936	8 6 9 5	8 684	8 930	8 640	8 640	9 2 4 2	8 936	8 6 4 0	N
	2 4	16 700	8 936 15 755	14 592	8 684 14 592	15 355	8 640 14 642	14592	9 242 16 304	14865	14 592	
heat-3d	8	25 883	24 326	20 608	20 608	23 307	14042 21190	21 300	25462	23 074	20 608	MI,IN
1000-00	16	42 137	*41 261	31 615	31 500	38 909	$\frac{21}{38053}$	35 909	40 148	37 659	31 500	M
	32	64 614	60 215	51 963	50 758	55 360	53525	51682	54 621	50 848		
				02000		00000	00.020	01 001		00010		

Table A.4: Detailed per instance results on the PolyBench benchmark set [51]. HOUKC refers to the algorithm developed by Herrmann et.al. [25]. Moreira refers to the algorithm developed by Moreira et al. [43]. mlDHGP + X refers to our multilevel algorithm and memDHGP + X refers to our memetic algorithm with X as undirected hypergraph partitioner for initial partitioning. For HOUKC, the Average column reports the better average from Table A.1 and Table A.2 in [25] and the Best column reports the best edge cut found during 8 hours of individual runs or the best edge cut reported in [25], if that is lower (marked with a star). For mlDHGP + X, the Average column reports the average edge cut of 5 individual runs and the Best column reports the best edge cut found during 8 hours. For memDHGP + X, the Best column reports the best result found after running for 8 hours. The Overall Best column shows the best cut found by any tool with the following identifiers: H: HOUKC, N: one of the new approaches, M: Moreira et.al. In general, lower is better.

		H	IOUKC	Moreira	et. al.			memDHGP			Overall	
							vith KaHy			PaToH	Bes	
Graph				Average	Best		Best (8h)	Best $(8h)$	Average	Best $(8h)$	Result	
	2	646	400	596	596	440	400	400	491	423	400	H,N
	4	1617	1 1 2 3	1 493	1 4 9 2	1188	1 0 4 6	1044	1 250	1 1 28	1044	N
jacobi-1d		2845	2 0 5 2	3 1 3 6	3136	2 0 2 8	1 754	1 750	2 170	1 855	1 750	N
	16	4519	3 5 1 7	6 3 4 0	6 3 38	3140	2912	2869	3 355	2 982	2869	N N
	32	6742	5 545 *3 342	8 923 2 994	8 750 2 991	4 776 3 878	4 565	4 498 2 986	4 910	4 587	4 498	N
	2 4	$\frac{3445}{7370}$	7 243	2 994 5 701	5 700	7 591	$3000 \\ 5979$	2 986 5 881	3 942 7 528	$3129 \\ 6245$	$2986 \\ 5700$	M
jacobi-2d		13168	12 134	9417	9 4 1 6	10 872	9 2 9 5	8 935	11 753	10 492	8 935	N
Jacobi-20	16	21565	18 394	16 274	16 231	15 605	14746	13 867	15 889	14 736	13 867	N
	32	29558	25 740	22 181	21 758	20 597	19647	18 979	21 653	19 530	18 979	N
	2	5 3 5 1	4 160	5 210	5 1 6 2	4160	4 160	4 160	4 1 6 0	4 160	4 160	H,N
	4	21258	12 214	13 5 28	13 510	12 720	12214	12 214	16 091	15 992	12 214	H,N
lu	8	53643	34074		33 211	42 963	33873	33954	41 1 13	38 318	33 211	M
	16	105289	81713	74 543	74006	81 224	74400	74448	83 980	75150	74 006	Μ
	32	156187	141 868	130674	129954	125 932	122977	121451	131850	127904	121451	Ν
	2	5731	5 337	5 380	5 337	5 3 37	5 337	5 337	5 3 37	5 3 37	5 3 37	H,N
	4	22368	15170	14744	14744	18 114	16889	17560	26 606	17113	14744	Ν
ludcmp	8	60255	41086	37 228	37069	46 268	37688	37790	52980	39362	37069	Ν
	16	106223	86959		78467	89 958	76074	80706	96275	85572	78467	N
	32	158619	144 224	134758			125957	127454	136 218	131 161	127454	Ν
	2	21281	16768	24 528	23091	23 798	16584	16596	32856	20016	16596	Ν
	4	38 215	29 229	74 386	73 035	41 156	29 318	30 070	52 353	42 870	29 229	H,N
mvt	8	46776	39 295	86 525	82 221	50 853	36 531	35 471	60 021	55 460	35 471	N
	16	54925	48 036		97 941	58 258	41 727	42 890	65 738	59 194	42 890	N
	32	62 584	54 293	105 066		58 413	45 958	46 122	69 221	64 611	46 122	N H,N
	2 4	$4374 \\ 11784$	3 401 10 872	$4991 \\ 12197$	4969 12169	4 036 11 352	$3578 \\ 10645$	3 504 10 404	4 206 11 480	$3786 \\ 10604$	$3401 \\ 10404$	H,N N
seidel-2d		21937	20711	21 4 197	12109 21 400	19954	10645 18528	10 404	20 309	18 482	10404 17770	N
seidei-2d	16	38065	33 647	38 222	38 110	29 930	27644	27 583	30 329	28 348	27 583	N
	32	58319	51 745	52 246	51 531	41 256	38 949	38 175	42 291	39 058	38 175	N
	2	26374	21 963	94357	94 214	22 000	21 840	21 836	29871	26 134	21 836	N
	4	59815	42 442			41 486	38 290	37 854	65 1 1 1	57 620	37 854	N
symm	8	91892	69554	152984		69 569	58084	60644	82 865	75151	60644	Ν
Ŭ	16	105418	89 320	167 822	167512	90 978	83 703	85508	96 932	89445	85 508	Ν
	32	108950	97174	174938	174843	110 495	104376	100337	108 814	104592	97174	Н
	2	4343	900	11 0 98	3894	900	900	900	900	900	900	H,N
	4	12192	3048	49662	48021	3 1 5 0	2978	2909	16589	9 9 9 1	2 909	Ν
syr2k	8	28787	12833	57584	57408	12 504	9 969	10154	21427	19507	10154	Ν
	16	29519	24457	59 780	59594	25 054	21626	21828	26 1 20	23588	21 828	Ν
	32	36 111	31 1 38	60 502	60 085	33 424	31 236	29 984	31 358	29 340	29 340	N
	2	11740	3 240	219 263		3 240	3 240	3 240	3 4 3 9	3 240	3 240	H,N
	4	56832	9 960	289 509		10 417	10 119	9970	89 457	80 801	9 960	H
syrk	8 16	$112236 \\ 179042$	30 602 147 058	329466 354223		83 000 117 357	46130 113 122	58 876 111 635	107 220 150 363	101516 135615	30602 111 635	H N
	32	179042 196173	173 550	362 016		158 590	113 122 154 818	154 921	182 222		111035 154921	N
	32 2	336	280	6788	3 5 4 9	280	279	279	308	279	279	N
	4	828	280 827	43 927	43549	823	821	821	865	823	821	N
trisolv	8	2156	1 907	66 1 48	65 662	2 112	1 893	1 895	2 0 3 5	1 897	1 895	N
	16	6 2 4 0	5 285	71 838	71 447	8 719	4 1 2 5	4 108	4 358	4 240	4 108	N
	32	13431	*13172	79125	79071	16 0 27	8 9 4 2	8784	9 2 1 0	8716	8716	Ν
	2	13659	3 4 4 0	138937		3 4 4 0	3440	3 4 4 0	3440	3 4 4 0	3 4 4 0	H,N
	4	58477	14543	192752	191492	14942	12622	12389	35964	35824	12389	Ń
trmm	8	92185	49 830	225192	223529	65 303	46059	45053	67011	61045	45053	Ν
	16	128838	103975	240788		92172	79507	80 186	96421	87275	80 186	Ν
	32	153644	131 899	246407	245173	120 839	115460	112 267	120753	113205	112267	N
Mean		25777	17897	44 923	43 200	18 887	15988	16 095	20 308	18 6 4 2		
		· · · ·									1	

B Detailed DAH Results

Table B.1: Detailed per instance results on the ISPD98 benchmark suite [3]. mlDHGP refers to our algorithm with KaHyPar as undirected hypergraph partitioner for initial partitioning. memDHGP refers to our memetic algorithm that uses mlDHGP equipped with KaHyPar as undirected hypergraph partitioner for initial partitioning to build an initial population. The *Best* column reports the best edge cut found during 8 hours of individual runs. For mlDHGP, the *Average* column reports the average edge cut of 5 individual runs. For memDHGP, the *Best* column reports the best result found after running for 8 hours. In general, lower is better.

							erPartKWa
Hypergra		Average			Best (8h)		Best (8h
	2	838	629	877	659	1267	66
	4	1835	1427		1684	4921	261
ibm01	8	2923	2136	3512	2670	6513	4 1 5
	16	3764	3049	4584	3710	8271	6 0 3
	32	4774	4013	5 6 2 6	4 706	9894	6 65
	2	2222	1869	2 6 2 9	1990	3048	2 31
	4	4391	3247		4017	7520	518
ibm02	8	6898	5674		6677		948
	16	9787	8 4 8 1	10678			1270
	32	12545	11448				14 59
	2	3 782	2242	3772	2862	4306	2 93
	4	5955	4 2 3 1	6335	4748	8661	674
ibm03	8	7679	5911	8478	6771	12510	1013
	16	9179	7386				1230
	32	10051	8 4 9 6	12271	10 116		1416
	2	3 080	717	4448	3044	5252	3 20
	4	5232	2467	6175	3 707		6 0 8
ibm04	8	7239	5339	9919	7304		9 9 1
	16	9415	7343	11868	10029		1258
	32	11129	9 259	13795	11947	21342	16 27
	2	4630	3954	4799	4 2 3 2	4952	424
	4	7629	5 930	9574			7 93
ibm05	8	10434	8612	13292			11 82
	16	13095	11285	16394	13566	21884	16 61
	32	15371	13837		15938		20 07
	2	4486	2 7 3 0	5624		7027	4 27
	4	8 1 8 9	5858				1197
ibm06	8	10203	8 2 8 1	11483	9590		1501
	16	12720	10157	14123			2059
	32	15155	12179	17588	14851	30019	24 37
	2	5355	3 680	8 2 7 3	3871	8 8 3 1	4 60
	4	10343	6250		7130		1169
ibm07	8	12386	8 993	13861	9482		1686
	16	13927	11870	17289	14109		23 06
	32	16880	14264	20418			26 76
	2	12865	9344		9331	12536	11 28
	4	18373	16860		15454	22071	2050
ibm08	8	22238	20526		20013		2483
	16	25572	23100		25072	34302	30 63
	32	29667	27425	31907	29150		34 42
	2	5593	3 357	11804	9159		9 65
	4	10610	6416		13852		1769
ibm09	8	12053	8 7 2 6		16942		2110
	16	14987	11588		20090		2884
	32	17802	14449	25313	22858	39478	3485

Table B.2: Detailed per instance results on the ISPD98 benchmark suite [3]. mlDHGP refers to our algorithm with KaHyPar as undirected hypergraph partitioner for initial partitioning. memDHGP refers to our memetic algorithm that uses mlDHGP equipped with KaHyPar as undirected hypergraph partitioner for initial partitioning to build an initial population. The *Best* column reports the best edge cut found during 8 hours of individual runs. For mlDHGP, the *Average* column reports the average edge cut of 5 individual runs. For memDHGP, the *Best* column reports the best result found after running for 8 hours. In general, lower is better.

			nemDHGP				
Hypergraph			Best $(8h)$				
	2		8 288			17445	11 3'
	4		11809				182
ibm10	8	22 141	16401	22146	18837		286
	16		20199				392
	32	30 100	26238	33799			513
	2	10 669	6550			13320	81
	4	16257	10447	18761	15073	25457	208
ibm11	8	17992	12341	23071	18430	36747	296
	16	20 197	16198	27404	23673	42915	359
	32	23409	19767	30592	27495	50761	436
	2	15449	11349	14881	10588	14858	127
	4	20 307	15652	20215	16538	23398	188
ibm12	8	23036	18 126	24916	21053	36501	314
	16	28437	23367	30176	26434	48537	351
	32	34536	27911	38183	33930	62718	537
	2	11 893	8 6 9 5	12790	8 2 8 4		105
	4	14791	10285	21166			344
ibm13	8	21405	14330	32543	21452	58298	483
	16	25313	16761	35524			588
	32	29676	26017			92014	813
	2	24 379	14713				183
	4	30 912	21613				333
ibm14	8	36 370	30710				447
	16	42321	35598				581
	32	48 741	43979				720
	2	27 810	19804				292
	4	44 069	33 151				745
ibm15	8	51 886	38 306	65971		102738	934
	16	58 961	49 687	74815			105 9
	32	66 287	56 374	82 252		141076	129 2
	2	25 941	11 494				30.8
	4	46 933	36 124				619
ibm16	8	57 761	45 328	61 1 32			890
	16	67 904	58 471	72 549		114337	1051
	32	80 591	69 325	83 631		135371	127 0
	2	36 934	32 507				296
	4	47 186	39 301				539
ibm17	8	62 896	54 230	61 043			784
omri	16	74 427	65 672	77 867		109982	1017
	32	86 597	80 152	93 864		109302 129758	121 2
	2	21 296	16 338	20 830			1212
	4	36 235	28 131	33642			390
ibm18	4 8	49 742	38947	45 337			590 659
101110	16		38947 47532	60.672	55 461		847
	32	67 770	47532 58061	$60672 \\ 76324$	69 222	94165 114532	847 1040
λ	02						
Mean		16 151	12245	18344	15030	26839	212

Table B.3: Detailed per instance results on the PolyBench benchmark set [51]. mlDHGP refers to our algorithm with KaHyPar as undirected hypergraph partitioner for initial partitioning. memDHGP refers to our memetic algorithm that uses mlDHGP equipped with KaHyPar as undirected hypergraph partitioner for initial partitioning to build an initial population. The *Best* column reports the best edge cut found during 8 hours of individual runs. For mlDHGP, the *Average* column reports the average edge cut of 5 individual runs. For memDHGP, the *Best* column reports the best result found after running for 8 hours. In general, lower is better.

							erPartKWa
Hypergraph		Average			Best $(8h)$		Best (8h
	2	212	200	224	200	344	21
_	4	633	608	905	840	1618	75
2mm	8	1376	1320	1 608	1440	3169	1 43
	16	2 2 3 9	2 1 5 3	2 6 9 5	2 248	4 6 9 1	263
	32	3 796	3624	4 562	3 934	7 015	4 22
	2	800	800	1112	805	1 564	1 09
0	4	2 419	2 000	3 1 5 5	2 480	5 036	3 56
3mm	8	3 950	3 540	5 940	4 689	9374	5 65
	$\frac{16}{32}$	$6264 \\ 9234$	$5999 \\ 8861$	9 099 12 719	7537 11 483	12996 19224	8 12 12 51
	32 2	9 234	460	460	460	19 224	5.82
	4	9200 9 4 3 8	4 9 4 3	7 162	1 719	24248	1946
atax	8	22036	17127	20110	9 2 9 1	24248 27736	20.98
atax	16	30917	28 378	29 675	24167	46 152	29.03
	32	43 936	41 981	40 637	39 098	52265	46 79
	2	2 930	2 590	3 3 4 3	3 160	3 1 9 0	3 05
	4	6 0 5 8	5 705	5 3 6 1	5 265	7 0 2 9	5 68
covariance	8	8 834	8 2 3 8	9 6 6 0	9 0 9 2	12815	10 47
	16	13406	12758	13 917	13 480	19825	16 52
	32	17605	17210	20 211	19833	29596	2464
	2	400	400	3134	2 9 2 7	3444	2 28
	4	1200	1 200	3 6 5 2	3 600	6760	311
doitgen	8	2892	2800	5 301	4613	11254	540
	16	6001	5800	7 263	6949	15725	8 24
	32	9566	9192	11405	11221	20172	1487
	2	349	349	349	349	352	34
	4	1024	1020	1 0 2 3	1020	1033	1 02
durbin	8	2361	2 3 3 9	2 362	2344	2375	2 34
	16	5030	4996	5 0 2 7	5000	5047	5 01
	32	10374	10364	10 366	10 358	10 396	10 37
	2	2 650	1 756	3 4 9 1	3 4 9 0	3 4 9 1	3 49
	4	5549	3 960	10 473	4 294	10 474	4 26
fdtd-2d	8	7 755	6351	13745	8 673	24 366	812
	16	10 971	8 959	19112	13 681	34 855	1510
	32	14 110	12759	24 248	18 726	42 703	22.03
	2	4200 12600	4200 12600	6179 18908	4758 14 781	5989 18579	4 50
	4 8	20 931	12000 19714	39 528	14781 39290	41055	14 58 35 13
gemm	16	33978	31355	63 1 39	63 1 39	77882	76 50
	32	53978 52721	50 300	89 660	89 660	117 319	115 71
	2	2 577	480	1824	480	4 800	2 94
	4	5341	2 070	6705	4 5 7 6	8 081	5 85
gemver	8	10 615	8 305	10 5 2 2	8 3 5 7	15 511	967
80111101	16	13432	12474	13 263	12618	20 260	1400
	32	17250	16 576	16 823	16 362	25050	21 08
	2	350	250	518	501	523	50
gesummv	4	975	750	927	760	1 1 9 1	1 05
	8	1 394	1 250	1 5 3 9	1 515	2128	2 05
	16	2247	2 2 4 6	2 600	2582	5403	29
	32	3526	3 4 2 8	3644	3454	4689	4 29
	2	1 280	1 280	1 3 4 7	1 280	1358	1 28
	4	3843	3840	3 9 4 7	3840	4190	384
heat-3d	8	9427	8 7 7 7	9 2 2 2	8 960	9776	8 96
	16	15406	14509	16 4 96	14325	19799	1431
	32	21102	19382	22 7 27	20483	28957	21 08

Table B.4: Detailed per instance results on the PolyBench benchmark set [51]. mlDHGP refers to our algorithm with KaHyPar as undirected hypergraph partitioner for initial partitioning. memDHGP refers to our memetic algorithm that uses mlDHGP equipped with KaHyPar as undirected hypergraph partitioner for initial partitioning to build an initial population. The *Best* column reports the best edge cut found during 8 hours of individual runs. For mlDHGP, the *Average* column reports the average edge cut of 5 individual runs. For memDHGP, the *Best* column reports the best result found after running for 8 hours. In general, lower is better.

		mIDHGP 1	nemDHGP	TopoOrc	lerPartRB	TopoOrd	erPartKWa
Hypergrap	h K	Average	Best (8h)	Average	Best (8h)	Average	Best (8)
	2	401	400	412	402	411	40
	4	926	793	1245	1 206	1279	1 20
jacobi-1d	8	1587	1467	2900	2814	3053	2.7
	16	2634	2423	6213	3 900	6676	3.3
	32	3992	3 786	8788	5540	13680	47
	2	1008	1008	1053	1008	1049	1 00
	4	3524	2981	3093	3024	3129	30
jacobi-2d	8	5786	4995	7184	6978	7419	68
	16	8 1 9 8	7215	13070	9282	15715	89
	32	11312	10326	16921	14002	24070	115
	2	3327	3 2 2 1	2966	2776	3644	31
	4	5922	5 735	6219	5898	9635	91
lu	8	10218	9831	10971	10837	20432	185
	16	15319	15145	15735	15152	27899	276
	32	22034	21652	23252	22984	36568	361
	2	2952	2887	3 0 2 0	2917	4364	38
	4	7546	7468	7631	7479	11193	107
ludcmp	8	12568	12494	12557	12322	22516	221
-	16	18211	17933	20093	19412	31115	304
	32	25273	24491	26447	26164	42154	421
	2	446	404	3247	558	11174	4
	4	1069	818	2988	1664	14887	85
mvt	8	2425	1648	6860	4187	20852	149
	16	2851	2586	13203	10041	32053	233
	32	6288	4295	14809	10009	40295	346
	2	838	838	996	935	1275	9
	4	2582	2473	2775	2672	3349	27
seidel-2d	8	4668	4274	6020	5403	7265	49
	16	7247	6580	10166	9045	14873	91
	32	10869	9 966	15649	13240	26383	156
	2	836	820	2915	2 3 4 6	2946	28
	4	2630	2540	4963	4370	7034	60
symm	8	6257	6107	9 0 2 3	8862	11819	96
0	16	10721	10445	13520	13251	20199	197
	32	15672	15282	18851	18594	33848	321
	2	900	880	1 1 3 9	900	1 356	9
	4	1 938	1820	2 3 2 7	1978	3 062	19
syr2k	8	3834	3 372	3913	3 1 9 8	6 0 1 0	37
-	16	5579	4967	5868	5294	10551	53
	32	7912	7590	8621	7 5 2 0	16163	96
	2	3 2 4 0	3 2 4 0	3 393	3 3 7 6	3854	3 2
	4	7 3 9 0	7 3 2 0	10 083	10079	10 431	94
syrk	8	13566	13 202	13924	13924	19379	171
-	16	20121	19674	31052	30 851	30102	287
	32	28222	27446	42805	42805	47622	467
	2	279	279	280	279	283	2
	4	620	595	777	600	643	5
trisolv	8	1 088	1054	1260	1 1 3 3	1 366	12
	16	1 788	1742	2008	1808	2622	24
	32	2 783	2 683	3 3 4 7	3 0 2 0	4 4 2 0	39
	2	2 704	1844	3 755	3 579	4 113	34
	4	6 2 2 6	5 673	7 311	7 167	11452	87
trmm	8	10 082	9914	13 669	13 484	19 559	124
	16	16002 16173	15472	20 933	20 933	30 0 26	203
	32	22126	21437	20355 27168	20333 27168	42503	41.8
	52						
Mean		4447	3 900	5698	4853	8247	6.0