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# L<sup>A</sup>T<sub>E</sub>X Report Template

Semester Project / Master Project

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# Acknowledgements

# Abstract

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# Declaration of Originality

I hereby confirm that I am the sole author of the written work here enclosed and that I have compiled it in my own words. Parts excepted are corrections of form and content by the supervisor. For a detailed version of the declaration of originality, please refer to Appendix ??

Pierre-Hugues BLELLY,  
Zurich, May 2020

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# List of Acronyms

AARCH64	. . .	.64 bit ARM architecture
API	. . . . .	.Application Programming Interface
CPU	. . . . .	.Central Processing Unit
CUDA	. . . . .	.Compute Unified Device Architecture
EEES	. . . . .	.Energy Efficient Embedded Systems
ETH Zürich	. . .	.Eidgenössische Technische Hochschule Zürich
FPGA	. . . . .	.Field Programmable Gate Array
GPU	. . . . .	.Graphic Processing Unit
HERO	. . . . .	.Heterogeneous Embedded Research Platform
IIS	. . . . .	.Integrated Systems Laboratory
ISA	. . . . .	.Instruction Set Architecture
llvm	. . . . .	.Low Level Virtual Machine
MIPS	. . . . .	.Microprocessor without Interlocked Pipelined Stages
MIT	. . . . .	.Massachusetts Institute of Technology
OpenCL	. . . . .	.Open Computing Language
OpenMP	. . . . .	.Open Multi-Processing



## *List of Acronyms*

PCMA . . . . .Programmable Many Core Accelerator

PULP . . . . .Parallel Ultra Low Power

RISC-V . . . . .RISC-V

SIMD . . . . .Single Instruction Multiple Data

SoC . . . . .System on Chip

ULP . . . . .Ultra Low Power

# Chapter 1

## Introduction

Thanks to the smaller nodes of modern lithography technologies and the transistor density we can achieve with them, modern low-power Central Processing Units (CPUs) can have a large amount of cores while keeping their power consumption under a few Watts. A single Raspberry Pi 3 has a peak performance of 6 gflop for a power consumption of only 7 Watts [1]. Embedded systems can take advantage of this increase in efficiency to become more autonomous and not rely on an external computer for heavy computation. Nano drones can now analyze in real-time a video signal and train a neural network for autonomous navigation, at a rate of 281MMAC/s on a power-enveloppe of 45mW [2].

To improve the energy efficiency and the computing power of Ultra Low Power (ULP) systems, new architectures are needed. To keep the power consumption low, an embedded system needs to power its subsystems only when needed. Autonomous drones [?] use a low-performance microcontroller to manage it coupled with a high-performance ULP cluster of RISC-V (RISC-V) cores for signal processing. System composed by one host processor and one or multiple coprocessors are called heterogeneous systems, they are suited for embedded applications as they can keep a low power consumption while using high-performance accelerator when needed.

This strategy has been used in the System on Chip (SoC) industry by ARM since 2011 [3]. The big.LITTLE architecture is based on two clusters of ARM Cortex A7(the “LITTLE”cores) and A15 (the “big” cores), and was designed to increase the computing power in low power systems such as smartphones while increasing the battery life of the device. This architecture relied on a single Instruction Set Architecture (ISA)(ARMv7). The goal was to use the more powerful cores during heavy computation or graphic rendering, and let the low power cores handle the background tasks or manage the device during sleep.

Researchers from the University of California [4] tried to leverage the advantages of multiple ISAs. Under heavy design constraints (such as die area or thermal dissipation)

## 1 Introduction

heterogeneous systems based on multiple ISAs performed better than the best homogeneous counterpart (in terms of energy efficiency and compute performance).

Even in data centers, where power consumption is also an issue, Graphic Processing Units (GPUs) are used thanks to their massive core count and the various Application Programming Interfaces (APIs) such as Compute Unified Device Architecture (CUDA) or Open Computing Language (OpenCL) which simplify the development process for GPU accelerators.

HERO [5] is a heterogeneous system developed by the Integrated Systems Laboratory (IIS) of ETH Zürich and the Energy Efficient Embedded Systems (EEES) of the University of Bologna. This platform is composed of a hard multicore ARM 64 Juno SoC (composed of two Cortex A57 and four Cortex A53 cores) and up to eight Parallel Ultra Low Power (PULP) clusters (composed of eight RI5CY cores [5]), running on an Field Programmable Gate Array (FPGA)(a Xilinx ZYNQ ZC706). The PULP cluster is based on the RISC-V ISA, an open source ISA designed to support a wide range of platform from embedded systems to supercomputer. The modularity of the ISA makes it interesting for Programmable Many Core Accelerators (PCMAAs).

This platform is designed to “facilitate rapid exploration on all software and hardware layers” [5], and includes a heterogeneous compilation toolchain with support for Open Multi-Processing (OpenMP), an API developed to make development of multi threaded applications easier [?]. This API implements new preprocessor instructions to tell the compiler how to execute the code on the system.

### 1.1 Design Issue with heterogeneous systems

During their conception, numerous design choices need to be made specify how the CPUs in the system will interact with each other. These choices will impact the peak performance of the design or its power consumption [4]. The computer architect has to choose how the different PCMAAs will interact, how they will share data, maybe extend the existing ISAs to distribute tasks, and so on.

The software design is challenging, when compiling for heterogeneous platforms. The compiler needs to create an executable that will run on the host processor, but also dedicate parts of the final binary to embed the code that will be distributed on the PCMAAs. Code distribution is handled by the programmer, APIs such as CUDA, OpenCL or OpenMP using function calls tell the compiler how to execute the code and on which PCMAA.

## 1.2 Currently Available Workflow for HERO

Currently, HERO supports OpenMP, which is an API which “defines a portable, scalable model with a simple and flexible interface for developing parallel applications on platforms from the desktop to the supercomputer” [6]. This API has been implemented on HERO to easily take advantage of the PULP clusters. The toolchain uses the Clang compiler [7] to compile the applications. HERO uses custom Clang front ends to supports all the available configurations (only the PULP cluster for simulation, with the ARM host CPU with a 64 bits RISC-V CPU).

To distribute the code, OpenMP uses preprocessor instructions to tell Clang where the code will run and how it will be executed. Exploring the design space using OpenMP’s directive can be time-consuming. For example, the developer must explicitly tell which part of the code to offload. Trying to change the order of multiple loops may cause bugs in the algorithmn, and complex schedules often impact code readability making them harder to debug

Halide [8] was proposed to explore the idea of separating the algorithm from the schedule. This separation makes testing different schedule easier on the developer, as the algorithm code will stay the same, and only the scheduling will be changed when testing. Every processing pipeline designed with Halide has two parts. The first part consists of the functional description of the processing kernel, i.e. the algorithm that will be executed. The second part is the schedule of the pipeline. The programer will explicitly tell Halide how the pipeline should be executed. Thanks to specific function calls, the developer can decide whether the code will be run on multiple threads or a single one, change the order of execution of different parts, split loops, unrolls them. The developer can still has the freedom to implement any schedule he wants but without having to change the main algorithm. This programming model is interesting because the developer can quickly implement the algorithm without having to take into account the boundaries of the inputs, and then work on an optimal schedule.

The intermediate variables can be bounded afterwards if needed, and the pricipal variables such as characteristics of the inputs are automatically bounded by Halide. An image processing pipeline will only compute the output on the pixels of the input. The scheduling process can even be done automatically during the compilation by the library, in order to find an optimal schedule on the target platform.

# Chapter 2

## Background

### 2.1 Hero

### 2.2 Halide Language

#### 2.2.1 Programing model

Halide is a functionnal programming language embedded into C++, designed to write high performance image and array-processing code [9]. This language uses a functionnal paradigm to describe the processing pipeline, and dissociate the array-processing code from the it's schedule (how the code will be compiled and run on the system).

Every pipeline is a function (`Halide::Func`) composed of other functions and expressions (`Halide::expr`). These two objects use special variables (`Halide::Vars`) to describe the operation executed on the array. The code listing 2.1 describe a basic pipeline which compute the distance of each coordinate of a two-dimensional array from on position specified by the vector (`center_x`, `center_y`). The creation of the pipeline is straightforward, we have to code the operation we want to execute on every element of the domain, for this example, we compute the square of the distance to the point (`center_x`, `center_y`).

```
Halide::Var x, y;
Halide::Param center_x, center_y;
Halide::Expr offset = Halide::pow(x - center_x, 2)
                      + Halide::pow(y - center_y, 2);

gradient(x, y) = offset;
```

Listing 2.1: Simple Pipeline Example

## 2 Background

This simple pipeline only has one stage, but it is possible to create multi-stage pipelines and schedule them as wanted. They can be transformed into a single-stage inlined pipeline or kept as is. The different stages can be scheduled to start as soon as they have enough data, or wait for the previous one to finish before starting to compute.

Scheduling is done via basic scheduling primitives implemented by Halide. They can be combined to create complex schedules. The primitives consists of basic code transformations such as loop unrolling or reordering, loop splitting or fusing variable together, more advanced instructions like parallelization or vectorization are also available. I will explain their behavior more in detail in the section Basic Scheduling Options.

In the example 2.2, we can see how the scheduling works. All instructions are a function of the pipeline object, they take the variables of the pipeline as argument to specify on which part of the pipeline this schedule will affect. During compilation, Halide will generate the code corresponding to the schedule specified by the programmer (if this schedule is valid).

```
gradient.parallel(x);  
gradient.unroll(y, 10);
```

Listing 2.2: Simple Pipeline Example

In the example 2.2, we will create as many task as the value x can take. They will then be distributed on the cores of the system. Every task will execute a single loop over the y axis, which will compute ten elements of the output every iterations (as we specified ten as argument of the `unroll` schedule).

The pipeline can be translated or compiled by Halide to be executed directly on the compilation computer or to be used in another application. The function `.realize(x_max, y_max)` is the fastest way to execute the pipeline, and is useful to debug the algorithm or the schedule. As halide was designed to work with different hardware platform, some of them being low power, cross compilation has been simplified to make the process as simple as possible.

Halide support translation to C code, Low Level Virtual Machine (llvm) assembly file, or already compiled object file specific to a given target(CUDA, ARM, RISC-V, Microprocessor without Interlocked Pipelined Stages (MIPS), PowerPc...), and a given operating system( Linux, Mac, Windows, Android). The pipeline can also be exported as a static library to use in another application.

### 2.2.2 Debugging Options

Halide has tools to debug the pipeline during it's compilation or during execution. First of all, the `print()` can be called any time to print a variable, the `print_when()` which print only when one boolean condition is met.

## 2 Background

Another useful tool is the `.trace_store()`, this function keeps a trace of every function evaluation in the pipeline. It is possible to get more informations during the compilation of the pipeline by setting the environmental variable `HL_DEBUG_CODEGEN` to 1, this will output information about every stages of the compilation and a pseudo code representation of the pipeline loops. Finally, variables and functions can have a label, which will be used by halide in it's internal representation or when printing the schedule, this function greatly reduce the debugging time of the schedules as Halide gives every variables a different name from the source code.

### 2.3 Basic Scheduling Options

Halide implement different scheduling instruction, and some of them just reshape the code in a different way. These scheduling instructions are useful to prepare the code for other instructions (such as parallelization or vectorization), but also to take advantage of memory locality.

#### 2.3.1 Non Platform Specific Schedule

##### Default Schedule

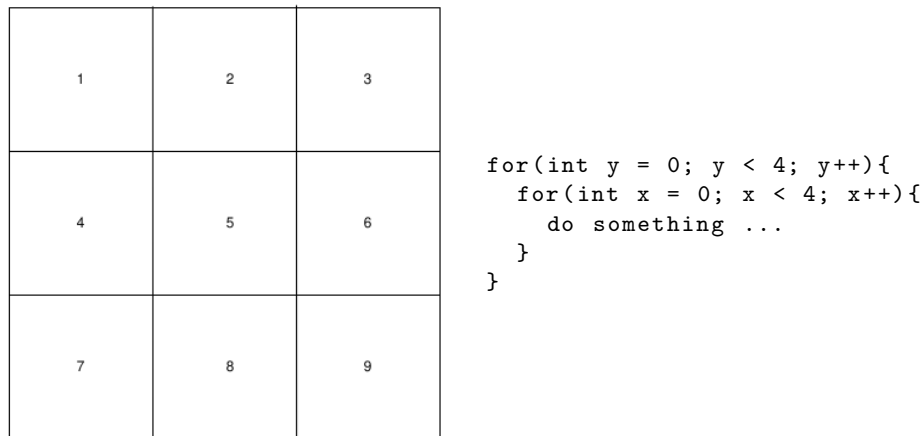


Figure 2.1: Base Schedule

If you don't specify any scheduling instructions, Halide will evaluate the pipeline in order. The first variable being the inner loop, and the last one the outer loop. In figure 2.1, we can see the schedule in action, the image is processed in a row major fashion.

## 2 Background

### Reorder

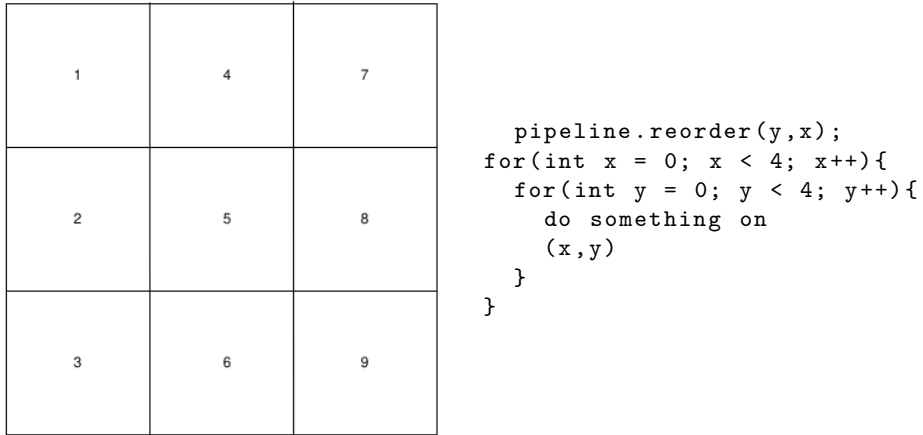


Figure 2.2: Schedule: Reorder

The `.reorder` instruction tells Halide how to traverse the domain space, using this instruction we can reorder the loops of the pipeline. In the example ??, we changed the way the array is being processed from row major to column major.

### Fuse

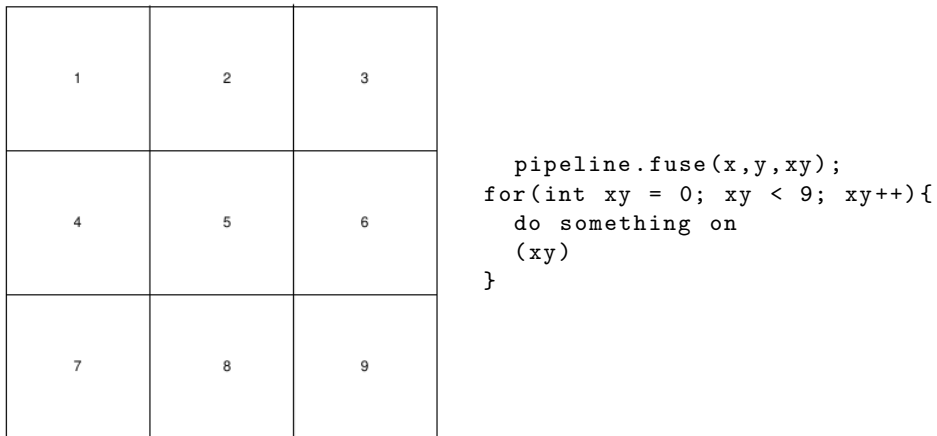
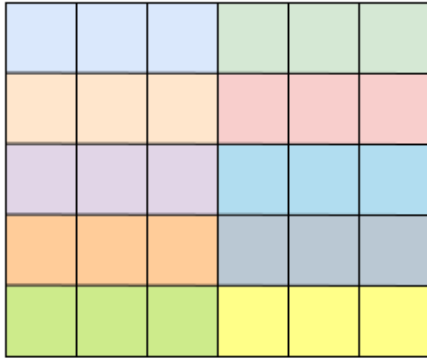


Figure 2.3: Schedule: Fused

The `.fused` instruction fuses two dimensions together, transforming a two-dimensional array into a one-dimensional array.



**Split**

```

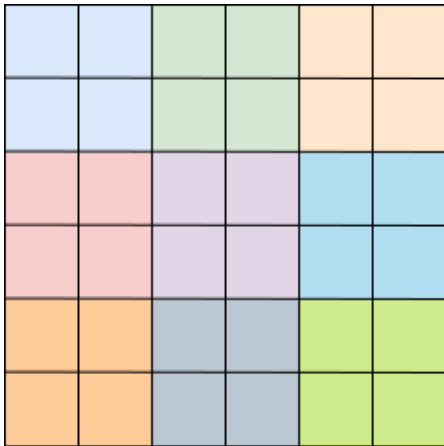
pipeline.split(x,x_o,x_i, 3);
for(int y = 0; y < 6; y++){
  for(int x_o = 0; x_o < 3; x_o++){
    for(int x_i = 0; x_i < 4; x_i++){

      do something on
      (x_o * 3 + x_i, y)
    }
  }
}

```

Figure 2.4: Schedule Split

This schedule replaces one loop over a dimension by two loops, an inner loop and an outer loop. This schedule is useful to cut the array in smaller pieces that will be computed in parallel or using Single Instruction Multiple Data (SIMD) instructions.

**Tile**

```

pipeline.split(x,y,x_o,y_o,x_i,y_i,2,2);
for(int y_o = 0; y_o < 6; y_o++){
  for(int x_o = 0; x_o < 3; x_o++){
    for(int y_i = 0; y_i < 2; y_i++){
      for(int x_i = 0; x_i < 2; x_i++){
        do something on
        (x_o * 2 + x_i, y_o * 2 + y_i)
      }
    }
  }
}

```

Figure 2.5: Schedule Tile

The Tile schedule is similar to the Split schedule, but along two dimensions. It creates multiples smaller rectangles which can be processed independently.

## 2 Background

### Unroll

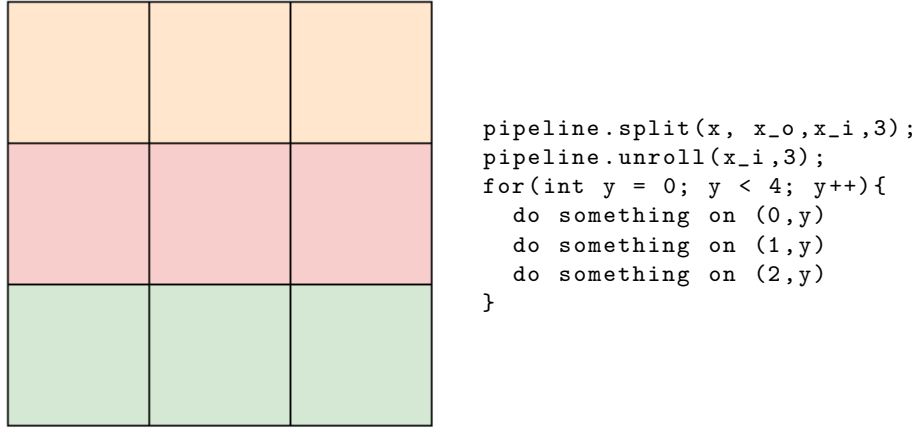


Figure 2.6: Unroll Schedule

The Unroll schedule unrolls the code along one dimension. This technique is often used when multiple computations share the same data, to prevent multiple memory access. In the example 2.6, we first split the x dimension before unrolling as Halide can't unroll a variable if it isn't bounded.

### 2.3.2 Platform Specific Schedules

#### Parallel

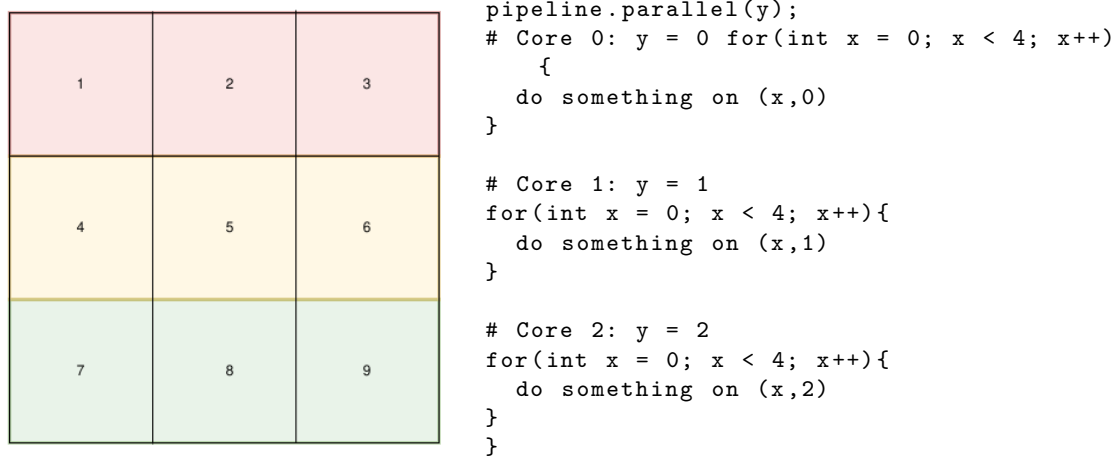


Figure 2.7: Schedule Parallel

## 2 Background

The parallel schedule allows the pipeline to be distributed to all the available cores. Halide will create for task for each value the variable can take, and these tasks will be executed with the `halide_do_par_for` function. This function has been overwritten on hero to execute on the PULP cluster. In the example 2.7, the code is distributed on three cores, each of them execute a single loop along the y axis.

### Vectorize

The goal of this schedule is to setup the code so to make use of the SIMD instructions of the CPU. Currently, llvm doesn't support the vector extension implemented in the pulp cluster, but the generated code will take advantages of all the registers available to compute the output values, and try to compute multiple values at the same time.

### 2.3.3 Porting Halide to new Platforms

In order to compile Halide, we need to compile llvm with the flag without shared libraries otherwise, Halide won't compile, and with support for the desired targets (x86, RISC-V and 64 bit ARM architecture (AArch64)).

Using the comment inside the pipeline header file, we can determine which function we need to implement to make Halide work on our target platform. The error messages during the linking phase are also a good source of information to find which function are needed to compile the code.

Currently only the memory allocation functions, the print functions and the task distribution functions are implemented, and they are enough to test basic pipelines such as matrix multiplications or light image modifications. After the implementation we can work on the compilation workflow for hero.

## 2.4 Compilation Workflow

Every application has at least two source files, one C++ file which will generate the object file of the pipeline, the main application. Currently, we can only compile the application to the hardware simulator. The compilation has two phases, during the first one, we compile the Halide application using llvm and run it on the host platform, this application will then generate an risc-V object file and a header. Then we compile the hero application using the already available Makefile, we include the header in the main application and the object file to the sources during the linking command.

## 2.5 The full hero platform

The hardware platform has a more complex compiling process, currently the code is distributed to the PULP thanks to OpenMp. The compilation first generate the llvm representation of the code, then assign space on the device via **hc-omp-space**, and also **clang-offload-bundler** to distribute generate the llvm assembly code for the right platform. Finally the program uses clang to compile the application, thanks to the special hero target, clang links every function correctly and then embed the riscV code inside the ARM application.

# Chapter 3

## Design Implementation

To test halide on hero, I used two benchmark. The first one was a basic gradient example, and the second one a matrix multiplication pipeline that I took in the provided examples and then adapted to be used in a hero application. The matrix example is more interesting, because it represent what a typical signal processing application may do. It is also quite easy to benchmark with different sizes to see the impact of the memory access on the execution time.

```
ImageParam A(type_of<int>(), 2);
ImageParam B(type_of<int>(), 2);
Var x, y;
Func matrix_mul("matrix_mul");
Func out;

RDom k( 0, A.width() );

matrix_mul(x, y) += A(x, k) * B(k, y);

out(x, y) = matrix_mul(x, y);
```

Listing 3.1: Matrix Multiplication Pipeline

### 3.1 Schedule Implementation

Most of the schedules implemented on halide doesn't require any platform specific implementation as they are only unrolling, splitting or swapping loops. During my project I used two platform specific schedules: vectorization and parallelization. The `.vectorize(x)` instruction unrolls one loop in assembly, and the vectorization is done by g++ using the SIMD instructions of the chip. For hero, the simd extention wasn't supported by g++

### *3 Design Implementation*

but, we could still use this instructions as it reduced the number of jumps and thus the total execution time.

The `.parallel(x)` instructions uses two functions: `halide_do_par_for` and `halide_do_par_for_fork`. `halide_do_par_for` adds the tasks to the task queue of the pulp cluster, every task will execute `halide_do_par_for_fork` on the corresponding core (if the core id is equal to the task number modulo the number of available cores). Every task consist of a part of the processing pipeline.

# Chapter 4

## Results

### 4.1 Test Setup

To compare Halide and OpenMp, I ran a matrix multiplication program coded using OpenMp and Halide. I ran the applications in the hardware simulation, using only the PULP cluster. The matrices for the Halide application were randomly generated and for the OpenMP application I generated them using a predefined pattern. As the multiplication operation always takes two cycles, the value inside the matrices doesn't matter as long as they have the same size. I made sure that the two matrices were stored in the L1 cache, to have the best access time possible. To measure the number of cycles needed to run the application, I used two functions available in the hero sdk: `hero_reset_clk_counter()` and `hero_get_clk_counter()`. These functions reset and output the value of a counter incremented every cycle, as they only take less than twenty cycles to execute, they are useful to get cycles accurate measurements of the execution time of the function. With this setup, we can easily compare the performances of Halide and OpenMp in a real world scenario for at least two basic schedules: Single threaded and Multi Threaded. I then experimented with different schedules with Halide to see the maximal performance I could get with this application.

To give the results more meaning, I converted the benchmark data in operations per cycles where one operation can either be an addition or a multiplication, so for a matrix of size  $n$ , the number of operations to finish the multiplication is :  $2n^3$ .

#### 4 Results

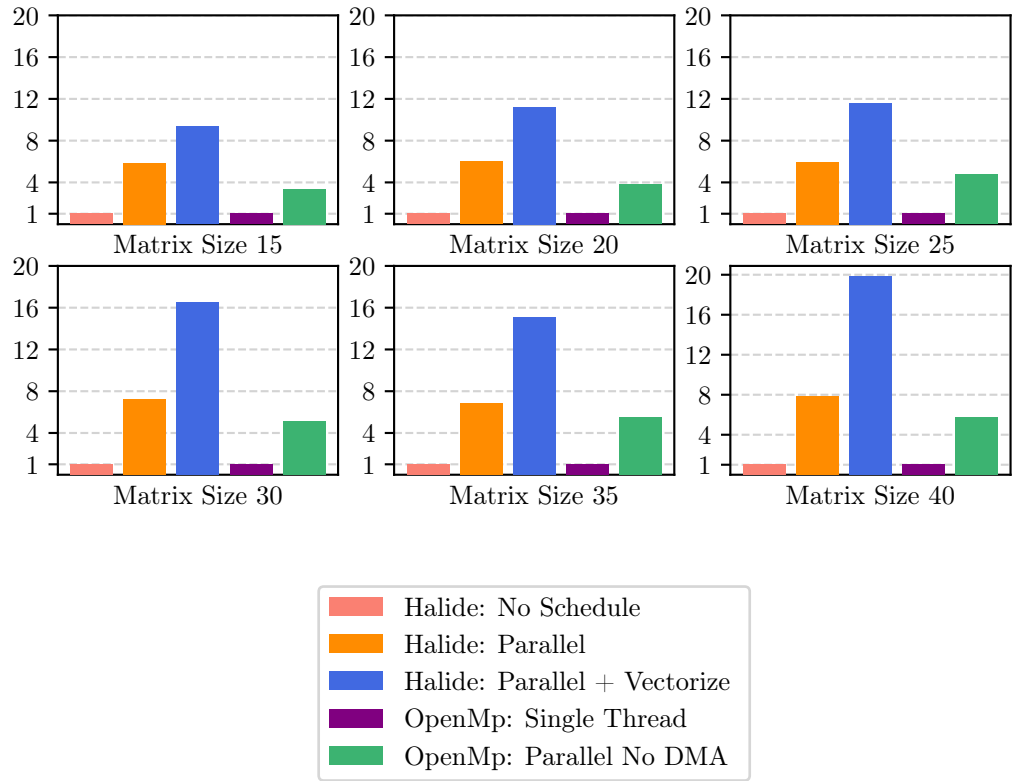


Figure 4.1: Halide vs OpenMP, relative to halide base schedule



## 4.2 Comparaison between OpenMp and Halide on the different platforms

Halide base schedule performs similarly from OpenMP single threaded, differing only by a few thousands cycles for the last test (with two matrices of size 35 by 35). The two implementation are prettty similar, and Halide overhead is in this case almost negligible. The second schedule I tested on both APIs was the parallel schedule; as paralellisation is the most efficient way to increase performance especially as there is no data dependancy in this pipeline. Often, paralellization instructions do have a lot of overhead which may impact the final performance of the code. To represent the code efficiency, I used the number of useful operations per cycles. As we have eight core at our disposal, capable at most of one operation per cycle, we can achieve at best eight operations per cycle. The graph 4.2, represent the performance of the schedule relative to the computing intensity of the program (The number of operations required to output one byte of data). The computational intensity for a matrix of size n is  $C_n = \frac{2n^3}{12n^2} = \frac{n}{6}$  where  $2n^3$  is the number of operations, and  $12n^2$  is the number of byte needed to complete the operation on 32 bits integers.

On the graph 4.2, we can see that Halide with it's parallel schedule achieves better results than OpenMP on all the sizes tested( from fifteen by fifteen up to thirty five by thirty five). The gaps in performance stays between .2 operations per cycles and .4 operations per cycles depending on the size of the matrices.

Schedule	15x15	20x20	25x25
Halide: No Schedule	40628 (0.166)	93818 (0.171)	180686 (0.173)
Halide: Parallel	6950 (0.971)	15585 (1.027)	30413 (1.028)
Halide: Parallel + Vectorize	4339 (1.556)	8358 (1.914)	15585 (2.005)
OpenMp: Single Thread	39820 (0.17)	92650 (0.173)	179030 (0.175)
OpenMp: Parallel No DMA	12079 (0.559)	24750 (0.646)	38090 (0.82)

Schedule	30x30	35x35	40x40
Halide: No Schedule	309426 (0.175)	488316 (0.176)	725606 (0.176)
Halide: Parallel	42659 (1.266)	71279 (1.203)	92536 (1.383)
Halide: Parallel + Vectorize	18776 (2.876)	32295 (2.655)	36487 (3.508)
OpenMp: Single Thread	307210 (0.176)	485440 (0.177)	721970 (0.177)
OpenMp: Parallel No DMA	59887 (0.902)	89283 (0.96)	126523 (1.012)

## 4 Results

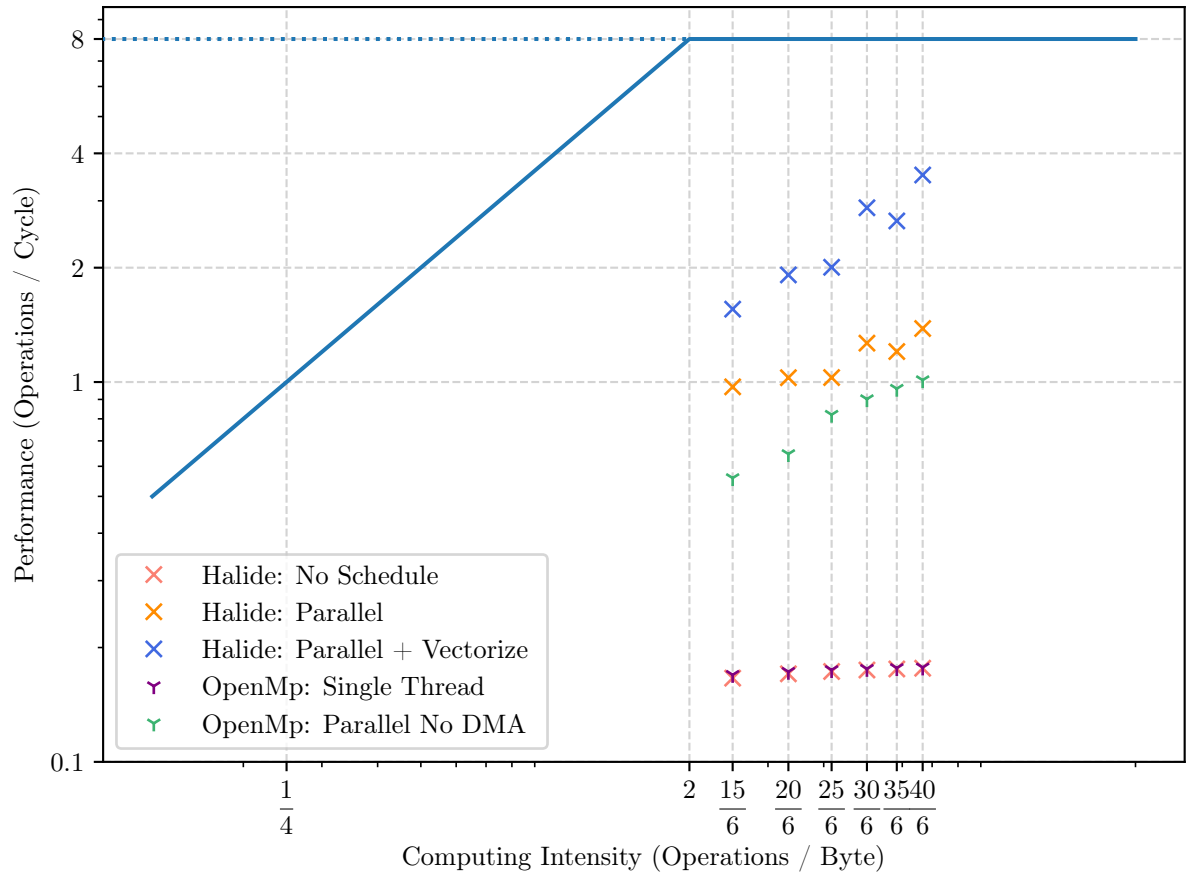


Figure 4.2: A PGF histogram from `matplotlib`.

# Chapter 5

## Conclusion and Future Work

Draw your conclusions from the results you achieved and summarize your contributions. Comparisons (e.g., of hardware figures) with related work are also appropriate here. Point out things that could or need to be investigated further.

### 5.1 First Section

### 5.2 Second Section

## 5 *Conclusion and Future Work*

# Bibliography

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