

## DEPARTMENT OF INFORMATION TECHNOLOGY AND ELECTRICAL ENGINEERING

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## LATEX Report Template

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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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### Introduction

Thanks to the smaller nodes of modern lithography technologies and the transistor density we can achieve with them, modern Central Processing Units (CPUs) in embedded systems, as capable enough to be used in data center. A single Raspberry Pi 3 has a peak performance of 6 Giga Floating Operations per Second Institute of Technologys (GFLOPs) for a power consumption of only 7 Watts ??. Embedded systems can take advantage to becomes more autonomous and not rely on external computer for heavy computation. Nano drones can now analyse in real time a video signal and train a neural network for autonomous navigation under 100mW ??.

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 it's subsystems only when needed. Autonomous drones ??, uses a low performance micro controller to manage the drone coupled with a high performance cluster foro the signal processing tasks. System using one host processor and one or multiple coprocessors are called heterogeneous systems. They are suited for embedded systems 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, The big.LITTLE architecture [?] is based on on two clusters of ARM Cortex A7 and A15, 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 also tried to leverage the advantages of multiple ISAs, according to this article ??, under heavy design constraints (such as die area or thermal dissipation) 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 comsumption 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 make them capable of heavy computation.

Hero [?] is a heterogeneous system developed by the Integrated Systems Laboratory (IIS) of ETHZ 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) and up to eight Parallel Ultra Low Power (PULP) clusters (composed of eight RI5CY cores), running on an Field Programmable Gate Array (FPGA).

#### 1.1 Design Issue with heterogeneous systems

Due to their heterogeneous nature, those systems are difficult to design and program. During their conception, numerous design choices need to be made to make sure that all the CPU in the system will interact will each other, these choices will impact the peak performance of the design or it's power consumption [?]. The architect has to choose how the different accelerators will interact, how they will share data, maybe extend the existing ISA to distribute tasks, and so on.

The software design isn't easy either, when compiling for such systems, the compiler needs to create an executable that will run on the host processor, but also integrate instructions that will be run on the coprocessors. Most of the time, the programmer assign which part of the code will run on which accelerator, OpenMP uses this model via the #pragma preprocessor instructions. Then during the compilation, llvm will compile the code bits to the right platform and link them using the adequate linker.

### 1.2 Currently Available Workflow for Halide

Currently Hero supports OpenMP, which is an APIwhich "defines a portable, scalable model with a simple and flexible interface for developing parallel applications on platforms from the desktop to the supercomputer" [?]. This API has been implemented on hero to easily take advantage of the clusters. The toolchain uses clang, and the clang-offload-bundler to compile the applications. Clang uses a custom front head for hero which supports all the available configurations for hero (only the pulp cluster for simulation, with the ARM CPU as the host for the FPGA or with a riscV 64 bits CPU).

Exploring the design space using OpenMp's directive isn't perfect, and require the developer to adapt its code to run with a specific schedule, for example, to take advantage of vectorization, the programmer has to manually unroll the loops and change it's code so

#### 1 Introduction

that the compiler knows what part of the code will be using the Single Instruction Multiple Data (SIMD) instructions. This approach leads to an important development time but also an extensive testing process whenever the schedule is changed to ensure that the resulting code works as intended. This approach makes testing for a new schedule pretty difficult and inefficient.

Starting from this idea that separating the algorithm from how it is run, researchers from the Massachusett Institute of Technology (MIT), created Halide. Halide [?] is a programming language that was designed to allow the developer to explore multiple design choices quickly by separating the algorithm from the execution schedule. This language was designed to be used in image or array processing applications. Every processing pipeline designed with Halide has two parts. The first part consist of the functional description of the processing kernel, this is the algorithm that will be executed on the array. The second part is the schedule of the pipeline. This schedule describes how the algorithm will be executed on the system. This programming model is interesting because the developer can implement the algorithm without having to take into account the boundaries of the functions or the border effects. Then he can quickly bound the different variables of the pipeline and design it's schedule afterward. All the constraints will be asserted during the compilation of the pipeline without any intervention from the developer. 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.



## Preliminaries / 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 [?]. This language uses a functionnal paradigm to describe the functionnalities of the processing pipeline. The code of the algorithm is separated from how it will be implemented on the target (schedule).

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 snippet 2.1 describe a basic pipeline which compute the distance of each coordinate of a two-dimentionnal array from on position specified by the vector (center\_x, center\_y).

This simple pipeline only has one stage, but it is possible to create multiple stage pipeline and transform it into a single stage inlined pipeline or keep it's multi stage structure, this operation will be done during the scheduling phase.

After designing the pipeline, we can define it's schedule via the different directive included in Halide. Halide implements all the basic scheduling option like parallelizing, unrolling the loops, splitting one loop into an inner and an outer loop... These options will be described in the section Basic Scheduling Options.

In the example 2.2, we can see how the scheduling works. All instructions are function of the pipeline object, and the final pipeline will implement these instructions. The example shows a simple schedule applied on our gradient, this schedule consists of parallelizing the execution over the x axis, and unrolling along the y axis.

To execute the pipeline, Halide provides a large range of options, we can execute it directly using the .realize(x\_max, y\_max) function, this is useful for debugging purpose, but most of the compile options are targeted at cross-compilation.

As the initial goal of Halide was to target devices such as CPU, the library is capable of compiling the pipeline to a lot of different platform and output format. Halide support translation to C code, llvm assembly file, or already compiled object file specific to a given target(Cuda, Arm, Risc-V, MIPS, PowerPc...), and a given operating system(Linux, Mac, Windows, Android). The developer can also chose to compile the pipeline to a library to use in another application.

#### 2.2.2 Debugging Options

Halide has tools to debug the pipeline during it's compilation or when it is executed. First of all, the print() and print\_when() functions can be called at any time in a pipeline and allow to print values of some variables. Another useful tool is the .trace\_store() function which prints the value every functions evaluated in the pipeline. It is possible to get more informations during the compilation of the pipeline by setting the HL\_DEBUG\_CODEGEN to 1, this will output the stages of the compilation and a pseudo code representation of the pipeline. Finally, variables and functions can have a label, which will be used by halide in it's internal representation, 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

#### 2 Preliminaries / Background

other instrutions (such as parallelization or vectorization), but also to take advantage of memory locality.

#### 2.3.1 Non Platform Specific Schedule

#### Default Schedule

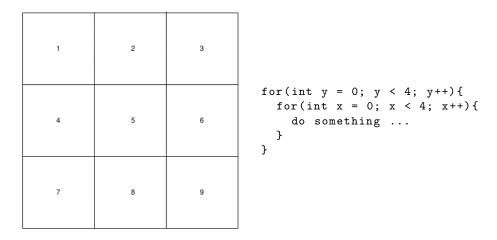


Figure 2.1: Base Schedule

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.

#### Reorder

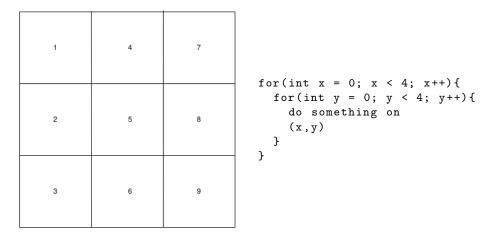


Figure 2.2: Schedule: Reorder

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

#### Unroll

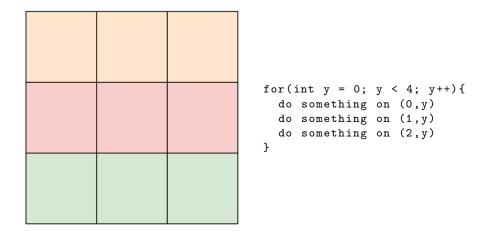


Figure 2.3: Unroll Schedule

The Unroll schedule unrolls the code along one dimension. This instruction sometimes improve the code execution speed by removing the jumps at the end of each loop. This technique is often used by compilers to optimize the speed of the executable.

#### Split

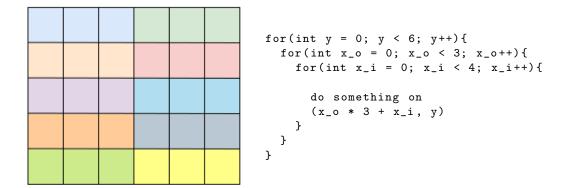


Figure 2.4: Schedule Split

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 SIMD instructions.

#### Tile

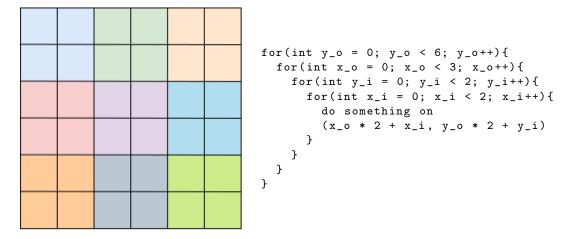
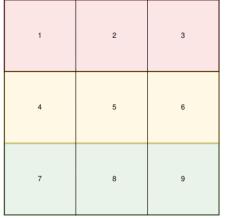


Figure 2.5: Schedule Tile

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

#### 2.3.2 Platform Specific Schedules

#### Parallel



```
# Core 0: y = 0
for(int x = 0; x < 4; x++) {
   do something on (x,0)
}
# Core 1: y = 1
for(int x = 0; x < 4; x++) {
   do something on (x,1)
}
# Core 2: y = 2
for(int x = 0; x < 4; x++) {
   do something on (x,2)
}
}</pre>
```

Figure 2.6: Schedule Parallel

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 2.6, 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, ans try to compute multiple values at the same time.

#### 2.3.3 Porting Halide to new Platforms

First of all, to compile to a specific platform, we need a build of llvm which support the desired architecture. Then we can look in the pipeline header file to list all the vital functions for our pipeline. We can also use the error message when linking the pipeline to determine which functions we need to implement on the target platform. Currently only the memory allocation functions, the print functions and the task distribution functions

#### 2 Preliminaries / Background

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 axample 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.



### Results

#### 4.1 Test Setup

I benchmarked two applications on two platforms. I benchmarked the halide port on the hardware simulation for the PULP cluste, and one openMp matrix multiplication application on the developpement platform on a Xilinx ZCU102. For the halide Application I generated random matrices, and for the openMp application I generated them using the same pattern every time. But as one multiplication takes two cycles every time, the execution time doesn't depend on the content of the matrices. The two matrices were stored in the L1 cache, to have the best access time we could, and to compare the code efficiency of both API. 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 resets 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 schedule 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 a multiplication or a memory access (which take 2 cycles each), so for a matrix of size n, the number of operations to finish the multiplication is : 2 \* n \* \*3.

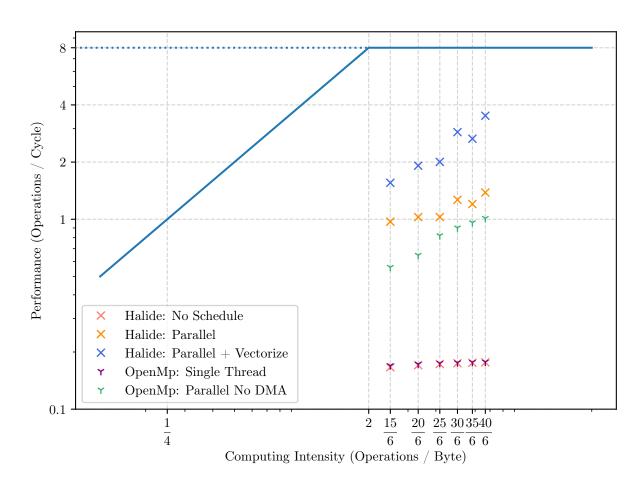


Figure 4.1: A PGF histogram from matplotlib.

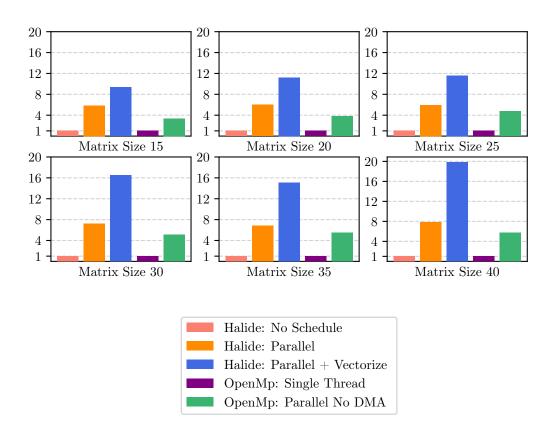


Figure 4.2: A PGF histogram from matplotlib.

#### 4 Results

4.2 Comparaison between OpenMp and Halide on the different platforms



### 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