

DEPARTMENT OF INFORMATION TECHNOLOGY AND ELECTRICAL ENGINEERING

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Acknowledgements

Ceci est un test de mon script

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

1.1 Heterogeneous systems

Heterogeneous Systems, relies on different accelerators to achieve the bbest possible efficiency. Most of them are composed of one Genereal Purpose Core and one or multiple Programmable Many Core Accelerator (PCMA). These systems are interesting because of their energy efficiency, the accelerators are designed to compute a lot of data in parallel, and the host allow for more flexibility and handle the program flow.

Hero is an heterogeneous system developped by the IIS (ETHZ) and the EEES (University of Bologna). This platform is composed of a hard multicore ARM 64 Juno SOC and one pulp cluster (eight RI5CY cores), running on an FPGA. During my project, I used this target to port Halide.

1.2 Design Issue with heterogeneous systems

Due to their heterogeneous nature, those systems are difficult to program. The compilation process is complex, requiering multiple passes on different toolchains to distribute the work to the host and the accelerator, and to allocate the memory. And this complexity is also forced on the end user, as he needs to know the target perfectly to make use of all the available resources.

1.3 Currently Available Workflow for Halide

Currently Hero support OpenMp, which is an API which "defines a portable, scalable model with a simple and flexible interface for developing parallel applications on plat-

1 Introduction

forms from the desktop to the supercomputer" (Ref: OpenMP website). This API has been implemented on hero, and is currently used to develop some test application. But exploring the design space using OpenMp's directive, isn't perfect, and require the developper to adapt it's code to run with a specific schedule. This approach leads to an important developpement time but also an extensive texting process whenever the schedule is changed to ensure that the resulting code works as intended.

Halide is a programming language that was designed to allow the developper 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 have two parts. Teh first part consist of the functionnal description of the processing kernel, this is the algorithm that will be executed on the arry. And the second part is the schedule of the pipeline. This schedule describe how the algorithm will be executed on the system. This programming model is interesting because the developper can in the first time 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 afterwards. All the constraints will be asserted during the compilation without any intervention from the developper.

 $^{\circ}$ Chapter $^{\circ}$

Preliminaries / Background

2.1 Hero

2.2 Halide Language

2.2.1 Programing model

Halide is a functionnal program embedded into C++ designed to write high performance image and array processing code (Halide Website). This language uses a functionnal paradigm to describe the functionnalities of the pipeline. The scheduling of the pipeline is described separately, which allow the developper to explore a wide range of schedule without having to rewrite most of the code. 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 describe a basic pipeline which compute the distance of each coordinate of the array from on position specified by the vector

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, splitting ... These options will be described in the section Basic Scheduling Options .

2 Preliminaries / Background

The snippet 2.2 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.

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

Listing 2.2: Simple Pipeline Example

To execute our pipeline, Halide provides a large range of options, we can execute it directly using the .realize(x_max, y_max) directive, this will execute the pipeline on a rectangle starting from it's top left corner in (0,0) to it's bottom right corner in (x_max, y_max).

But Halide also gives the programmer a lot off options to execute the pipeline, it's able to convert the code to C code, llvm assembly file, or already compiled object file specific to a given target. More over, Halide support a wire variety of CPU architecture (X86, ARM, MIPS, PowerPc, Risc-V), operating systems (Linux, Windows, Android, Mac) and also Gpu Api's (Cuda, OpenCL, OpenGl, DirectX ...). Halide support for new architecture is getting better and better, and is by design targeted for cross compilation and Heterogeneous systems.

2.2.2 Debugging Options

2.2.3 Basic Scheduling Options

Halide implement different scheduling instruction, and most of them aren't architecture specific.

Non Architecture Specific Instrutions

- Reorder: Tells halide how to traverse the domain of the pipeline stage (ffor example in a column major or row major way)
- Split: Split a dimension along inner and outer subdimensions.
- Tile: Cut the domain in tiles.
- Fuse: Join two dimensions in a single fused dimension.
- Unroll: Unroll along one dimension.

Some of the primitives such as vectorize or parallel. needs to be implemented on the target platform as they take advantage of the specificities of the system. To do so Halide uses some functions which are defined in a header file we can get when compiling the pipeline.

2.2.4 Porting Halide to new Platforms

From the header file, we can find the functions vital for halide to work and implement them in the pulp runtime. From then, we have to compile the pipeline to a risc-V object file, and then compile the main application using this object file and the provided header. When we will compile the final application using the hero toolchain, the linker will link the halide function calls to the implemented functions in the pulp runtime. Currently, only the memory allocation, print, and fork primitive are implemented but they are sufficient to try some basic parallel schdeules.

2.3 Compilation Workflow

Hero currently have different platforms, and also different workflow. I started by working on the simulation platform which simulate an eight-core pulp cluster. Then I tried to port it to the hardware platform (One hard ARM core and a PULP cluster implemented on the FPGA). The compilation is slightly different for both platform, so I started by working on the simulation platform which is not heterogeneous.

2.4 Simulation

The compilation process for the simulation platform is quite easy, we start by compiling the C++ code which will generate the final pipeline, we then run this application, to generate the pipeline and the matching header file. Then we compile the true application which will call the pipeline. During the process, we add the object file of the pipeline in the source file, g++ will then link the pipeline with the main application and also the halide calls to the pulp runtime functions.

2.5 The full hero platform

The hardware platform has a more complex compiling process, currently, the applications use OpenMp to distribute the code to the pulp cluster. With instructions such as #pragma omp target device(BIGPULP_MEMCPY) to explicitly tell the compiler to distribute the following part of the code to the pulp cluster. The application is compiled using Clang and a the clang-offlad bundler to create an heterogeneous application that will run on both platform.

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 both application, I generated random matrices of différent sizes, and for each sizes multiplication I counted the number of cycles needed to complete the operation. 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. To give the results a more meaning I also calculated the number of 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 : 6*n**3+n**2 (each coefficient needs 2n memory accesses, 2n additions and multiplications and one memory store).

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

Glossary

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