Sorbonne Université – SESI M2

#### MU5IN160 – Parallel Programming

## Hands-on Session 6 – Multi-threading for MOTION Application

**Short introduction** In this session, we will work on the MOTION streaming application. In the last session, we focused on converting the MOTION application through an explicit dataflow representation with the AFF3CT DSEL. Today, we will benefit from the code transformations we made before to parallelize the application over the CPU multi-core architecture.

Very important, about the submission of your work At the end of this session you will have to upload the following files on Moodle: 1) a zip of the src folder and 2) a zip of the include folder. After that you will have 2 weeks (until November, 5) to complete your work and update your first submission. You have to work in group of two people but each of you will have to upload the files on Moodle. Finally, please write your name plus the name of your pair at the top of all these files.

# 1 Appetizer

First you need to update the repository of the MOTION project:

git pull origin master && git submodule update --init --recursive

Then, you have to re-generate the Makefike and to re-compile a sub-part of the application:

```
cd build && cmake .. && make -j4
```

# 2 MOTION Parallelization

In this section, we will focus on reducing the execution time by using different types of parallelism. We will mainly rely on the multi-core architecture ( $\approx$  multi-thread programming) of the CPU. To measure the execution time, we will not consider the video decoding time, the visualization time and the logs time.

MOTION comes with the --vid-in-buff MOTION parameter to hide the video decoding time: the frames are decoded and put into a buffer during the video initialization. This behavior is representative of a real embedded application, where there would be no video decoding. Instead, a camera would send frames at a fixed rate. Be careful, buffering frames can take a lot of memory (and time). Thus, we will limit this buffer to 100 frames.

Here is the command line to use for the measurements:

```
./bin/motion2 --vid-in-buff --vid-in-path ./traffic/1080p_day_street_top_view_snow.mp4 \
--vid-in-stop 100 --flt-s-min 2000 --knn-d 50 --trk-obj-min 5
```

As a consequence, we will not take into account the whole execution time of the MOTION application but only the sequence or the pipeline execution time. And, we will focus on increasing its throughput (= number of frames per second or FPS).

Of course for debugging and validation purpose, you will need to use the logs (--log-path) and the visualization (--vid-out-play --vid-out-id) parameters.

Work to do Run the motion2 and motion2-aff3ct executables and note their throughput in FPS.

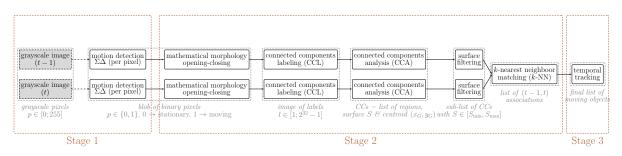


Figure 1: MOTION detection and tracking processing graph. In gray and italic: the output of each processing. The graph is split into a 3-stage pipeline.

### 2.1 Pipeline Strategy

AFF3CT comes with a round-robin *producer/consumer* implementation to support the *pipeline* parallelism (and more specifically the synchronizations between the pipeline stages). This type of parallelism is well adapted to the streaming applications like MOTION. It improves the throughput of the system.

We will instantiate the 3-stage pipeline given in Fig. 1. Other stages decomposition is possible but it will not lead to significant improvements (and it will not help for the next sections).

To instantiate a pipeline, we need to delimit its stages (= sub-parts of the graph). In AFF3CT, a stage is defined by 3 vectors of tasks (in the exact following order):

- 1. A list of the first tasks in the stage,
- 2. A list of the last tasks in the stage (these tasks are included in the stage),
- 3. A list of exceptions that represents the tasks that should not be included in the current stage (can be empty).

In other words, for each stage, AFF3CT will build the sub-graph from the first and last tasks and it will exclude the exception tasks. When building the sub-graph, if some of the tasks are bound to tasks that are in a previous stage, then AFF3CT will not automatically add these tasks to the current stage. In this case, you need to explicitly add these tasks into the list of first tasks.

Additionally, in order to facilitate the debugging process, the pipeline constructor requires the vectors of first and last tasks of the corresponding sequence. This way AFF3CT can compare if the stages match the initial sequence. If it does not match, AFF3CT will automatically return an error and will output two debug graphs (in fact, it will store these two files on the file system):

- dbg\_ref\_sequence.dot: the graph corresponding to the sequence (= without parallelization),
- dbg\_cur\_pipeline.dot: the graph corresponding to the pipeline (= with the stages).

Then, you can visually compare the two graphs and find the mismatch. Common mistakes are:

- A task is missing in the pipeline graph: this is often the case because this task needs to be explicitly declared in the list of first tasks,
- A same task is present in two (or more) different stages: this is generally because the task has been automatically added to a previous stage. In this case, using the exclusion list of tasks can solve the problem.

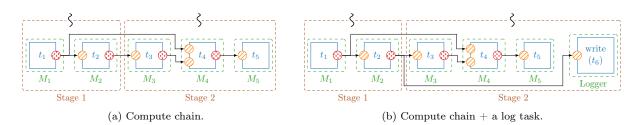


Figure 2: Example of a chain decomposed into two pipeline stages.

```
// 1) creation of the module objects
1
    M1 m1(); M2 m2(); M3 m3(); M4 m4(); M5 m5(); Logger log();
2
з
    // 2) binding of the tasks
^{4}
    m2["t2::in" ] = m1["t1::out"];
5
    m3["t3::in" ] = m2["t2::out"];
6
    m4["t4::in1"] = m3["t1::out"];
7
    m4["t4::in2"] = m3["t3::out"];
8
    m5["t5::in" ] = m4["t4::out"];
9
    m6["t6::in" ] = m5["t5::out"];
10
    // conditionnal binding of the log task
11
12
    if (is_log)
13
      log["write::in"] = m2["t2::out"];
14
    // definition of the pipeline stages in the 'pip_stages' variable
15
    std::vector<std::tuple<std::vector<runtime::Task*>,
16
                            std::vector<runtime::Task*>.
17
                            std::vector<runtime::Task*>>> pip_stages =
18
        { // pipeline stage 1
19
          std::make_tuple<std::vector<runtime::Task*>,
20
                           std::vector<runtime::Task*>,
21
                           std::vector<runtime::Task*>>(
^{22}
            { &m1("t1"), },
                                                                 // first tasks of stage 1
23
            { &m2("t2"), },
                                                                 // last tasks of stage 1
^{24}
            { /* no exclusion in this stage */ }
                                                                 // tasks excluded from stage 1
^{25}
26
          ).
          // pipeline stage 2
27
          std::make_tuple<std::vector<runtime::Task*>,
^{28}
                           std::vector<runtime::Task*>.
29
                           std::vector<runtime::Task*>>(
30
            { &m3("t3"), &m4("t4"), },
                                                                 // first tasks of stage 2
31
            { /* last tasks will be automatically found */ }, // last tasks of stage 2
32
            { /* no exclusion in this stage */ } ),
                                                                  // tasks excluded from stage 2
33
34
        /* if you'd had more stages, you could have continued their declaration here */
35
        };
36
37
    // if there is the log task
38
    if (is_log) {
39
        // in stage 1 ('pip_stages[0]'): add the log task at the end of the exclusion list ('get<2>')
40
        std::get<2>(pip_stages[0]).push_back(&log("write"));
41
        // in stage 2 ('pip_stages[1]'): add the log task at the end of the first tasks list ('get\langle 0 \rangle')
^{42}
        std::get<0>(pip_stages[1]).push_back(&log("write"));
43
    }
44
^{45}
    // the first tasks of the sequence, used to check the pipeline stages
46
    std::vector<runtime::Task*> seq_first_tasks = { &m1("t1") };
47
48
    // declaration (= construction) of a pipeline object 'pip'
49
    runtime::Pipeline pip(seq_first_tasks, pip_stages,
50
        {
              1,
                           1 }, // number of threads per stage -> one thread per stage
51
                                 // buffer size between stages -> size 1 between stage 1 and 2
        ſ
                   1.
                             },
52
                             }, // active waiting between stage 1 and stage 2 -> no
        {
                 false,
53
                      false }, // enable pinnig -> no
        { false,
54
                         {1} }); // pinning to threads -> ignored because pinning is disabled
55
        ſ
            {0},
56
    // pipeline execution, note that two different functions can be used for the stop condition
57
    pip.exec({ [] (const std::vector<const int*>& statuses) { return false; },
                                                                                     // stage 1 stop
58
                [] (const std::vector<const int*>& statuses) { return false; }, }); // stage 2 stop
59
```

Source code 1: Source code corresponding to the 2-stage pipeline described in Fig. 2. Depending on the is\_log value (true or false), the code will build either the Fig. 2a pipeline or the Fig. 2b pipeline.

Fig. 2 introduces a simple example with a 2-stage pipeline. The compute chain (= tasks that do not include logs and visualization) is given in Fig. 2a. Fig.2b shows a graph where we added a log task for debugging purpose. The main goal of this graphs is to help you to bootstrap in your work.

The corresponding source code is given in Code 1. It is interesting to note that for the first time we introduce a **conditional binding depending on the is\_log boolean value** (lines 11-13). Also, the list of the stages is declared in a dedicated **pip\_stages** variable (line 18) before to call the pipeline constructor (line 50). It makes possible to define the "compute" graph, and its decomposition in stages, only once. After that, depending on the **is\_log** value, the **pip\_stages** variable can be modified to add the log task (lines 39-44). Note that the  $t_4$  is added to the first tasks list of stage 2. This is because  $t_4$  depends on  $t_1$  which is declared in the stage 1. Finally, the stop condition of a pipeline can be set as one stop condition per stage (see lines 57-59). Generally, we will only customize the stop condition of the last stage.

Work to do #1 Implement the 3-stage pipeline given in Fig. 1 in the motion2\_aff3ct.cpp file (it is strongly recommended to save your previous working implementation with the sequence in an other file...). Moreover, your code should be able to enable/disable the logs and the visualization. Fig. 2 and Code 1 are here to help you and you must take inspiration from it. Validate that your new pipeline version is working (see Note #2 in the previous session). Measure and report the achieved FPS in the motion2\_aff3ct.cpp file.

Note #1 If you implemented a sequence with a Switcher (to skip the first execution of  $\Sigma\Delta$ ), there is a known problem: a select or a commute task cannot be directly at the interface of a synchronization between two pipeline stages. To make this work, you need to add a Relayer (relay task) after the select task. If you did not implemented a sequence with a Switcher then you can ignore this note.

The code for displaying pipeline task statistics differs slightly from the code for a sequence. Indeed, you first need to iterate over each stage before to print the statistics.

```
const bool ordered = true, display_throughput = false;
1
   auto stages = pip.get_stages();
2
   for (size_t s = 0; s < stages.size(); s++) {</pre>
3
       const int n_threads = stages[s]->get_n_threads();
4
       std::cout << "#" << std::endl << "# Pipeline stage " << (s + 1) << " ("
5
                  << n_threads << " thread(s)): " << std::endl;
6
       tools::Stats::show(stages[s]->get_tasks_per_types(), ordered, display_throughput);
7
   }
8
```

Source code 2: Display the statistics of the tasks for a pipeline.

Work to do #2 Modify the code that displays the task statistics in motion2\_aff3ct.cpp (see Code 2).

When using the --stats option, you should have an output similar to the Code 3. As you can see, the pipeline automatically added new push\_x and pull\_x tasks for you. These tasks are taking care of the communications and synchronizations between the pipeline stages. In other words, they represent a *producer/consumer* implementation. For instance, the push\_1 task of the Adp\_1\_to\_n\_0 module sends/produces output data from stage 1 to stage 2. And, in stage 2, the pull\_n task of the Adp\_1\_to\_n\_0 module receives/consumes the input data from stage 1. Sometimes push\_x and pull\_x tasks take a lot of time in the stage. It means that they are waiting.

Work to do #3 From the output statistics on the Jetson TX2, which stage limits the overall throughput of the pipeline? Note your answer in the motion2\_aff3ct.cpp file.

#	Statistics for the //			Basic statistics /			Measured latency 		
#	given task //			on the task /					
#		//-			/	/			
#	//	//-				· ·	/		
#	MODULE /	TASK	CALLS /	TIME					
#	1	11	1	(s)		(us)			
	//		/-		,				
	_to_n_0 / _delta0 /	push_1 // compute //	100   101	1.66 0.81		16592.90     8060.59		18569.21   8568.33	
0	_delta1 /	compute //	101	0.79					
# 009ma	Video /	generate //	100 /	0.01				· · · · · · · · · · · · · · · · · · ·	
	Delayer /	produce //	101 /	0.01			54.88		
	Delayer /	memorize	100 /	0.00					
#	//	//-	/-			/			
#	TOTAL /	* //	100	3.29	100.00	32932.16	16019.56	36316.68	
#									
# Pipel	ine stage 2 (	(1 thread(s)):							
#		//-			/	/			
#	Statistics for the //			Basic statistics /			/ Measured latency		
#	given task //			on the task /					
#		//-	,		/	/			
#	// MODULE /	//- TASK //	· · · · ·	TIME		 			
# #	MUDULE   I	1A5K    	CALLS	(s)	,				
#		, , -	1	(3)			(43)		
#	Morpho1	compute	100 /	1.32		13180.10			
	Morpho0	compute //	100 /	1.31	1 .	13134.11			
#	CCLO /	apply //	100 /	0.22		2183.68			
#	CCL1 /	apply //	100 /	0.22					
#	CCA1 /	extract //	100	0.13		1325.75		, 1960.42	
#	CCAO /	extract //	100 /	0.13	3.92	1314.09	1 724.04	1920.88	
$# Adp_1$	to_n_0 /	pull_n //	100	0.02	0.52	172.73	0.04	17246.83	
# $Ftr_{-}$	filter1 /	filter //	100 /	0.00	0.02	/ 7.02	4.38	1 27.46	
# $Ftr_{-}$	filter0 /	filter //	100 /	0.00	0.02 /	1 6.48	4.38	16.71	
#	KNN /	match	100	0.00			1		
$# Adp_1.$	_to_n_1 /	push_1 //	100	0.00					
#	//	//-	/-			/			
#	TOTAL /	* //	100 /	3.35	100.00	33499.82	32188.67	54816.59	
# # Dime1.		(1, 4b = a d(a))							
# ripel	ine stage 3 (	(1 thread(s)):			,	1			
#	Statistics for the //			Basic statistics /			isured later	200	
#	given task //			on the task /			/		
#		//-			· /	/			
#	//		/-	,	, 	/	/	/	
#	MODULE	TASK	CALLS /	TIME	I PERC	AVERAGE	MINIMUM	MAXIMUM	
#	1	11		(s)					
#	/	//-	/-			/		/	
# Adp_1	_to_n_1 /	pull_n //	100 /	3.35	99.98	33472.32	32689.25	52477.75	
# T1	racking /	perform	100	0.00	0.02	6.25	2.08	44.42	
#	/	//-	/ -			/	/	/	
#	TOTAL /	* //	100 /	3 35	1 100 00 1	33478.57	32691 31	1 52522 17	

Source code 3: Task statistics output for the pipeline described in Fig. 1. The automatically added Adaptor tasks (push\_x and pull\_x) are highlighted. The code ran on the Apple Silicon M1 Pro CPU. For each stage, the tasks are ordered by their execution time.

### 2.2 Tasks Replication

In Fig. 1, the  $\Sigma\Delta$  and the *tracking* tasks cannot be replicated because they have a data dependency over time. As a consequence, the tasks of the stages 1 et 2 cannot be replicated. These dependencies are not explicit in Fig. 1 because they come from inner data:

- In sigma\_delta\_compute function: the sd\_data->M and sd\_data->V buffers,
- In tracking\_perform function: tracking\_data->history and tracking\_data->tracks buffers.

However, in theory, stage 2 tasks can all be replicated. For instance, in the Code 1 you can perform the replication of stage 2 by replacing the second "1" line 51. If you replace { 1, 1 } with { 1, 4 }, the pipeline will replicate stage 2 four times and it will run a total of 5 threads (1 for stage 1 and 4 for stage 2).

Work to do #1 Modify the pipeline in the motion2\_aff3ct.cpp file to run the second stage over two threads (2 times replication). What is happening? Is it normal?

You should not be able to replicate stage 2 because some of the tasks are stateful. To overcome this problem you need to implement the **clone** method (and maybe the **deep\_copy** method) on the involved modules.

```
Morpho* Morpho::clone() const {
1
        auto m = new Morpho(*this);
2
        m->deep_copy(*this); // we override this method just after
3
        return m:
4
    }
\mathbf{5}
    // in the deep_copy method, 'this' is the newly allocated object while 'm' is the former object
6
7
    void Morpho::deep_copy(const Morpho& m) {
         // call the 'deep_copy' method of the Module class
8
        Module::deep_copy(m);
9
        // simply copy the 'i0, i1, j0, j1' variables to have shorter names later in the method
10
        int i0 = m.morpho_data->i0, i1 = m.morpho_data->i1;
11
        int j0 = m.morpho_data->j0, j1 = m.morpho_data->j1;
12
        // allocate new morpho inner data
13
        this->morpho_data = morpho_alloc_data(i0, i1, j0, j1);
14
        // initialize the previously allocated data
15
        morpho_init_data(this->morpho_data);
16
        // allocate the 2D containers for converting 1D to 2D images in the task % \mathcal{A} = \mathcal{A} = \mathcal{A}
17
        this->in_img = (const uint8_t**)malloc((size_t)(((i1 - i0) + 1) * sizeof(const uint8_t*)));
18
        this->in_img -= i0;
19
        this->out_img = (uint8_t**)malloc((size_t)(((i1 - i0) + 1) * sizeof(uint8_t*)));
20
        this->out_img -= i0;
21
    }
22
```

Source code 4: clone and deep\_copy methods of the Morpho module.

Code 4 give you the implementation of the clone and deep\_copy method for the Morpho module.

Work to do #2 Implement the missing clone and deep\_copy methods for stage 2 tasks. Then, run the pipeline with 2 replications in stage 2. Is it faster? Increase the number of threads in stage 2, what do you observe? Note the new FPS in your motion2\_aff3ct.cpp file.

### 2.3 Tasks Graph Simplification

Fig. 3 presents an optimized version of the tasks graph shown in Fig. 1. One may have noticed that in Fig. 1 some computations are performed twice! Indeed, the computations on the t - 1 frame could be memorized from a stream to the next one. In Fig. 3, the list of CCs at t, after the surface filtering, are

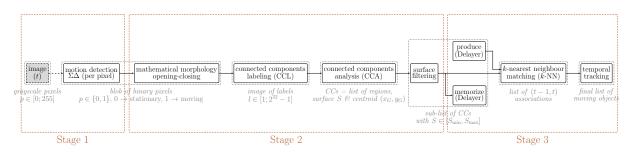


Figure 3: MOTION simplified graph. In gray and italic: the output of each processing. The graph is split into a 3-stage pipeline.

memorized with a Delayer. This way, in the next stream at t + 1, they can be reused (produce task of the Delayer) to perform the k-NN algorithm.

The produce and memorize tasks are stateful and by definition they have a data dependency over time. Thus, they cannot be replicated. Fig. 3 gives you the new stages decomposition.

Work to do #1 Implement this simplified tasks graph version in a new motion\_aff3ct.cpp file. This will produce the motion-aff3ct executable. Verify that this new version is working by comparing its logs output with the motion2 executable logs. Then, report the achieved FPS in the motion\_aff3ct.cpp file. It is better? How much and why?

Work to do #2 The k-NN task could have been replicated and a 5-stage pipeline could have been implemented. Why this does not sound like a good idea? Comment your answer in the motion\_aff3ct.cpp file.

### 2.4 Data Parallelism with OpenMP

Now, the limiting stage should be the stage 1 because of the  $\Sigma\Delta$  task. This task cannot be replicated because of its data dependency over time. However, for one given frame, each pixel is computed independently of the others! This means that the loop over the pixels can be split in multiple sub-domains and each sub-domain can be affected to a thread.

Work to do #1 Write a multi-threaded version the code of the  $\Sigma\Delta$  algorithm (sigma\_delta\_compute function). For this you will use the OpenMP library.

Now you will have 3 different types of parallelism in MOTION: the pipeline, the replication and the data parallelism inside the  $\Sigma\Delta$  task (with OpenMP). Remember that you can control the number of OpenMP threads with the OMP\_NUM\_THREADS environment variable.

Work to do #2 Find a configuration of threads that gives the best throughput. You can change the number of OpenMP threads and the number of replications in stage 2 of the pipeline. Report this configuration and the achieved throughput in the motion\_aff3ct.cpp file.

### 2.5 Task Vectorization

An other way to speedup the  $\Sigma\Delta$  task is to vectorize it with the SIMD instructions. As each pixel computation is independent, the  $\Sigma\Delta$  task is a very good candidate for vectorization.

Work to do #1 Write a vectorized version the  $\Sigma\Delta$  algorithm (inside the sigma\_delta\_compute function). For this you will use the MIPP wrapper (do not forget to add the #include <mipp.h> header).

Now you can combined the 3 previous types of parallelism with a new one: the vectorization of  $\Sigma\Delta$ .

Work to do #2 Find a configuration of threads that gives the best throughput (with the  $\Sigma\Delta$  vectorized task). Report this configuration and the achieved throughput in the motion\_aff3ct.cpp file.

### 2.6 [Bonus] Data Parallelism versus Pipeline+Replication

The most compute intensive tasks of the MOTION application are the  $\Sigma\Delta$  and the morphology. We saw before that  $\Sigma\Delta$  can easily be speedup over its data parallelism. In fact, this is the same for the morphology algorithms!

Work to do #1 First you will create a new motion.c file where you will implement the simplified graph given in Fig. 3. It will produced a new motion executable without AFF3CT (no pipeline, no replication). You can start from a copy-paste of the motion2.c file.

Work to do #2 Parallelize the morphology tasks with OpenMP. Compare the throughput of the motion executable (only data parallelism) with the throughput of the motion-aff3ct executable (data, pipeline and replication parallelisms). Which one gives the highest throughput? Why? Report the achieved results in the corresponding source files.

**Final note** We could also have vectorized the morphology task, it is combination of stencils like the **blur** kernel we worked on in previous sessions... But we will keep this for an other time ;-).