Mali™-T600 Series GPU OpenCL
Version 1.1.0
Developer Guide
Mali-T600 Series GPU OpenCL
Developer Guide

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Release Information

The following changes have been made to this book.

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Product Status

The information in this document is final, that is for a developed product.

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Preface

This preface introduces the Mali-T600 Series GPU OpenCL Developer Guide. It contains the following sections:

- *About this book* on page vi.
- *Feedback* on page viii.
About this book

This book is for the Mali-T600 Linux OpenCL Software Development Kit (SDK).

Product revision status

The rnpm identifier indicates the revision status of the product described in this book, where:

- r: Identifies the major revision of the product.
- n: Identifies the minor revision or modification status of the product.

Intended audience

This guide is written for software developers with experience in C or C-like languages who want to develop OpenCL applications for Mali-T600 Series GPUs.

Using this book

This book is organized into the following chapters:

Chapter 1 Introduction
Read this for an introduction to OpenCL and the Mali-T600 Series OpenCL SDK.

Chapter 2 Parallel Processing Concepts
Read this for an introduction to parallel processing concepts and how these work in OpenCL.

Chapter 3 OpenCL Concepts
Read this for a description of the OpenCL concepts.

Chapter 4 Stages in an OpenCL Program
Read this for a description of the stages in an OpenCL program.

Chapter 5 Converting Existing Code to OpenCL
Read this for a description of how to convert existing code to OpenCL.

Chapter 6 Retuning Existing OpenCL Code for Mali GPUs
Read this for a description of how to retune existing OpenCL code for the Mali-T600 Series GPUs.

Chapter 7 Optimizing OpenCL for Mali GPUs
Read this for a description of how to optimize OpenCL for the Mali-T600 Series GPUs.

Chapter 8 The Mali OpenCL SDK
Read this for an introduction to the Mali OpenCL SDK.

Appendix A OpenCL Data Types
Read this for a description of the OpenCL memory model and the data types available.

Appendix B OpenCL Built-in Functions
Read this for a list of the OpenCL built-in functions implemented in the Mali-T600 Series Linux OpenCL driver.
Appendix C OpenCL Extensions

Read this for a list of extensions the Mali-T600 Series Linux OpenCL driver supports.

Glossary

The ARM Glossary is a list of terms used in ARM documentation, together with definitions for those terms. The ARM Glossary does not contain terms that are industry standard unless the ARM meaning differs from the generally accepted meaning.


Typographical conventions

This book uses the following typographical conventions:

- **italic**: Introduces special terminology, denotes cross-references, and citations.
- **bold**: Highlights interface elements, such as menu names. Denotes signal names. Also used for terms in descriptive lists, where appropriate.
- **monospace**: Denotes text that you can enter at the keyboard, such as commands, file and program names, and source code.
- **mono space**: Denotes a permitted abbreviation for a command or option. You can enter the underlined text instead of the full command or option name.
- **monospace italic**: Denotes arguments to monospace text where the argument is to be replaced by a specific value.
- **monospace bold**: Denotes language keywords when used outside example code.
- `< and >`: Enclose replaceable terms for assembler syntax where they appear in code or code fragments. For example:

  ```
  MRC p15, 0 <Rd>, <CRn>, <CRm>, <Opcode_2>
  ```

Additional reading

This section lists publications by ARM and by third parties.


Other publications

This section lists relevant documents published by third parties:

- OpenCL 1.1 Specification, www.khronos.org
Feedback

ARM welcomes feedback on this product and its documentation.

Feedback on this product

If you have any comments or suggestions about this product, contact your supplier and give:
• The product name.
• The product revision or version.
• An explanation with as much information as you can provide. Include symptoms and diagnostic procedures if appropriate.

Feedback on content

If you have comments on content then send an e-mail to errata@arm.com. Give:
• The title.
• The number, DUI0538E.
• The page numbers to which your comments apply.
• A concise explanation of your comments.

ARM also welcomes general suggestions for additions and improvements.

Note

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Chapter 1
Introduction

This chapter introduces GPU compute, OpenCL, the Mali-T600 Series Linux OpenCL driver, and the Mali OpenCL SDK. It contains the following sections:

- About GPU compute on page 1-2.
- About OpenCL on page 1-3.
- About the Mali-T600 Series Linux OpenCL driver on page 1-4.
- About the Mali-T600 Series OpenCL SDK on page 1-5.
1.1 **About GPU compute**

GPU compute, or *General Purpose computing on Graphics Processing Units* (GPGPU), is the practice of using the parallel computing power of a GPU for tasks other than 3D graphics rendering.

Application processors are designed to execute a single thread as quickly as possible. This sort of processing typically includes scalar operations and control code.

GPUs are designed to execute many threads at the same time. They run compute intensive data processing tasks in parallel that contain relatively little control code. GPUs typically contain many more processing elements than application processors so can compute at a much higher rate than application processors.

OpenCL is the first open standard language to enable developers to run general purpose computing tasks on GPUs, application processors, and other types of processors.
1.2 About OpenCL

The Open Computing Language (OpenCL) is an open standard for writing programs to run on heterogeneous multi-processor systems. OpenCL provides a single programming environment for programs that can run on different processors.

OpenCL includes a platform-independent C99-based language for writing functions called kernels that execute on OpenCL devices, and APIs that define and control the platforms.

OpenCL enables you to execute some programs faster by moving intensive data processing routines to the GPU instead of the application processor.

OpenCL is an open standard developed by the Khronos Group, http://www.khronos.org.
1.3 About the Mali-T600 Series Linux OpenCL driver

The Mali-T600 Series Linux OpenCL driver is the implementation of OpenCL for the Mali-T600 Series GPUs. In this document, it is known as the Mali OpenCL driver.

The Mali OpenCL driver:

• Supports OpenCL version 1.1, Full profile.

• Is binary-compatible with OpenCL 1.0 programs. This includes compatibility with the APIs deprecated in OpenCL 1.1.

Note

The Mali OpenCL driver is for the Mali-T600 Series GPUs. It does not support the Mali-300, Mali-400, or Mali-450 GPUs.
1.4 About the Mali-T600 Series OpenCL SDK

The Mali-T600 Series OpenCL SDK contains code examples and tutorials to help you understand OpenCL development.

See Chapter 8 The Mali OpenCL SDK.
Parallel processing is the processing of computations on multiple processors simultaneously. OpenCL enables applications to use hardware resources such as GPUs to accelerate computations with parallel processing.

This chapter introduces the main concepts of parallel processing. It contains the following sections:

- *Types of parallelism* on page 2-2.
- *Concurrency* on page 2-4.
- *Limitations of parallel processing* on page 2-5.
- *Embarrassingly parallel applications* on page 2-6.
- *Mixing different types of parallelism* on page 2-7.
2.1 Types of parallelism

There are the following types of parallelism:

Data parallelism

In a data-parallel application, data is divided into data elements that can be processed in parallel. Multiple data elements are read and processed simultaneously by different processors.

The data must be in data structures that can be read from and written to in parallel. An example of a data parallel application is rendering three dimensional graphics. The generated pixels are independent so the computations required to generate them can be performed in parallel. This sort of parallelism is very fine grained and there can be hundreds or thousands of threads active simultaneously.

OpenCL is primarily used for data parallel processing.

Task parallelism

Task parallelism is where the application is broken up into tasks and these tasks are executed in parallel. Task parallelism is also known as functional parallelism.

An example of an application that can use task parallelism is playing an on-line video. To display a web page your device must do several tasks:

- Run a network stack that performs communication.
- Request data from external server.
- Read data from external server.
- Parse data.
- Decode video data.
- Decode audio data.
- Draw video frames.
- Play audio data.

Figure 2-1 shows parts of an application and operating system that operate simultaneously when playing an on-line video.

Pipelines

Pipelines process data in a series of stages. In a pipeline the stages can operate simultaneously but they do not process the same data. A pipeline typically has a relatively small number of stages.

An example of a pipeline is a video recorder application that must perform the following stages:

1. Capture image data from an image sensor and measure light levels.
2. Modify the image data to correct for lens effects.
3. Modify the contrast, color balance, and exposure of the image data.
4. Compress the image
5. Add the data to video file.
6. Write the video file to storage.

These stages must be performed in order but they can be all be operating on data from a different video frame at the same time.

Figure 2-2 shows parts of an application that can operate simultaneously as a pipeline playing a video from the internet.

![Figure 2-2 Pipeline processing](image-url)
2.2 Concurrency

Concurrent applications have multiple operations in progress at the same time. These can operate in parallel or in serial using a time sharing system.

In a concurrent application multiple tasks are typically trying to share the same data. Access to this data must be managed carefully otherwise there can be complex problems such as:

**Race conditions**

A race condition occurs when two or more threads try to modify the value of one variable at the same time. The final value of the variable should always be the same, but when a race condition occurs the variable can get a different value depending on the order of the writes.

**Deadlocks**

A deadlock occurs when two threads become blocked by each other and neither thread can progress with their operations. This can happen when the threads each obtain a lock that the other thread requires.

**Live locks**

Live locks are similar to deadlocks but the threads keep running. However, because of the lock the threads can never complete their task.

A concurrent data structure is a data structure that can be accessed by multiple tasks without causing concurrency problems.

Data parallel applications use concurrent data structures. These are the sorts of data structures that you typically use in OpenCL.
2.3 Limitations of parallel processing

There are limitations of parallel processing that you must consider when developing parallel applications.

For example, if your application parallelizes perfectly, executing the application on 10 processors makes it run 10 times faster.

Applications rarely parallelize perfectly because part of the application is serial. This serial component imposes a limit on the amount of parallelization the application can use.

Amdahl’s law describes the speedup you can get from parallel processing. The formula for Amdahl’s law is shown in Figure 2-3 where the terms in the equation are:

- $S$: Fraction of the application that is serial.
- $P$: Fraction of the application that is parallelizable.
- $N$: Number of processors.

\[
\text{Speedup} = \frac{1}{S + \frac{P}{N}}
\]

Figure 2-3 Formula for Amdahl’s law

Figure 2-4 shows the speedup that different numbers of processors provide for applications with different serial components.

The biggest speedups are achieved with relatively small numbers of processors. However, as the number of processors rises the gains reduce.

You cannot avoid Amdahl’s law in your application but you can reduce the impact. See Reducing the effect of serial computations on page 7-8.

For high performance with a large number of processors the application must have a very small serial component. These sorts of applications are said to be embarrassingly parallel. See Embarrassingly parallel applications on page 2-6.
2.4 Embarrassingly parallel applications

If an application can be parallelized across a large number of processors easily, it is said to be embarrassingly parallel.

An example of an embarrassingly parallel application is rendering three dimensional graphics. The pixels are completely independent so they can be computed and drawn in parallel.

OpenCL is ideally suited for developing and executing embarrassingly parallel applications.

Figure 2-5 shows an image divided into many small parts. These parts can all be processed simultaneously.
2.5 Mixing different types of parallelism

You can mix these types of parallelism in your applications. For example, an audio synthesizer might use a combination of all three types of parallelism:

- Task parallelism computes the notes independently.
- A pipeline of audio generation and processing modules creates the sound of an individual note.
- Within the pipeline some stages can use data parallelism to accelerate the computation of processing.
Chapter 3
OpenCL Concepts

This chapter describes the OpenCL concepts. It contains the following sections:

• *About OpenCL* on page 3-2.
• *OpenCL applications* on page 3-3.
• *OpenCL execution model* on page 3-4.
• *OpenCL data processing* on page 3-5.
• *The OpenCL memory model* on page 3-8.
• *The Mali GPU memory model* on page 3-9.
• *Summary* on page 3-10.
3.1 About OpenCL

OpenCL is an open standard that enables you to use the parallel processing capabilities of multiple types of processors including application processors, GPUs, and other computing devices.

OpenCL specifies an API for parallel programming that is designed for portability:
• It uses an abstracted memory and execution model.
• There is no requirement to know the application processor or GPU instruction set.
• There is scope for specific hardware optimizations.

Functions executing on OpenCL devices are called kernels. These are written in a language called OpenCL C that is based on C99.

The Mali-T600 Series GPUs supports OpenCL 1.1 Full Profile.
3.2 OpenCL applications

OpenCL applications consist of two parts:

**Application, or host, side code**
- Calls the OpenCL APIs.
- Compiles the CL kernels.
- Allocates memory buffers to pass data into and out of the OpenCL kernels.
- Sets up command queues.
- Sets up dependencies between the tasks.
- Sets up the NDRanges that the kernels execute over.

**OpenCL kernels**
- Written in OpenCL C language.
- Perform the parallel processing.
- Runs on the compute devices such as GPU shader cores.

You must write both of these parts correctly to get the best performance.
3.3 OpenCL execution model

The OpenCL execution model includes:
• Kernels that run on compute devices.
• A host program that runs on the application processor.

The host program

The host program manages the execution of the kernels by setting up command queues for:
• Memory commands.
• Kernel execution commands.
• Synchronization.

The context

The host program defines the context for the kernels. The context includes:
• The kernels.
• Compute devices.
• Programs objects.
• Memory objects.

Operation of OpenCL kernels

A kernel is a block of code that is executed on a compute device in parallel with other kernels. Kernels operate in the following sequence:
1. A kernel is defined in a host application.
2. The host application submits the kernel for execution on a compute device.
   A compute device can be an application processor, GPU, or another type of processor.
3. When the application issues a command to submit a kernel, OpenCL creates the NDRange of work-items.
4. An instance of the kernel is created for each element in the NDRange. This enables each element to be processed independently in parallel.
3.4 OpenCL data processing

This section describes OpenCL data processing. It contains the following sections:

- Work-items and the NDRange.
- OpenCL work-groups on page 3-6.
- Identifiers in OpenCL on page 3-7.

3.4.1 Work-items and the NDRange

The data processed by OpenCL is in an index space of work-items. The work-items are organized in an N-Dimensional Range (NDRange) where:

- $N$ is the number of dimensions minus one.
- $N$ can be zero, one, or two.

One kernel instance is executed for each work-item in the index space.

Figure 3-1 shows NDranges with one, two and three dimensions.

![Figure 3-1 NDranges and work-items](image)

You group work-items into work-groups for processing. Figure 3-2 on page 3-6 shows a three dimensional NDRange that is split into 16 work-groups each with 16 work-items.
3.4.2 OpenCL work-groups

Work-groups have a number of properties and limitations:

**Properties of work-groups**
- Work-groups are independent of each other.
- You can issue multiple work-groups for execution in parallel.
- The work-items in a work-group can communicate with each other using shared data buffers, you must synchronize access to these buffers.

**Limitations of work-groups**
Work-groups typically do not directly share data. They can share data using global memory.

The following are not supported across different work-groups:
- Barriers.
- Dependencies.
- Ordering.
- Coherency.

Global atomics are available but these can be slower than local atomics.

**Work-items in a work-group**
The work-items in a work-group can do the following:
- Perform barrier operations to synchronize execution points.
  For example:
  ```
  barrier(CLK_LOCAL_MEM_FENCE); // Wait for all kernels in
  // this work-group to catch up
  ```
- Use local atomic operations.
- Access shared memory.
3.4.3 Identifiers in OpenCL

There are a number of identifiers in OpenCL:

- **global ID** Every work-item has a unique global ID that identifies it within the index space.

- **local ID** Within each work-group, each work-item has a unique local ID that identifies it within its work-group.

- **work-group ID** Each work-group has a unique work-group ID.
3.5 The OpenCL memory model

The OpenCL memory model contains a number of components. Figure 3-3 shows the OpenCL memory model.

- **Private memory**
  - Private memory is specific to a work-item.
  - It is not visible to other work-items.

- **Local memory**
  - Local to a work-group.
  - Accessible by the work-items in the work-group.
  - Consistent by all work-items in the work-group.
  - Accessed with the __local keyword.

- **Constant memory**
  - A memory region used for objects allocated and initialized by the host.
  - Accessible as read-only by all work-items.

- **Global memory**
  - Accessible to all work-items executing in a context,
  - Accessible to the host using read, write, and map commands.
  - Consistent across work-items in a single work-group.
  - Implements a relaxed consistency, shared memory model.
  - There are no guarantee of memory consistency between different work-groups.
  - Accessed with the __global keyword.
3.6 The Mali GPU memory model

The Mali GPU has a different memory model to desktop workstations:

**Desktop**  
Traditional desktop workstations have physically separate global, local and private memories.  
Typically a graphics-card has its own local memory.  
Data must be copied to the local memory and back again.

**Mali GPU**  
Mali GPUs have a unified memory system.  
Local and private memory is physically global memory.  
Moving data from global to local or private memory typically does not improve performance.  
The traditional copying of data is not required.

Each compute device, that is the shader cores, have their own data caches.
3.7 Summary

OpenCL primarily uses data parallel processing. OpenCL uses the following terminology:

- Computations in OpenCL are performed by pieces of code called *kernels* that execute on *compute devices*. Compute devices can be application processors, GPUs, or other types of processors.

- The data processed by OpenCL is in an *index space of work-items*. The work-items are organized in an NDRange.

- One kernel instance is executed for each work-item in the index space.

- Kernels execute in parallel.

- Work-items group together to form *work-groups*. The work-items in a work-group can communicate with each other using shared data buffers, but access to the buffers must be synchronised.

- Work-groups typically do not directly share data. They can share data using global memory.

- Issue multiple work-groups for execution in parallel.
Chapter 4

Stages in an OpenCL Program

This chapter describes the stages in an OpenCL program. It contains the following sections:

• Software required for OpenCL development on page 4-2.
• Development stages on page 4-3.
• Finding the available compute devices on page 4-5.
• Initializing and creating OpenCL contexts on page 4-6.
• Creating a command queue on page 4-7.
• Creating program objects on page 4-8.
• Building a program executable on page 4-9.
• Creating kernel and memory objects on page 4-10.
• Executing the kernel on page 4-11.
• Reading the results on page 4-13.
• Cleaning up on page 4-14.
4.1 Software required for OpenCL development

To develop OpenCL programs for Mali GPUs, you require:
• A platform with a Mali-T600 Series GPU.
• An implementation of OpenCL for the Mali-T600 Series GPU.

You can develop on other hardware platforms with implementations of OpenCL but you cannot use them to estimate performance on a Mali-T600 series GPU.

Implementations of OpenCL are available for the following operating systems:
• Mac OS X 10.6 and higher.
• Microsoft Windows.
• Linux.
4.2 Development stages

These are the stages for developing an OpenCL application:

1. Determine what you want to parallelize.

   The first step when deciding to use OpenCL is to look at what your program does and identify the parts of the program that can run in parallel. This is often the hardest part of developing an OpenCL program. See Analyzing code for parallelization on page 5-3.

   **Note**

   Only convert the parts of a program to OpenCL where there is likely to be a benefit. Profile your application to find the most active parts and consider these parts for conversion.

2. Write kernels.

   OpenCL programs consists of a set of kernel functions. You must write the kernels that perform the computations.

3. Find out what OpenCL devices are available.

   Query to find out what OpenCL devices are available on the system using OpenCL platform layer functions. See Finding the available compute devices on page 4-5.

4. Set up the context.

   Create and set up a OpenCL context and one or more command queues to schedule execution of your kernels. See Initializing and creating OpenCL contexts on page 4-6.

5. Write code to compile and build your OpenCL program.

   Write code in your program that includes the commands to compile and build your source code and extracts kernel objects from the compiled code. The sequence of commands you must follow is:

   a. The program object is created either by calling clCreateProgramWithSource() or clCreateProgramWithBinary(). clCreateProgramWithSource() creates the program object from the kernel source code. clCreateProgramWithBinary() creates the program with a pre-compiled binary file.

   b. Call the clBuildProgram() function to compile the program object for the specific devices on the system.

   c. Call the clCreateKernel() function for each kernel, or call the clCreateKernelsInProgram() function to create kernel objects for all the kernels in the OpenCL program.

6. Allocate memory buffers.

   Use the OpenCL API to allocate memory buffers. You can use the map() and unmap() operations to enable both the application processor and the Mali GPU to access the data.

7. Enqueue commands and kernels.

   Enqueue to the command queues the commands that control the sequence and synchronization of kernel execution, reading and writing of data, and manipulation of memory objects.

   To execute a kernel function, you must do the following:

   a. Call clSetKernelArg() for each parameter in the kernel function definition to set the kernel parameter values.

   b. Determine the work-group size and index space to use to execute the kernel.
c. Enqueue the kernel for execution in the command queue.

8. Enqueue commands that make the results from the work-items available to the host.
4.3 Finding the available compute devices

To set up OpenCL you must choose compute devices. Call clGetDeviceIDs() to query the OpenCL driver for a list of devices on the machine that support OpenCL. You can restrict your search to a particular type of device or to any combination of device types. You must also specify the maximum number of device IDs that you want returned.
4.4 Initializing and creating OpenCL contexts

After you know the available OpenCL devices on the machine and have at least one valid device ID, you can create an OpenCL context. The context groups devices together to enable memory objects to be shared across different compute devices.

Pass the device information to the `clCreateContext()` function. For example:

```c
// Create an OpenCL context

context = clCreateContext( NULL, 1, &device_id, notify_function, NULL, &err );
if (err != CL_SUCCESS)
{
    Cleanup();
    return 1;
}
```

You can optionally specify an error notification callback function when creating an OpenCL context. Leaving this parameter as a NULL value results in no error notification being registered.

Providing a callback function can be useful if you want to receive runtime errors for the particular OpenCL context. For example:

```c
// Optionally user_data can contain contextual information
// Implementation specific data of size cb, can be returned in private_info

void context_notify( const char *notify_message, const void *private_info, size_t cb, void *user_data )
{
    printf("Notification:\n\t%s\n", notify_message);
}
```
4.5 Creating a command queue

After creating your OpenCL context, use `clCreateCommandQueue()` to create a command queue. For example:

```c
// Create a command-queue on the first device available
// on the created context

commandQueue = clCreateCommandQueue( context, &device);
if (commandQueue == NULL)
{
    Cleanup();
    return 1;
}
```

If you have multiple OpenCL devices, such as an application processor and a GPU, you must:
1. Create a command queue for each device.
2. Divide up the work.
3. Submit commands separately to each device.
4.6 Creating program objects

Load the OpenCL C kernel source and create a program object from it. The program object is loaded with the kernel source code and then the code is compiled for execution on the devices attached to the context. All kernel functions must be identified in the program source with the __kernel qualifier. OpenCL programs can also include functions you can call from your kernel functions.

A program object encapsulates:
- Your OpenCL program source.
- The latest successfully built program executable.
- The build options.
- The build log.
- A list of devices the program is built for.

To create a program object use the clCreateProgramWithSource() function. For example:

```c
// Create OpenCL program

program = clCreateProgramWithSource( context, device, "<kernel source>" );
if (program == NULL)
{
    Cleanup();
    return 1;
}
```

There are different options for building OpenCL programs:

- You can create a program object directly from the source code of an OpenCL program and compile it at runtime. Do this at application startup to save compute resources while the application is running.

- To avoid compilation at runtime, you can build a program object with a previously built binary.

  ——- Note ———-

  Applications with pre-built program objects are not portable.

  ————

Creating a program object from a binary is a similar process to creating a program object from source code, except that you must supply the binary for each device that you want to execute the kernel on. Use the clCreateProgramWithBinary() function to do this.

Use the clGetProgramInfo() function to obtain the binary after you have generated it.
4.7 Building a program executable

After you have created a program object, you must build a program executable from the contents of the program object. Use the `clBuildProgram()` function to build your executable.

Compile all kernels in the program object. In the following code, there is only one kernel:

```c
err = clBuildProgram( program, 1, &device_id, "", NULL, NULL );
if (err == NULL)
{
    Cleanup();
    return 1;
}
```
4.8 Creating kernel and memory objects

This section describes creating kernel and memory objects. It contains the following sections:

- Creating kernel objects.
- Creating memory objects.

4.8.1 Creating kernel objects

Call the clCreateKernel() function to create a single kernel object, or call the clCreateKernelsInProgram() function to create kernel objects for all the kernels in the OpenCL program. For example:

```c
// Create OpenCL kernel
kernel = clCreateKernel(program, "<kernel_name>", NULL);
if (kernel == NULL)
{
    Cleanup();
    return 1;
}
```

4.8.2 Creating memory objects

After you have created and registered your kernels, send the program data to the kernels:

1. Package the data in a memory object.
2. Associate the memory object with the kernel.

There are two types of memory objects:

**Buffer objects**

Simple blocks of memory.

**Image objects**

Opaque structures, specifically for representing 2D or 3D images.

To create buffer objects, use the clCreateBuffer() function. To create image objects, use the clCreateImage2D() or clCreateImage3D() functions.
4.9 Executing the kernel

This section describes the stages in executing the kernel. It contains the following sections:

• Determining the data dimensions.
• Determining the optimal global work size.
• Determining the local work-group size.
• Enqueuing kernel execution on page 4-12.
• Executing kernels on page 4-12.

4.9.1 Determining the data dimensions

If your data is an image $x$ pixels wide by $y$ pixels high, it is a two-dimensional data set. If you are dealing with spatial data that involves the $x$, $y$, and $z$ position of nodes, it is a three-dimensional data set.

The number of dimensions in the original data set does not have to be the same in OpenCL. You can for example, process a three dimensional data set as a single dimensional data set in OpenCL.

4.9.2 Determining the optimal global work size

The global work size is the total number of work-items required for all dimensions combined.

You can change the global work size by processing multiple data items in a single work-item. The new global worksize is then the original global work size is divided by the number of data items processed by each work-item.

Use the formula in Table 4-1 to work out the global work size for data with different dimensions where $n$ is the number of data elements processed in a single work-item.

<table>
<thead>
<tr>
<th>Dimensions</th>
<th>Formula</th>
</tr>
</thead>
<tbody>
<tr>
<td>One</td>
<td>$x / n$</td>
</tr>
<tr>
<td>Two</td>
<td>$(x \times y) / n$</td>
</tr>
<tr>
<td>Three</td>
<td>$(x \times y \times z) / n$</td>
</tr>
</tbody>
</table>

The optimal global work size must be large if you want to ensure high performance. Typically the number is several thousand but the ideal number depends on the number of shader cores in your device.

To calculate the optimal global work size use the following equation:

$$\text{global work size} = \text{<maximum_work-group_size>} \times \text{<number of shader cores>} \times \text{<constant>}$$

Where constant is typically 4 or 8 for the Mali-T604 GPU.

4.9.3 Determining the local work-group size

You can specify the size of the work-group that OpenCL uses when you enqueue a kernel to execute on a device. To do this, you must know the maximum work-group size permitted by the OpenCL device your work-items execute on. To find the maximum work-group size, use the clGetKernelWorkGroupInfo() function and request the CL_KERNEL_WORK_GROUP_SIZE property.
If your application is not required to share data among work-items, set the local_work_size parameter to NULL when enqueuing your kernel. This enables the OpenCL driver to determine the most efficient work-group size for your kernel.

To get the maximum work-group size in each dimension, call clGetDeviceInfo() with CL_DEVICE_MAX_WORK_ITEM_SIZES. To get the total work-group size call clGetKernelWorkGroupInfo() with CL_KERNEL_WORK_GROUP_SIZE. If the maximum work-group size for a kernel is lower than 128, try simplifying the kernel.

The work-group size for each dimension must divide evenly into the total data-size for that dimension. That is, the x size of the work-group must divide evenly into the x size of the total data. If this requirement means padding the work-group with extra work-items, ensure the additional work-items return immediately and do no work.

4.9.4 Enqueuing kernel execution

When you have identified the dimensions necessary to represent your data, the necessary work-items for each dimension, and an appropriate work-group size, enqueue the kernel for execution using clEnqueueNDRangeKernel(). For example:

```c
size_t globalWorkSize[1] = { ARRAY_SIZE };
size_t localWorkSize[1] = { 4 } ;

// Queue the kernel up for execution across the array
errNum = clEnqueueNDRangeKernel(commandQueue, kernel, 1, NULL, globalWorkSize, localWorkSize, 0, NULL, NULL);
if (errNum != CL_SUCCESS)
{
    printf( "Error queuing kernel for execution.\n" );
    Cleanup();
    return 1;
}
```

4.9.5 Executing kernels

Queuing the kernel for execution does not mean that it executes immediately. The kernel execution is put into the command queue for later processing by the device. After making the call to clEnqueueNDRangeKernel() the kernel might not yet have executed on the device.

It is possible to make a kernel wait for execution until previous events are finished. You can specify certain kernels wait until other specific kernels are completed before executing.
4.10 Reading the results

After your kernels have finished execution, you must make the result accessible to the host.

To access the results from the kernel, use `clEnqueueMapBuffer()` to map the buffer into host memory. For example:

```c
local_buffer = clEnqueueMapBuffer( queue, buffer, CL_FALSE, CL_MAP_READ, 0,
        ( sizeof(unsigned char) * buffer_size), num_deps, deps[1], NULL, &err );

ASSERT( CL_SUCCESS == err );
```

**Note**

This call does not guarantee to make the buffer available until you call `clFinish()`. If you change the third parameter `CL_FALSE` to `CL_TRUE`, the call becomes a blocking call and it is executed immediately.
4.11 Cleaning up

When the program no longer requires the various objects associated with the OpenCL runtime and context, you must free these resources. Use the following functions to release your OpenCL objects. These functions decrement the reference count for the associated object:

- clReleaseMemObject().
- clReleaseKernel().
- clReleaseProgram().
- clReleaseCommandQueue().
- clReleaseContext().

Ensure the reference counts for all OpenCL objects reach zero when your program no longer requires them. You can obtain the reference count by querying the object. For example, by calling clGetMemObjectInfo().
Chapter 5
Converting Existing Code to OpenCL

This section describes converting existing code to OpenCL. It contains the following sections:

• Profile your application on page 5-2.
• Analyzing code for parallelization on page 5-3.
• Parallel Processing Techniques on page 5-5.
• Using parallel processing with non-parallelizable code on page 5-10.
• Dividing data for OpenCL on page 5-11.
5.1 Profile your application

Profile your application to find the most compute intensive parts. These are the parts that might be worth porting to OpenCL.

The proportion of an application that requires high performance is typically a relatively small part of the code. This is the part of the code that can probably make best use of OpenCL. Porting any more of the application to OpenCL is unlikely to provide a benefit.

You can use profilers such as DS-5™ to profile your application. You can download DS-5 from the Mali developer web site, http://www.malideveloper.arm.com
5.2 Analyzing code for parallelization

This section describes how to analyze compute intensive code for parallelization. It contains the following sections:

- About analyzing code for parallelization.
- Look for data parallel operations.
- Look for operations with few dependencies on page 5-4
- Analyze loops on page 5-4.

5.2.1 About analyzing code for parallelization

When you have identified the most compute intensive parts of your application, analyze the code to see if you can run it in parallel.

Parallelizing code can be the following:

**Easy**

Parallelizing the code requires little or no modification.

**Straight forward**

Parallelizing the code requires small modifications. See Use the global ID instead of the loop counter on page 5-5.

**Difficult**

Parallelizing the code requires complex modifications. See Compute values in a loop with a formula instead of using counters on page 5-6.

**Difficult and includes dependencies**

Parallelizing the code requires complex modifications and the use of special techniques to avoid dependencies. See the following sections:

- Compute values per frame on page 5-6.
- Perform computations with dependencies in multiple-passes on page 5-7.
- Pre-compute values to remove dependencies on page 5-8.

**Apparently impossible**

If parallelizing the code appears to be impossible, investigate parallel alternatives to the algorithms and data structures the code uses. These might make parallelization possible. See Using parallel processing with non-parallelizable code on page 5-10.

**Impossible**

This only means the implementation cannot be parallelized. Do not think that code is the definitive solution to a problem. The code is only one possible implementation of a solution. There might be multiple different solutions and some of these might be parallelizable. See Using parallel processing with non-parallelizable code on page 5-10.

5.2.2 Look for data parallel operations

Look for tasks that do large numbers of operations that:

- Complete without sharing data.
- Do not depend on the results from each other.

These types of operations are data parallel so are ideal for OpenCL.
5.2.3 Look for operations with few dependencies

If tasks have few dependencies, it might be possible to run them in parallel.

Dependencies between tasks prevent parallelization because it forces tasks to be performed sequentially. If the code has dependencies consider:

• Is there a way to remove the dependencies?
• Can you delay the dependencies to later in execution?

5.2.4 Analyze loops

Loops are good targets for parallelization because they repeat computations many times, often independently.

**Loops that process a small number of elements**

If the loop only processes a relatively small number of elements it might not be appropriate for data parallel processing. It might be better to parallelize these sorts of loops with task parallelism on one or more application processors.

**Perfect loops**

Look for loops that:

• Process thousands of items.
• Have no dependencies on previous iterations.
• Accesses data independently in each iteration.

These types of loops are data parallel so are ideal for OpenCL.

**Simple loop parallelization**

If the loop includes a variable that is incremented based on a value from the previous iteration, this is a dependency between iterations that prevents parallelization.

See if you can work out a formula that enables you to compute the value of the variable based on the main loop counter.

In OpenCL kernels process work-items parallel. There is no loop counter to reference because work-items are not processed in a loop.

Every work-item has a unique global id that identifies it. You can use this value in place of a loop counter. See *Use the global ID instead of the loop counter on page 5-5*.

**Loop requires data from previous iteration**

If your loop involves dependencies based on data processed by a previous iteration this is a more complex problem.

Can the loop be restructured to remove the dependency? If not, it might not be possible to parallelize the loop.

There are a number of techniques to divide the dependencies apart. See if you can use these techniques to parallelize the loop. See *Parallel Processing Techniques on page 5-5* for a description of some of these techniques.

**Non-parallelizable loops**

If the loop contains dependencies that you cannot remove, investigate alternative methods of performing the computation. These might be parallelizable.

See *Using parallel processing with non-parallelizable code on page 5-10*. 
5.3 Parallel Processing Techniques

This section describes parallel processing techniques you can use in OpenCL. It contains the following sections:

- Use the global ID instead of the loop counter.
- Compute values in a loop with a formula instead of using counters on page 5-6.
- Compute values per frame on page 5-6.
- Perform computations with dependencies in multiple-passes on page 5-7.
- Pre-compute values to remove dependencies on page 5-8.
- Use software pipelining on page 5-8.
- Use task parallelism on page 5-9.

5.3.1 Use the global ID instead of the loop counter

In OpenCL you use kernels to perform the equivalent of loop iterations. This means there is no loop counter to use in computations.

The global ID of the work-item provides the equivalent of the loop counter. Use the global ID to perform any computations based on the loop counter.

Note
You can include loops in OpenCL kernels but they can only iterate over the data for that work-item, not the entire NDRange.

The following example shows a simple loop in C that assigns the value of the loop counter to each array element.

Loop example in C:

The following loop fills an array with numbers.

```c
void SetElements(void)
{
  int loop_count;
  int my_array[4096];

  for (loop_count = 0; loop_count < 4096; loop_count++)
  {
    my_array[loop_count] = loop_count;
  }

  printf("Total %d\n", loop_count);
}
```

This loop is parallelizable because the loop elements are all independent. There is no main loop counter `loop_count` in the OpenCL kernel so it is replaced by the global ID.

The equivalent code in an OpenCL kernel:

```c
__kernel void example(__global int * restrict my_array)
{
  int id;
  id = get_global_id(0);
  my_array[id] = id;
}
```
5.3.2 **Compute values in a loop with a formula instead of using counters**

OpenCL kernels perform the equivalent of loop iterations so it is not possible to increment or decrement variables based on loop iterations.

In place of incrementing and decrementing variables, devise a formula that computes the value of the variable based on the value of the global ID. The global ID of the work-item provides the equivalent of the loop counter.

5.3.3 **Compute values per frame**

If your application requires continuous updates of data elements and there are dependencies between them, try breaking the computations into discrete units and perform one iteration per image frame displayed.

For example, the image shown in Figure 5-1 is of an application that runs a continuous physics simulation of a flag.

![Figure 5-1 Flag simulation](image_url)

The flag is made up of a grid of nodes that are connected to the neighboring nodes. These are shown in Figure 5-2 on page 5-7.
The simulation runs as a series of iterations. In one iteration all the nodes are updated and the image is redrawn.

The following operations are performed in each iteration:
1. The node values are read from a buffer A.
2. A physics simulation computes the forces between the nodes.
3. The position and forces on the nodes are updated and stored into buffer B.
4. The flag image is drawn.
5. Buffer A and buffer B are switched.

In this case splitting the computations into iterations also splits the dependencies. The data required for one frame is computed in the previous frame.

Some types of simulation require many iterations for relatively small movements. If this is the case try computing multiple iterations before drawing frames.

5.3.4 Perform computations with dependencies in multiple-passes

If your application requires continuous updates of data elements and there are dependencies between them, try breaking the computations into discrete units and perform the computations in multiple stages.

This technique extends the technique described in Compute values per frame on page 5-6 by breaking up computations more.

Divide the data elements into odd and even fields. This divides the dependencies so the entire computation can be performed in stages. The processing alternates between computing the odd then the even fields.

For example, this technique can be used in neural network simulation.

The individual neurons are arranged in a three dimensional grid. Computing the state for a neuron involves reading inputs from the surrounding neurons. This means each neuron has dependencies on the state of the surrounding neurons.

To execute the simulation, the three dimensional grid is divided into layers and executed in the following manner:
1. The even node values are read.
2. The odd layers are computed and the results stored.
3. The odd node values are read.
4. The even layers are computed and the results stored.

5.3.5 **Pre-compute values to remove dependencies**

If part of your computation is serial, see if it can be removed and performed separately.

For example, The audio synthesis technique *Frequency Modulation* (FM) works by reading an audio waveform called the carrier. The rate the waveform is read at is dependent on another waveform called the modulator.

The carrier values are read by a pointer to generate the output waveform. The position of the pointer is computed by taking the previous value and moving it by an amount determined by the value of the modulator waveform.

The position of the pointer has a dependency on the previous value and that value has a dependency on the value before it. This series of dependencies makes the algorithm difficult or impossible to parallelize.

Another approach is to consider that the pointer is moving through the carrier waveform at a fixed speed and the modulator is adding or subtracting an offset. This can be computed in parallel but the offsets are incorrect because they do not take account of the dependencies on previous offsets.

The computation of the correct offsets is a serial process. If you pre-compute these values the remaining computation can be parallelized. The parallel component reads from the generated offset table and uses this to read the correct value from the carrier waveform.

There is a potential problem with this example. The offset table must be re-computed every time the modulating waveform changes. This is an example of Amdahl’s law. The amount of parallel computation possible is limited by the speed of the serial computation.

5.3.6 **Use software pipelining**

Software pipelines are a parallel processing technique that enable multiple data elements to be processed simultaneously by breaking the computation into a series of sequential stages.

Pipelines are common in both hardware and software. For example, application processors and GPUs use hardware pipelines. The graphics standards OpenGL ES is based on a virtual pipeline.

In a pipeline a complete process is divided into a series of stages. A data element is processed in a stage and the results are then passed to the next stage.

Because of the sequential nature of a pipeline only one stage is used at a time by a particular data element. This means the other stages can process other data elements.

You can use software pipelines in your application to process different data elements.

For example, a game requires many different operations to happen. A game might use a similar pipeline to this:

1. Input read from player.
2. Game logic computes the progress of the game.
3. Scene objects moved based on the results of the game logic.
4. Physics engine computes positions of all objects in the scene.
5. Game uses OpenGL ES to draw objects on screen.
5.3.7 Use task parallelism

Task or functional parallelism involves breaking an application up by function into different tasks.

For example, an online game can take advantage of task parallelism. To run an online game your device performs several functions:

- Communicate with an external server.
- Read player input.
- Update the game state.
- Generate sound effects.
- Play music.
- Update the display.

These tasks require synchronisation but are otherwise largely independent operations. This means you can execute the tasks in parallel on separate processors.

Another example of task parallelism is Digital Television (DTV). At any time the television might be performing several of the following operations:

- Downloading program.
- Recording program.
- Updating program guide.
- Displaying options.
- Reading from media storage device.
- Playing program.
- Decoding video stream.
- Playing audio.
- Scaling image to correct size.
5.4 Using parallel processing with non-parallelizable code

If you cannot parallelize your code there is still the possibility that you can use parallel processing.

Most code is written to run on application processors that run sequentially. The code uses serial algorithms and non-concurrent data structures. Parallelizing this sort of code can be difficult or impossible.

The fact the code cannot be parallelized only means this specific implementation cannot be parallelized. It does not mean the problem cannot be solved in a parallel way.

Investigate the following approaches:

Use parallel versions of your data structures and algorithms

Many common data structures and algorithms that use them are non-concurrent. This prevents you from parallelizing the code.

There are parallel versions of many common data structures and algorithms. You might be able to use these in place of the originals to parallelize the code.

See Use concurrent data structures on page 5-11.

Solve the problem in a different way

Take a step back and think about what problem the code solves.

Look at the problem and investigate alternative ways of solving it. There might be alternative solutions that use algorithms and data structures that are parallelizable.

To do this think in terms of the purpose of the code and data structures.

Typically the aim of code is to process or transform data. It takes a certain input and produces a certain output.

• Can the data you want to process be broken up into small data elements?
• Can these data elements be placed into a concurrent data structure?
• Can you process the data elements independently?

If the answer to these are yes, then you can probably solve your problem with OpenCL.
5.5 Dividing data for OpenCL

This section describes dividing data for processing with OpenCL. It contains the following sections:

- About dividing data for OpenCL.
- Use concurrent data structures.
- Data division examples.

5.5.1 About dividing data for OpenCL

Data is divided up so it can be computed in parallel with OpenCL. The data is divided into the following hierarchy of levels:

- At the highest level the data is divided into an NDRange. The total number of elements in the NDRange is known as the global work size.
- The NDRange is divided into work-groups.
- Each work-group is divided into work-items.

See Chapter 3 OpenCL Concepts.

5.5.2 Use concurrent data structures

OpenCL executes hundreds or thousands of individual kernel instances so the processing and data structures must be parallelizable to that degree.

This means you must use data structures that permit multiple data elements to be read and written simultaneously and independently. These are known as concurrent data structures.

Many common data structures are non-concurrent. This prevents parallelizing the code. For example, the following data structures are non-concurrent:

- Linked list.
- Hash table.
- Btree.
- Map.

There are parallel versions of many commonly used data structures.

5.5.3 Data division examples

The following are examples of data in different dimensions that you can process with OpenCL:

Note These examples map the problems into the NDRanges with the same number of dimensions. OpenCL does not require that you do this. You can map a one-dimensional problem onto a two or three-dimensional NDRange.

One dimensional data

An example of one dimensional data is audio. Audio is represented as a series of samples. Changing the volume of the audio is a parallel task because the operation is performed independently per sample.

In this case the NDRange is the total number of samples in the audio. Each work-item can be one sample and a work-group is a collection of samples.
Audio can also be processed with vectors. If your audio samples are 16-bit, you can make a work-item represent 8 samples and process 8 of them at a time with vector instructions.

**Two dimensional data**

An image is a natural fit for OpenCL because it is a two dimensional array of pixels. You can process a 1600 by 1200 pixel image by mapping it onto an NDRange of 1600 by 1200.

The total number of work-items is the total number of pixels in the image, that is, 1920000.

The NDRange is divided into work-groups where each work-group is also a two dimensional array. The number of work-groups must divide into the NDRange exactly.

If each work-item processes a single pixel, a work-group size of 8 by 16 has the size of 128. This work-group size fits exactly into the NDRange on both the x and y axis. To process the image you require 15000 work-groups of 128 work-items each.

You can vectorize this example by processing all the color channels in a single vector. If the channels are 8-bit values you can process multiple pixels in a single vector. If each vector processes 4 pixels, this means each work-item processes 4 pixels and you require 4 time fewer work-items to process the entire image. This means your NDRange can be reduced to 400 by 1200 and you only require 3750 work-groups to process the image.

**Three dimensional data**

You can use three dimensional data to model the behavior of materials in the real world. For example, you can model the behavior of concrete for building by simulating the stresses in a three dimensional data set. You can use the data produced to determine the size and design of the concrete you require to hold a specific load.

You can use this technique in games to model the physics of objects. When an object is broken the physics simulation makes the process of breaking more realistic.
Chapter 6
Retuning Existing OpenCL Code for Mali GPUs

This chapter describes how to optimize existing OpenCL code for Mali GPUs. It contains the following sections:

• About optimizing existing OpenCL code for Mali GPUs on page 6-2.
• Procedure for optimizing existing OpenCL code for Mali GPUs on page 6-3.
6.1 About optimizing existing OpenCL code for Mali GPUs

OpenCL is a portable language but it is not always performance portable. This means that OpenCL can work on many different types of compute device but performance is not preserved.

Existing OpenCL is typically tuned for specific architectures such as desktop GPUs. To achieve better performance on Mali GPUs you must retune the code for the Mali GPUs.

The procedure to convert OpenCL code to run optimally on Mali GPUs is:

1. Analyze the code.
2. Locate and remove optimizations for alternative compute devices.
3. Vectorize the code.
4. Optimize the code for the Mali GPU.
6.2 Procedure for optimizing existing OpenCL code for Mali GPUs

This section describes the procedure for optimizing existing OpenCL code for Mali GPUs. It contains the following sections:

- Analyze code.
- Locate and remove device optimizations.
- Optimizing your OpenCL code for Mali GPUs on page 6-4.

6.2.1 Analyze code

If you did not write the code yourself, you must analyze it to find out exactly what it does.

Try to understand the following:

- What is the purpose of the code?
- How does the algorithm work?
- What would the code look like if there were no optimizations?

The answers to these questions can act as a guide to help you remove the device specific optimizations.

These questions can be difficult to answer because highly optimized code can be very complex.

6.2.2 Locate and remove device optimizations

There are optimizations for alternative compute devices that have no effect on Mali GPUs or reduce performance.

To optimize the code for a Mali GPU you must first remove all of the following types of optimizations to create a non device-specific reference implementation:

**Use of local or private memory**

Mali GPUs use caches instead of local or private memories so you are not required to manually move data in or out of them. Remove all allocations and copies to or from local or private memory.

**Barriers**

Data transfers to or from local memory are typically synchronized with barriers. Remove any barriers used for this purpose.

**Cache size optimizations**

Some code optimizes reads and writes so data fits into cache lines. This is a useful optimization but the cache line sizes are likely to be different from a Mali GPU. Remove these sorts of optimizations.

**Use of scalars**

Many GPUs work with scalars whereas Mali GPUs use vectors. If there are optimizations based around scalars, remove them.

**Modifications for bank conflicts**

Some code includes optimizations to avoid bank conflicts. Remove these.

**Warps or wavefronts**

Mali GPUs do not use warps or wavefronts. Remove any optimizations for these.

**Optimizations for divergent threads**

Threads on a Mali GPU are independent and cannot diverge. If your code contains optimizations or workarounds for divergent threads, remove them.
6.2.3 Optimizing your OpenCL code for Mali GPUs

To optimize the code for a Mali GPU see Chapter 7 Optimizing OpenCL for Mali GPUs.
Chapter 7
Optimizing OpenCL for Mali GPUs

This chapter describes a number of optimizations to use when writing OpenCL for the Mali-T600 series GPUs. It contains the following sections:

- General optimizations on page 7-2.
- Code optimizations on page 7-3.
- Memory optimizations on page 7-4.
- Kernel optimizations on page 7-6.
- Execution optimizations on page 7-7.
- Reducing the effect of serial computations on page 7-8.
7.1 General optimizations

ARM recommends the following:

- Process large amounts of data. You must be processing a relatively large data set before you see any benefits from using OpenCL. The exact point where you start to see benefits depends on the OpenCL implementation and the Mali GPU you are using.

- Use the best language and processor for the job. Application processors are designed for high speed serial computations. GPUs and OpenCL are designed for parallel processing.

- Experiment to see how fast you can get your algorithm to execute. Operations that perform multiple computations per element of data loaded are typically good for programming in OpenCL.
7.2 Code optimizations

ARM recommends the following:

- Avoid writing kernels that operate on single bytes or other small values. Write kernels that work on vectors. The shader cores in the Mali-T600 Series GPUs contain 128-bit wide vector registers. Vectorize the algorithms in your kernels to make best use of the Mali GPU hardware.

- Use vector load VLOAD instructions on arrays of data even if you do not process the data as vectors. This enables you to load multiple data elements with a single instruction.

- Align data on 128-bit boundaries. This can improve the speed of loading data and ensures data fits evenly into the cache.

- Use the built-in functions. Many of these are implemented as fast hardware instructions. See Appendix B *OpenCL Built-in Functions* for a list of built-in functions with relative speed ratings.

- Use the precise versions of built-in functions. Using the half_ or native_ versions of built-in functions provide no extra performance.

- Avoid writing kernels that use a large number of variables. Using too many variables can impact performance and limits the maximum workgroup size.
7.3 Memory optimizations

OpenCL is typically run in systems where the application processor and the GPU have separate memories. To use OpenCL in these systems you must allocate buffers to copy data to and from the separate memories.

Systems with Mali GPUs typically have a shared memory so you are not required to copy data. However, OpenCL assumes the memories are separate and buffer allocation involves memory copies. This is wasteful because copies take time and consume power.

To avoid the copies, use the OpenCL API to allocate memory buffers and use map() and unmap() operations. These operations enable both the application processor and the Mali GPU to access the data without any copies.

ARM recommends the following:

• Use ALLOC_HOST_PTR to avoid copying memory.
• Do not create buffers with USE_HOST_PTR if possible.
• Do not use private or local memory to improve memory read performance.
• If your kernel is memory bandwidth bound try using a simple formula to compute variables instead of reading from memory. This saves memory bandwidth and might be faster.
• If your kernel is compute bound try reading from memory instead of computing variables. This saves computations and might be faster.

The Mali GPU cannot access the memory buffers created by malloc() because they are not mapped into the memory space of the Mali GPU. This is shown in Figure 7-1.

![Figure 7-1 Memory buffer created by malloc()](image)

The Mali GPU can access the memory buffers created by clCreateBuffer(CL_MEM_USE_HOST_PTR) but buffers created this way must have data copied into them by the application processor. These copy operations are computationally expensive so it is best to avoid this method of allocating buffers if possible. This method of allocating buffers is shown in Figure 7-2 on page 7-5.
Figure 7-2 Memory buffer created by `clCreateBuffer(CL_MEM_USE_HOST_PTR)`

The Mali GPU can access the memory buffers created by `clCreateBuffer(CL_MEM_ALLOC_HOST_PTR)`. This is the preferred method to allocate buffers because data copies are not required. This method of allocating buffers is shown in Figure 7-3.

---

**Note**

- The OpenCL driver must be initialized before allocating memory buffers.
- You must make the initial memory allocation through the OpenCL API.
- If OpenCL calls are repeatedly interleaved with application processor activity, the pointers that access buffers on the CPU might change. The pointer can change to point to a different virtual address that represents the correct physical address.
7.4 Kernel optimizations

ARM recommends the following:

• If your kernel has no preference for the work-group size, pass NULL to the local work size argument of the clEnqueueNDRangeKernel.

• Use work-group sizes that are a multiple of 4. These are more efficient on the Mali-T600 Series GPUs. The maximum work-group size is typically 256.

• If possible, use a work-group size of 128 or 256. These make optimal use of the hardware in the Mali-T600 Series GPUs. If the maximum work-group size is below 128, your kernel might be too complex.

• Some kernels require work-groups for synchronisation of the work-items within the work-group with barriers. These typically require a specific work-group size.

• If your kernels are small, use data with a single dimension and ensure the work-groups size is a power of two.

• Use CL_MAX_WORKGROUP_SIZE to check if the device can execute a kernel that requires a minimum of inter-thread communication. If the device cannot execute the kernel, the algorithm must be implemented as a multi-pass algorithm. This involves enqueuing multiple kernels.

• If you have multiple kernels that work in a sequence, consider combining them into a single kernel. If you combine kernels be careful of dependencies between them.

• Avoid splitting kernels. If you are required to split a kernel, split it into as few kernels as possible.
7.5 Execution optimizations

ARM recommends the following:

- If you are building from source, cache binaries on the storage device.
- If the kernels in use are known at initialization time, creating them once initiates the finalizing compile. Creating the same kernels in the future is faster because the finalized binary is used.
- If you use callbacks to prompt the processor to continue processing data resulting from the execution of a kernel, ensure the callbacks are set before you flush the queue. If you do not do this, the callbacks might occur at the end of a larger batch of work, later than they might have based on actual completion of work.
7.6 Reducing the effect of serial computations

You can reduce the impact of serial components in your application by reducing and optimizing the computations:

- Use memory mapping instead of memory copies to transfer data.
- Optimize the communication code that sends and receives data to reduce latency.
- Keep messages small. Reduce communication overhead by sending only the data that is absolutely required.
- Ensure the size of memory blocks used for communication are a power of 2. This makes the data more cacheable.
- If possible, send more data in a smaller number of transfers.
- Compute values instead of reading them from memory. A simple computation is likely to be faster than reading from memory.
- Do serial computations on the application processors. These are optimized for low latency tasks.
Chapter 8
The Mali OpenCL SDK

The Mali OpenCL SDK includes the following tutorials to help you understand OpenCL development:

**Hello World Tutorial**
This tutorial provides a basic introduction to OpenCL and vectorization.

**Template Tutorial**
This tutorial provides an OpenCL template that you can use as a starting point to develop an OpenCL application.

**Memory Optimizations**
The Memory Optimizations directory contains a Data Sharing tutorial that demonstrates efficient sharing of memory between a Mali-T600 series GPU and an application processor.

**Sobel Filter Tutorial**
This tutorial provides demonstrates the use of the Sobel image filter. This is a simple convolution filter used primarily for edge detection algorithms.

**FIR Float Filter Tutorial**
This tutorial provides demonstrates the use of a floating point *Finite Input Response* (FIR) image filter. You can use this for pixelization or noise reduction.

**Mandelbrot Tutorial**
This tutorial provides demonstrates the use of calculating the Mandelbrot set to produce fractal patterns.
SGEMM Tutorial

This tutorial provides demonstrates the use of Single-Precision General Matrix Multiplication (SGEMM) in OpenCL.

The OpenCL SDK is available from the Mali developer center, http://www.malideveloper.arm.com
Appendix A
OpenCL Data Types

This appendix lists the data types available in OpenCL. Table A-1 shows built-in scalar data types.

<table>
<thead>
<tr>
<th>OpenCL Type</th>
<th>API Type</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td>bool</td>
<td>-</td>
<td>true (1) or false (0)</td>
</tr>
<tr>
<td>char</td>
<td>cl_char</td>
<td>8-bit signed</td>
</tr>
<tr>
<td>Unsigned char, uchar</td>
<td>cl_uchar</td>
<td>8-bit unsigned</td>
</tr>
<tr>
<td>Short</td>
<td>cl_short</td>
<td>16-bit signed</td>
</tr>
<tr>
<td>Unsigned short, ushort</td>
<td>cl_ushort</td>
<td>16-bit unsigned</td>
</tr>
<tr>
<td>Int</td>
<td>cl_int</td>
<td>32-bit signed</td>
</tr>
<tr>
<td>Unsigned int, uint</td>
<td>cl_uint</td>
<td>32-bit unsigned</td>
</tr>
<tr>
<td>Long</td>
<td>cl_long</td>
<td>64-bit signed</td>
</tr>
<tr>
<td>Unsigned long, ulong</td>
<td>cl_ulong</td>
<td>64-bit unsigned</td>
</tr>
<tr>
<td>Float</td>
<td>cl_float</td>
<td>32-bit float</td>
</tr>
<tr>
<td>Half</td>
<td>cl_half</td>
<td>16-bit float, for storage only</td>
</tr>
<tr>
<td>Size_t</td>
<td>-</td>
<td>32-bit or 64-bit unsigned integer</td>
</tr>
<tr>
<td>Ptrdiff_t</td>
<td>-</td>
<td>32-bit or 64-bit unsigned integer</td>
</tr>
</tbody>
</table>
### Table A-2 Built-in vector data types

<table>
<thead>
<tr>
<th>OpenCL Type</th>
<th>API Type</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>char</code>(n)</td>
<td><code>cl_char</code></td>
<td>8-bit signed</td>
</tr>
<tr>
<td><code>uchar</code>(n)</td>
<td><code>cl_uchar</code></td>
<td>8-bit unsigned</td>
</tr>
<tr>
<td><code>short</code>(n)</td>
<td><code>cl_short</code></td>
<td>16-bit signed</td>
</tr>
<tr>
<td><code> ushort</code>(n)</td>
<td><code>cl_ushort</code></td>
<td>16-bit unsigned</td>
</tr>
<tr>
<td><code>int</code>(n)</td>
<td><code>cl_int</code></td>
<td>32-bit signed</td>
</tr>
<tr>
<td><code>uint</code>(n)</td>
<td><code>cl_uint</code></td>
<td>32-bit unsigned</td>
</tr>
<tr>
<td><code>long</code>(n)</td>
<td><code>cl_long</code></td>
<td>64-bit signed</td>
</tr>
<tr>
<td><code>ulong</code>(n)</td>
<td><code>cl_ulong</code></td>
<td>64-bit unsigned</td>
</tr>
<tr>
<td><code>float</code>(n)</td>
<td><code>cl_float</code></td>
<td>32-bit float</td>
</tr>
</tbody>
</table>

\(n\) Where \(n = 2, 3, 4, 8, \text{ or } 16\).

### Table A-3 Other built-in data types

<table>
<thead>
<tr>
<th>OpenCL Type</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>image2d_t</code></td>
<td>2D image handle</td>
</tr>
<tr>
<td><code>image3d_t</code></td>
<td>3D image handle</td>
</tr>
<tr>
<td><code>sampler_t</code></td>
<td>sampler handle</td>
</tr>
<tr>
<td><code>event_t</code></td>
<td>event handle</td>
</tr>
</tbody>
</table>

### Table A-4 Reserved data types

<table>
<thead>
<tr>
<th>OpenCL Type</th>
<th>Description</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>bool</code>(n)</td>
<td>boolean vector</td>
</tr>
<tr>
<td><code>double, double\(n\)</code></td>
<td>64-bit float, vector</td>
</tr>
<tr>
<td><code>half\(n\)</code></td>
<td>16-bit, vector</td>
</tr>
<tr>
<td><code>quad, quad\(n\)</code></td>
<td>128-bit float, vector</td>
</tr>
<tr>
<td>OpenCL Type</td>
<td>Description</td>
</tr>
<tr>
<td>---------------------------------------------------------------------------</td>
<td>---------------------------------</td>
</tr>
<tr>
<td>complex half, complex half(n), imaginary half, imaginary half(n)</td>
<td>16-bit complex, vector</td>
</tr>
<tr>
<td>complex float, complex float(n), imaginary float, imaginary float(n)</td>
<td>32-bit complex, vector</td>
</tr>
<tr>
<td>complex double, complex double(n), imaginary double, imaginary double(n)</td>
<td>64-bit complex, vector</td>
</tr>
<tr>
<td>complex quad, complex quad(n), imaginary quad, imaginary quad(n)</td>
<td>128-bit complex, vector</td>
</tr>
<tr>
<td>float(nxm)</td>
<td>(n\times m) matrix of 32-bit floats</td>
</tr>
<tr>
<td>double(nxm)</td>
<td>(n\times m) matrix of 64-bit floats</td>
</tr>
<tr>
<td>long double, long double(n)</td>
<td>64-bit - 128-bit float, vector</td>
</tr>
<tr>
<td>long long, long long(n)</td>
<td>128-bit signed</td>
</tr>
<tr>
<td>unsigned long long, ulong long, ulonglong(n)</td>
<td>128-bit unsigned</td>
</tr>
</tbody>
</table>
Appendix B
OpenCL Built-in Functions

This appendix lists the OpenCL built-in functions. It contains the following sections:

- **Work-item functions** on page B-2.
- **Math functions** on page B-3.
- **half_ and native_ math functions** on page B-4.
- **Integer functions** on page B-5.
- **Common functions** on page B-6.
- **Geometric functions** on page B-7.
- **Relational functions** on page B-8.
- **Vector data load and store functions** on page B-9.
- **Synchronisation** on page B-10.
- **Asynchronous copy functions** on page B-11.
- **Atomic functions** on page B-12.
- **Miscellaneous vector functions** on page B-13.
- **Image read and write functions** on page B-14.

The functions listed have a relative speed rating. The ratings are from A to C, where A is the fastest. For maximum performance always use vectors and try to use functions rated with an A or B.

---

**Note**

Ratings for memory accesses are separate from arithmetic operations. An A rated memory operation might be equivalent to a C rated arithmetic operation.
B.1 Work-item functions

Table B-1 lists the work-item functions.

<table>
<thead>
<tr>
<th>Function</th>
<th>Speed</th>
</tr>
</thead>
<tbody>
<tr>
<td>get_work_dim()</td>
<td>A</td>
</tr>
<tr>
<td>get_global_size()</td>
<td>A</td>
</tr>
<tr>
<td>get_global_id()</td>
<td>A</td>
</tr>
<tr>
<td>get_local_size()</td>
<td>A</td>
</tr>
<tr>
<td>get_local_id()</td>
<td>A</td>
</tr>
<tr>
<td>get_num_groups()</td>
<td>A</td>
</tr>
<tr>
<td>get_group_id()</td>
<td>A</td>
</tr>
<tr>
<td>get_global_offset()</td>
<td>A</td>
</tr>
</tbody>
</table>
### B.2 Math functions

Table B-2 lists the math functions.

<table>
<thead>
<tr>
<th>Function</th>
<th>Speed</th>
<th>Function</th>
<th>Speed</th>
<th>Function</th>
<th>Speed</th>
</tr>
</thead>
<tbody>
<tr>
<td>fabs()</td>
<td>A</td>
<td>acos()</td>
<td>B</td>
<td>acosh()</td>
<td>C</td>
</tr>
<tr>
<td>ceil()</td>
<td>A</td>
<td>acospi()</td>
<td>B</td>
<td>asinh()</td>
<td>C</td>
</tr>
<tr>
<td>fdim()</td>
<td>A</td>
<td>asin()</td>
<td>B</td>
<td>atanh()</td>
<td>C</td>
</tr>
<tr>
<td>fmax()</td>
<td>A</td>
<td>asinpi()</td>
<td>B</td>
<td>copysign()</td>
<td>C</td>
</tr>
<tr>
<td>fmin()</td>
<td>A</td>
<td>atan()</td>
<td>B</td>
<td>erf()</td>
<td>C</td>
</tr>
<tr>
<td>mad()</td>
<td>A</td>
<td>atan2()</td>
<td>B</td>
<td>erf()</td>
<td>C</td>
</tr>
<tr>
<td>maxmag()</td>
<td>A</td>
<td>atanpi()</td>
<td>B</td>
<td>fmod()</td>
<td>C</td>
</tr>
<tr>
<td>minmag()</td>
<td>A</td>
<td>atan2pi()</td>
<td>B</td>
<td>fract()</td>
<td>C</td>
</tr>
<tr>
<td>rint()</td>
<td>A</td>
<td>cbrt()</td>
<td>B</td>
<td>frexp()</td>
<td>C</td>
</tr>
<tr>
<td>round()</td>
<td>A</td>
<td>cos()</td>
<td>B</td>
<td>hypot()</td>
<td>C</td>
</tr>
<tr>
<td>trunc()</td>
<td>A</td>
<td>cosh()</td>
<td>B</td>
<td>ilogb()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>exp()</td>
<td>B</td>
<td>ldexp()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>exp2()</td>
<td>B</td>
<td>log()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>exp10()</td>
<td>B</td>
<td>log1()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>expm1()</td>
<td>B</td>
<td>log10()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>floor()</td>
<td>B</td>
<td>log1p()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>fma()</td>
<td>B</td>
<td>logb()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>log2()</td>
<td>B</td>
<td>modf()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>pow()</td>
<td>B</td>
<td>nan()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>pown()</td>
<td>B</td>
<td>nextafter()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>powr()</td>
<td>B</td>
<td>remainder()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>rsqrt()</td>
<td>B</td>
<td>remquo()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>sin()</td>
<td>B</td>
<td>rootn()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>sincos()</td>
<td>B</td>
<td>sinh()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>sinpi()</td>
<td>B</td>
<td>tan()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>sqrt()</td>
<td>B</td>
<td>tanh()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>-</td>
<td>-</td>
<td>tanpi()</td>
<td>C</td>
</tr>
<tr>
<td>-</td>
<td>-</td>
<td>-</td>
<td>-</td>
<td>tgamma()</td>
<td>C</td>
</tr>
</tbody>
</table>
B.3 half_ and native_ math functions

Typically, on most architectures there is a trade-off between accuracy and speed. The Mali-T600 Series GPUs implements the full precision variants of the math functions at full speed so you are not required to make this trade-off.

The half_ and native_ variants of the math functions are provided for portability. They are no faster than the precise variants. See Math functions on page B-3.

Table B-3 lists the half_ and native_ math functions.

<table>
<thead>
<tr>
<th>half_ functions</th>
<th>native_ functions</th>
</tr>
</thead>
<tbody>
<tr>
<td>half_cos()</td>
<td>native_cos()</td>
</tr>
<tr>
<td>half_divide()</td>
<td>native_divide()</td>
</tr>
<tr>
<td>half_exp()</td>
<td>native_exp()</td>
</tr>
<tr>
<td>half_exp2()</td>
<td>native_exp2()</td>
</tr>
<tr>
<td>half_exp10()</td>
<td>native_exp10()</td>
</tr>
<tr>
<td>half_log()</td>
<td>native_log()</td>
</tr>
<tr>
<td>half_log2()</td>
<td>native_log2()</td>
</tr>
<tr>
<td>half_log10()</td>
<td>native_log10()</td>
</tr>
<tr>
<td>half_powr()</td>
<td>native_powr()</td>
</tr>
<tr>
<td>half_recip()</td>
<td>native_recip()</td>
</tr>
<tr>
<td>half_rgba()</td>
<td>native_rgba()</td>
</tr>
<tr>
<td>half_sin()</td>
<td>native_sin()</td>
</tr>
<tr>
<td>half_sqrt()</td>
<td>native_sqrt()</td>
</tr>
<tr>
<td>half_tan()</td>
<td>native_tan()</td>
</tr>
</tbody>
</table>
### B.4 Integer functions

Table B-4 lists the integer functions.

<table>
<thead>
<tr>
<th>Function</th>
<th>Speed</th>
</tr>
</thead>
<tbody>
<tr>
<td>abs()</td>
<td>A</td>
</tr>
<tr>
<td>abs_diff()</td>
<td>A</td>
</tr>
<tr>
<td>add_sat()</td>
<td>A</td>
</tr>
<tr>
<td>hadd()</td>
<td>A</td>
</tr>
<tr>
<td>rhadd()</td>
<td>A</td>
</tr>
<tr>
<td>clz()</td>
<td>A</td>
</tr>
<tr>
<td>max()</td>
<td>A</td>
</tr>
<tr>
<td>min()</td>
<td>A</td>
</tr>
<tr>
<td>sub_sat()</td>
<td>A</td>
</tr>
<tr>
<td>mad24(), identical to 32-bit multiply accumulate</td>
<td>A</td>
</tr>
<tr>
<td>mul24(), identical to 32-bit multiplies</td>
<td>A</td>
</tr>
<tr>
<td>clamp()</td>
<td>B</td>
</tr>
<tr>
<td>mad_hi()</td>
<td>B</td>
</tr>
<tr>
<td>mul_hi()</td>
<td>B</td>
</tr>
<tr>
<td>mad_sat()</td>
<td>B</td>
</tr>
<tr>
<td>rotate()</td>
<td>B</td>
</tr>
<tr>
<td>upsample()</td>
<td>B</td>
</tr>
</tbody>
</table>
B.5 Common functions

Table B-5 lists the common functions.

<table>
<thead>
<tr>
<th>Function</th>
<th>Speed</th>
</tr>
</thead>
<tbody>
<tr>
<td>max()</td>
<td>A</td>
</tr>
<tr>
<td>min()</td>
<td>A</td>
</tr>
<tr>
<td>step()</td>
<td>A</td>
</tr>
<tr>
<td>clamp()</td>
<td>B</td>
</tr>
<tr>
<td>degrees()</td>
<td>B</td>
</tr>
<tr>
<td>mix()</td>
<td>B</td>
</tr>
<tr>
<td>radians()</td>
<td>B</td>
</tr>
<tr>
<td>smoothstep()</td>
<td>B</td>
</tr>
<tr>
<td>sign()</td>
<td>B</td>
</tr>
</tbody>
</table>
B.6 Geometric functions

Table B-6 lists the geometric functions.

<table>
<thead>
<tr>
<th>Function</th>
<th>Speed</th>
</tr>
</thead>
<tbody>
<tr>
<td>dot()</td>
<td>A</td>
</tr>
<tr>
<td>normalize()</td>
<td>B</td>
</tr>
<tr>
<td>fast_distance()</td>
<td>B</td>
</tr>
<tr>
<td>fast_length()</td>
<td>B</td>
</tr>
<tr>
<td>fast_normalize()</td>
<td>B</td>
</tr>
<tr>
<td>cross()</td>
<td>B</td>
</tr>
<tr>
<td>distance()</td>
<td>B</td>
</tr>
<tr>
<td>length()</td>
<td>B</td>
</tr>
</tbody>
</table>
### B.7 Relational functions

Table B-7 lists the relational functions.

<table>
<thead>
<tr>
<th>Function</th>
<th>Speed</th>
</tr>
</thead>
<tbody>
<tr>
<td>any()</td>
<td>A</td>
</tr>
<tr>
<td>all()</td>
<td>A</td>
</tr>
<tr>
<td>bitselect()</td>
<td>A</td>
</tr>
<tr>
<td>select()</td>
<td>A</td>
</tr>
<tr>
<td>isequal()</td>
<td>A</td>
</tr>
<tr>
<td>isnotequal()</td>
<td>A</td>
</tr>
<tr>
<td>isgreater()</td>
<td>A</td>
</tr>
<tr>
<td>isgreaterequal()</td>
<td>A</td>
</tr>
<tr>
<td>isless()</td>
<td>A</td>
</tr>
<tr>
<td>islessequal()</td>
<td>A</td>
</tr>
<tr>
<td>islessgreater()</td>
<td>A</td>
</tr>
<tr>
<td>isfinite()</td>
<td>B</td>
</tr>
<tr>
<td>isinf()</td>
<td>B</td>
</tr>
<tr>
<td>isnan()</td>
<td>B</td>
</tr>
<tr>
<td>isnormal()</td>
<td>B</td>
</tr>
<tr>
<td>isordered()</td>
<td>B</td>
</tr>
<tr>
<td>isunordered()</td>
<td>B</td>
</tr>
<tr>
<td>signbit()</td>
<td>B</td>
</tr>
</tbody>
</table>
B.8 Vector data load and store functions

Table B-8 lists the vector data load and store functions. These are all speed A.

<table>
<thead>
<tr>
<th>Function</th>
<th>Speed</th>
</tr>
</thead>
<tbody>
<tr>
<td>vload()</td>
<td>A</td>
</tr>
<tr>
<td>vstore()</td>
<td>A</td>
</tr>
<tr>
<td>vload_half()</td>
<td>A</td>
</tr>
<tr>
<td>vstore_half()</td>
<td>A</td>
</tr>
<tr>
<td>vloada_half()</td>
<td>A</td>
</tr>
<tr>
<td>vstorea_half()</td>
<td>A</td>
</tr>
</tbody>
</table>
B.9 Synchronisation

Table B-9 lists the synchronisation functions. These have no speed rating because synchronisation functions wait for multiple threads to complete. The time this takes determines the length of time the functions take in your application.

<table>
<thead>
<tr>
<th>Function</th>
</tr>
</thead>
<tbody>
<tr>
<td>barrier()</td>
</tr>
<tr>
<td>mem_fence()</td>
</tr>
<tr>
<td>read_mem_fence()</td>
</tr>
<tr>
<td>write_mem_fence()</td>
</tr>
</tbody>
</table>
### B.10 Asynchronous copy functions

Table B-10 lists the asynchronous copy functions. These have no speed rating because the copy speed depends on the size of the data copied.

<table>
<thead>
<tr>
<th>Function</th>
</tr>
</thead>
<tbody>
<tr>
<td>async_work_group_copy()</td>
</tr>
<tr>
<td>async_work_group_strided_copy()</td>
</tr>
<tr>
<td>wait_group_events()</td>
</tr>
<tr>
<td>prefetch()</td>
</tr>
</tbody>
</table>
# B.11 Atomic functions

Table B-11 lists the atomic functions.

<table>
<thead>
<tr>
<th>Function</th>
<th>Speed</th>
</tr>
</thead>
<tbody>
<tr>
<td>atomic_add()</td>
<td>B</td>
</tr>
<tr>
<td>atomic_sub()</td>
<td>B</td>
</tr>
<tr>
<td>atomic_xchg()</td>
<td>B</td>
</tr>
<tr>
<td>atomic_inc()</td>
<td>B</td>
</tr>
<tr>
<td>atomic_dec()</td>
<td>B</td>
</tr>
<tr>
<td>atomic_cmpxchg()</td>
<td>B</td>
</tr>
<tr>
<td>atomic_min()</td>
<td>B</td>
</tr>
<tr>
<td>atomic_max()</td>
<td>B</td>
</tr>
<tr>
<td>atomic_and()</td>
<td>B</td>
</tr>
<tr>
<td>atomic_or()</td>
<td>B</td>
</tr>
<tr>
<td>atomic_xor()</td>
<td>B</td>
</tr>
</tbody>
</table>
## B.12 Miscellaneous vector functions

Table B-12 lists the miscellaneous vector functions.

<table>
<thead>
<tr>
<th>Function</th>
<th>Speed</th>
</tr>
</thead>
<tbody>
<tr>
<td>vec_step()</td>
<td>A</td>
</tr>
<tr>
<td>shuffle()</td>
<td>A</td>
</tr>
<tr>
<td>shuffle2()</td>
<td>B</td>
</tr>
</tbody>
</table>
## B.13 Image read and write functions

Table B-13 lists the image read and write functions.

<table>
<thead>
<tr>
<th>Function</th>
<th>Speed</th>
</tr>
</thead>
<tbody>
<tr>
<td><code>read_imagef()</code></td>
<td>A</td>
</tr>
<tr>
<td><code>read_imagei()</code></td>
<td>A</td>
</tr>
<tr>
<td><code>read_imageui()</code></td>
<td>A</td>
</tr>
<tr>
<td><code>write_imagef()</code></td>
<td>A</td>
</tr>
<tr>
<td><code>write_imagei()</code></td>
<td>A</td>
</tr>
<tr>
<td><code>write_imageui()</code></td>
<td>A</td>
</tr>
<tr>
<td><code>get_image_width()</code></td>
<td>B</td>
</tr>
<tr>
<td><code>get_image_height()</code></td>
<td>B</td>
</tr>
<tr>
<td><code>get_image_depth()</code></td>
<td>B</td>
</tr>
<tr>
<td><code>get_image_channel_data_type()</code></td>
<td>B</td>
</tr>
<tr>
<td><code>get_image_channel_order()</code></td>
<td>B</td>
</tr>
<tr>
<td><code>get_image_dim()</code></td>
<td>B</td>
</tr>
</tbody>
</table>
Appendix C
OpenCL Extensions

The Mali OpenCL driver supports the following extensions:

- `cl_khr_int64_base_atomics`.
- `cl_khr_int64_extended_atomics`.
- `cl_khr_global_int32_base_atomics`.
- `cl_khr_global_int32_extended_atomics`.
- `cl_khr_local_int32_base_atomics`.
- `cl_khr_local_int32_extended_atomics`.
- `cl_khr_byte_addressable_store`. 