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RELEASE NOTES

v1.3.0:

• ONNXRuntime updated to v1.11.1.
• Backend updated to 0.4.0
  o Misc models speed up
  o Bug fixes

v1.2.0:

v1.1.0:

• Updated to use Ampere Optimized ONNXRuntime 0.3.0
  o Misc models speed up

OVERVIEW

Ampere Optimized ONNXRuntime inference acceleration engine is fully integrated with the ONNX Runtime framework. ONNX models and ONNX Runtime software written with the ONNX Runtime API can run as-is, without modifications.

ONNX RUNTIME FRAMEWORK

Python is installed with Ampere Optimized ONNXRuntime and all dependencies. No additional installation steps are needed.

Versions Compatibility

This release is based on ONNX Runtime 1.10.0. Please refer to ONNX Runtime version compatibility documentation, found at ONNX Runtime and ONNX Versioning Guide, to check the compatibility of models built with older versions of ONNX Runtime.

PYTHON

ONNX Runtime 1.10.0 is built for Python 3.8. Regarding other Python versions, contact your Ampere sales representative. If you are using the software through a third party, contact their customer support team for help. You can also contact the Ampere AI team at ai-support@amperecomputing.com.

CONFIGURATIONS

Ampere Optimized ONNXRuntime inference engine can be configured by a set of environment variables for performance and debugging purposes. They can be set in the command line when running ONNX models (e.g., AIO_NUM_THREADS=16 python run.py) or set in the shell initialization script.
**AIO_PROCESS_MODE**
This variable controls whether Ampere Optimized ONNXRuntime inference engine is used to run the ONNX model:
- 0: disabled.
- 1: enabled (Default).

**AIO_CPU_BIND**
Enables core binding. If enabled, each Ampere Optimized ONNXRuntime thread will bind itself to a single core:
- 0: Core binding disabled.
- 1: Core binding enabled (Default).

**AIO_MEM_BIND**
Binds memory to NUMA (Non-uniform memory access) node 0. For optimal performance, `numactl` ([https://linux.die.net/man/8/numactl](https://linux.die.net/man/8/numactl)) is preferred. `numactl bind` will affect both the ONNX Runtime framework and framework and the optimized framework buffers, while the optimized framework is unable to affect buffers allocated by the ONNX Runtime framework:
- 0: Membind disabled.
- 1: Membind to node 0 (Default).

**AIO_NUMA_CPUS**
Select cores that Ampere Optimized ONNXRuntime should bind to (if CPU_BIND is enabled):
- Not set: use the first N cores of the machine, excluding hyper-threaded machines (Default).
- Set: try to use N first cores from the list of cores for N threads. The list is in space-separated, 0-based number format. For example, selecting cores 0 to 1: `AIO_NUMA_CPUS="0 1"`.

**AIO_NUM_THREADS**
Specifies the number of cores that Ampere Optimized ONNXRuntime should use:
- Not set: use one core (Default).
- "all": use all cores, as specified by `AIO_NUMA_CPUS`.
- N: use N cores.
**AIO_DEBUG_MODE**

Control verbosity of debug messages:

- 0: No messages
- 1: Errors only
- 2: Basic information, warnings, and errors (Default)
- 3: Most messages
- 4: All messages

**QUICKSTART**

The following instructions run on Altra/Altra Max Linux machines installed with Docker. To initialize Ampere Optimized ONNXRuntime environment run:

```bash
$ wget -O aio-onnxrt.tar.gz "<your_unique_url>"
$ docker load < aio-onnxrt.tar.gz
$ docker run --privileged=true --rm --name onnxrt-aio --network host -it aio-onnxrt-1.11.1:1.3.0
```

Skip the above steps if running **without a Docker container**.

You can try Ampere Optimized ONNXRuntime by either running the Jupyter Notebook examples or Python scripts on the CLI level.

To run the Jupyter Notebook QuickStart examples follow the instructions below:

Set `AIO_NUM_THREADS` to the requested value first.

```bash
$ export AIO_NUM_THREADS=16; export OMP_NUM_THREADS=16
$ cd /workspace/aio-examples/
$ bash download_models.sh
$ bash start_notebook.sh
```

If you run the Jupyter Notebook Quickstart on a cloud instance, make sure your machine has port 8080 open and on your local device run:

```bash
$ ssh -N -L 8080:localhost:8080 -l <ssh_key> your_user@xxx.xxx.xxx.xxx
```

Use a browser to point to the URL printed out by the Jupyter Notebook launcher.

You will find the Jupyter Notebook examples (examples.ipynb) under the /classification and /object detection folders.

The examples run through several inference models, visualize results they produce and present the performance numbers.
To use CLI-level scripts:

Set AIO_NUM_THREADS to the requested value first.

```bash
$ export AIO_NUM_THREADS=16; export OMP_NUM_THREADS=16
$ cd /workspace/aio-examples/
$ bash download_models.sh
$ pip install torch
```

Go to the directory of choice, e.g.,

```bash
$ cd classification/resnet_50_v1.5
```

Evaluate the model.

```bash
$ python3 run.py -m resnet_50_v1.5_fp32.onnx -p fp32
```

AMPERE OPTIMIZED FRAMEWORKS PROGRAMMING GUIDE

Overview

Ampere Optimized TensorFlow is powered by Ampere® AI backend which accelerates Deep Learning (DL) operations on the Ampere Altra family of processors. Ampere Optimized Frameworks accelerate DL operations through model optimization, highly vectorized compute kernels and multi-thread operations that are automatically tuned to deliver the best latency and throughput on Ampere Altra processors. It delivers 2-5x gains over alternative backend solutions.
**Supported Inference Ops**

Ampere Optimized Tensorflow accelerates most common Tensorflow ops that are used in various types of models. Here is a list of accelerated ops and formats (Note: non-accelerated ops will still run without problem, at the original framework operator speed):

<table>
<thead>
<tr>
<th>Op</th>
<th>FP32</th>
<th>FP16</th>
<th>Remarks</th>
</tr>
</thead>
<tbody>
<tr>
<td>Conv2D</td>
<td>Y</td>
<td>Y</td>
<td></td>
</tr>
<tr>
<td>Conv3D</td>
<td>Y</td>
<td>N</td>
<td>NDHWC only</td>
</tr>
<tr>
<td>__FusedConv2D</td>
<td>Y</td>
<td>Y</td>
<td></td>
</tr>
<tr>
<td>FusedBatchNorm</td>
<td>Y</td>
<td>N</td>
<td>NHWC only</td>
</tr>
<tr>
<td>FusedBatchNormV2</td>
<td>Y</td>
<td>N</td>
<td></td>
</tr>
<tr>
<td>FusedBatchNormV3</td>
<td>Y</td>
<td>N</td>
<td></td>
</tr>
<tr>
<td>MaxPool</td>
<td>Y</td>
<td>Y</td>
<td>NHWC only</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td>2D Max Pooling only</td>
</tr>
<tr>
<td>AvgPool</td>
<td>Y</td>
<td>Y</td>
<td>NHWC only</td>
</tr>
<tr>
<td></td>
<td></td>
<td></td>
<td>2D Average Pooling only</td>
</tr>
<tr>
<td>MatMul</td>
<td>Y</td>
<td>Y</td>
<td>transpose_a == 0 only</td>
</tr>
<tr>
<td>__FusedMatMul</td>
<td>Y</td>
<td>Y</td>
<td>transpose_a == 0 only</td>
</tr>
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<td>Y</td>
<td>adj_x == 0 only</td>
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<td>BatchMatMulV2</td>
<td>Y</td>
<td>Y</td>
<td>adj_x == 0 only</td>
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<td>Mean</td>
<td>Y</td>
<td>Y</td>
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<td>Mul</td>
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<td>Y</td>
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<tr>
<td>Add</td>
<td>Y</td>
<td>Y</td>
<td></td>
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<tr>
<td>AddV2</td>
<td>Y</td>
<td>Y</td>
<td></td>
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<tr>
<td>BiasAdd</td>
<td>Y</td>
<td>Y</td>
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<td>Y</td>
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<td>Y</td>
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<td>RealDiv</td>
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<tr>
<td>Tanh</td>
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<td>Y</td>
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</tr>
<tr>
<td>Sqrt</td>
<td>Y</td>
<td>Y</td>
<td></td>
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<tr>
<td>Square</td>
<td>Y</td>
<td>Y</td>
<td></td>
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<tr>
<td>Rsqrt</td>
<td>Y</td>
<td>Y</td>
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</tr>
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<td>SquaredDifference</td>
<td>Y</td>
<td>Y</td>
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<td>LeakyRelu</td>
<td>Y</td>
<td>Y</td>
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<td>Softmax</td>
<td>Y</td>
<td>Y</td>
<td></td>
</tr>
<tr>
<td>AddN</td>
<td>Y</td>
<td>Y</td>
<td></td>
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<tr>
<td>Pad</td>
<td>Y</td>
<td>Y</td>
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</tr>
<tr>
<td>Concat</td>
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<td>Y</td>
<td>axis_constant only</td>
</tr>
<tr>
<td>ConcatV2</td>
<td>Y</td>
<td>Y</td>
<td>axis constant only</td>
</tr>
<tr>
<td>Operation</td>
<td>Y</td>
<td>N</td>
<td>Details</td>
</tr>
<tr>
<td>-------------------------</td>
<td>---</td>
<td>---</td>
<td>-------------------------------------------------------------------------</td>
</tr>
</tbody>
</table>
| Gather                  |   |   | indices int32 only
|                         |   |   | axis constant only                                                      |
| GatherV2                |   |   | indices int32 only
|                         |   |   | axis constant only                                                      |
|                         |   |   | batch_dim = 0 only                                                      |
| StridedSlice            |   |   | index int32 only
|                         |   |   | begin_mask and end_mask only                                           |
| Squeeze                 | Y | Y |                                                                        |
| DepthwiseConv2dNative   | Y | Y |                                                                        |
| Reshape                 | Y | Y |                                                                        |
| ExpandDims              | Y | Y |                                                                        |
| Transpose               | Y | Y | perm constant only                                                      |
| Erf                     | Y | Y |                                                                        |
| SplitV                  | Y | Y | axis constant only                                                      |
| Conv3dBackpropInputV2   | Y | N | NDHWC only                                                              |

Ampere AI continues to expand the coverage of TensorFlow ops. If your model has any op that is not listed in the table or custom ops that need acceleration, please contact ai-support@amperecomputing.com.

Ampere Optimized TensorFlow also supports acceleration of TensorFlow Lite int8 models. Please contact us for information about TensorFlow Lite model support.

**TensorFlow Eager Execution and Graph Execution**

While TensorFlow Eager Execution provides excellent model building, programming, and debugging experience, it is slower than graph execution. So, graph execution is typically used for inference deployment. In current version Ampere Optimized TensorFlow only accelerates Graph Execution mode.

After building your model in Eager mode, you can use tf.function() to compile you eager function into callable graph. More details can be found in TensorFlow documentation at: https://www.tensorflow.org/api_docs/python/tf/function.

Ampere model library also provides some sample code in how to run eager model efficiently, access AML at: https://github.com/AmpereComputingAI/ampere_model_library/blob/main/natural_language_processing/extractive_question_answering/electra_large/run.py.

**Threading**

Ampere Optimized TensorFlow controls the number of intra_op threads of AIO with tensorflow.config.threading.set_intra_op_parallelism_threads() (Or in the case of TF v1 session, set config.intra_op_parallelism_threads). This controls both the number of threads used for ops delegated to AIO as well as the ops running on default CPU backend.
Some default CPU backend ops (non-AIO) also need to set OMP_NUM_THREADS environment variable to control the intra_op threads.

**Programming Tips**

- In the first inference pass, Ampere Optimized Tensorflow performs runtime compilation of TF graphs. So, the latency of the first pass is expected to be longer. Subsequent passes will be accelerated.

- Frozen TF models will provide slightly better performance. Please see TF documentations in how to generate frozen graphs.

- Ampere Optimized TensorFlow provides much better latency scaling as core count increase, comparing to other platforms. You can easily try the optimal number of cores with the above intra_op_parallelism_threads configurations that can give you the best performance/$, while meeting your latency requirements.

- If any issues occur, Ampere AI team is ready to help. Typically, the first step is to get more debug logs and send it to ai-support@amperecomputing.com. Please set environment variable AIO_DEBUG_MODE=5 to capture low level logs.

We can also provide more in-depth profiling of your model to help enhancing performance to meet your need.