SOFTWARE STACK: TENSORFLOW PYTORCH POPLAR

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GRAPHCORE





AGENDA

- Architecture Refresher
- Software Ecosystem
- TensorFlow2/Keras
- PyTorch
- Poplar



ARCHITECTURE REFRESHER



IPU – ARCHITECTURED FOR AI

Massive parallelism with ultrafast memory access



Massively parallel MIMD. Designed for fine-grained, highperformance computing

PROVEN IPU ADVANTAGE SELECT CASE STUDIES ACROSS MANY INDUSTRIES & FIELDS





EXECUTION MODEL





OUTPUT FROM POPVISION GRAPH ANALYSER

GRAPHCORE

BOW-2000 IPU MACHINE

IU blade form factor delivering 1.4 PetaFLOPS AI Compute

BOW IPU-2000





BOW-2000 TOPOLIGY





4x Bow IPUs

- 1.4 PFLOP₁₆ compute
- 5,888 processor cores
- > 35,000 independent parallel threads



COMMUNICATIONS

Exchange Memory

- 3.6GB In-Processor Memory @ 260 TB/s
- 128GB Streaming Memory DRAM (up to 256GB)

IPU-Fabric managed by IPU-GW

- Host-Link 100GE to Poplar Server for standard data center networking
- IPU-Link 2D Torus for intra-POD64 communication
- GW-Link 2x 100Gbps Gateway-Links for rack-torack - flexible topology

BOW-POD64 TOPOLOGY



IPU DEVELOPER ECOSYSTEM



GRAPHCORE

GRAPHCORE SOFTWARE ECOSYSTEM

WORLD CLASS DEVELOPER RESOURCES FOR IPU USERS



WWW.GRAPHCORE.AI/DEVELOPER

GRAPHCORE **GRAPHCORE DOCUMENTS** Graphcore Documents Version: Latest Software Hardware **Getting Started** Search docs Documents Documents Background information and **Getting Started** quick-start guides for Documentation for the Documentation for installing Graphcloud and Pod and using IPU-Machines and Poplar SDK and other Software Documents Pod systems systems software Hardware Documents Technical Notes and White Papers Technical Notes and Examples and **Document Updates** Examples and Tutorials White Papers Tutorials The latest news about new Document Updates Technical notes and white documents and examples Tutorials and application papers on Graphcore examples for running on the Alphabetical List of All Documents technology IPU Graphcore License Agreements



Getting started with PyTorch for the IPU

Running a basic model for training and inference

Al Customer Engineer, Chris Bogdiukiewicz introduces PyTorch for the IPU. With PopTorch™ - a simple Python wrapper for PyTorch programs, developers can easily run models, directly on Graphcore IPUs with a few lines of extra code.

 $\textbf{Get the Code} \quad \rightarrow \quad$

In this video, Chris provides a quick demo on running a basic model for both training and inference using a MNIST based example.

Read the Guide \rightarrow



OPEN SOURCE

github.com/graphcore

- As part of our ethos to put power in the hands of AI developers, Graphcore open sourced in 2020
- PopLibs[™], PopART, PyTorch & TensorFlow for IPU fully open source and available on GitHub
- Our code is public and open for code contributions from the wider ML developer community





VIDEO + GITHUB TUTORIALS

A comprehensive set of online developer training materials and educational content





Running PyTorch on the IPU: NLP



Bulk Synchronous Parallel E







| Learn how to create and run program PopLibs with our hands-on programm | ns using Poplar and ning tutorials. | | |
|---|--|-----|------------------------------|
| Programs and Variables | Using PopLibs | | Writing Vertex Code |
| Profiling Output | Basic Machine Learning Example | | Matrix-Vector Multiplication |
| Matrix-Vector Multiplication Optimisation | Simple PyTorch for the IPU | NEW | |

Tutorial 1: programs and variables

Copy the file tut1 variables/start_here/tut1.cpp to your working directory and open it in an editor. The file contains the outline of a C++ program including some Poplar library headers and a namespace.

Graphs, variables and programs

All Poplar programs require a Graph object to construct the computation graph. Graphs are always created for a specific target (where the target is a description of the hardware being targeted, such as an IPU). To obtain the target we need to choose a device.

The tutorials use a simulated target by default, so will run on any machine even if it has no Graphcore hardware attached. On systems with accelerator hardware, the header file poplar/DeviceManager.hpp contains API calls to enumerate and return Device objects for the attached hardware.

Simulated devices are created with the IPUModel class, which models the functionality of an IPU on the host. The createDevice function creates a new virtual device to work with. Once we have this device we can create a Graph object to target it.

· Add the following code to the body of main :

// Create the IPU Model device IPUModel ipuModel: Device device = ipuModel.createDevice(); Target target = device.getTarget();

// Create the Graph object Graph graph(target):

Any program running on an IPU needs data to work on. These are defined as variables in the graph.

· Add the following code to create the first variable in the program:

Tutorial 5: a basic machine learning example

This tutorial contains a complete training program that performs a logistic regression on the MNIST data set, using gradient descent. The files for the demo are in tut5_ml. There are no coding steps in the tutorial. The task is to understand the code, build it and run it. You can build the code using the supplied makefile.

Before you can run the code you will need to run the get_mnist.sh script to download the MNIST data.

The program accepts an optional command line argument to make it use the IPU hardware instead of a simulated IPU.

As you would expect, training is significantly faster on the IPU hardware.

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RESOURCES CENTRE

graphcore.ai/resources

- Central source of research papers, white papers, videos, on-demand webinars and documentation
- Product resources for ML Engineers & IT / Infrastructure Managers now available





GRAPHCORE DEVELOPER ECOSYSTEM

СС



STANDARD ML FRAMEWORK SUPPORT

Develop models using standard high-level frameworks or port existing models





Platforms



GRAPHCORE SOFTWARE







СC

ENHANCED MODEL GARDEN

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PUBLIC ACCESS TO WIDE VARIETY OF MODELS, READY TO RUN ON IPU

NEW FILTER/SEARCH CAPABILITY

DIRECT ACCESS TO GITHUB

https://www.graphcore.ai/resources/model-garden

MODEL GARDEN COVERAGE



20

POPVISION®

INDUSTRY LEADING AI APPLICATION PERFORMANCE ANALYSIS TOOLS

Introduced in Q2 2020 our PopVision analysis tools provide detailed observability of IPU applications

- Poplar Graph Analyser allows visual inspection of IPU execution down to the individual tile level
- Poplar System Analyser gives users the ability to view host side application and IPU interaction
- Both tools extend debug information back up into Tensorflow and Pytorch for developers

SUPPORTED PLATFORMS







POPVISION TOOLS

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IPU MEMORY ANALYSIS

Capture memory information from your ML models when executed on IPUs. Inspect variable placement, size and liveness throughout the execution.



EXECUTION TRACE REPORT

View the output of instrumenting a Poplar program, capturing cycle counts for each step. See execution statistics, tile balance, cycle proportions and compute-set details.

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REPORT COMPARISONS

Open two reports at once to compare their memory, execution, liveness and operations. Visualise where efficiencies can be made with different model parameters.



HOST EXECUTION ANALYSIS

Understand the execution of IPU-targeted software on your host system processors. Identify any bottlenecks between CPUs and IPUs across a visual interactive timeline.



GRAPH DATA

Plot graph data of any numerical data points from the host or IPU processor systems, such as board temperature, power consumption and IPU utilisation.

| open file | |
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| Enter your usemaner and the host to con | ned to. |
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LOCAL + REMOTE REPORTS

Ability to open reports either on your local machine, or remotely on the host machine. The Graph Analyser also supports local and remote report access.



POPVISION PERFORMANCE ANALYSER

Navigation 🥥 🔍 65787% 🔍 « » <



Utilization/memory map of every tile/every IPU



TF2/KERAS ON IPU



KERAS ON IPU

- IPU optimized Keras Model and Sequential are available for the IPU. These have the following features:
 - * On-device training loop for reduction of communication overhead.
 - * Gradient accumulation for simulating larger batch sizes.
 - * Automatic data-parallelisation of the model when placed on a multi-IPU device.



| ✿ gpu_cnn_keras.py ↔ ipu_cnn_keras.py tf_keras | |
|---|---|
| and the second se | import tensorflow as tf |
| Crac ras.layers import * | from tensorflow.keras.layers import * |
| ICIAS //////////////////////////////////// | + from tensorflow.python import ipu |
| | |
| | + cfg = ipu.config.IPUConfig() |
| | + cfg.auto_select_ipus = 1 |
| | + cfg.configure_ipu_system() |
| | + with ipu.ipu_strategy.iPUStrategy().scope(): |
| (x_train, y_train), (x_test, y_test) = tf.keras.datasets.citar10.load_data() | (x_train, y_train), (x_test, y_test) = tr.keras.datasets.citar10.load_data() |
| $x_{train} = x_{train.astype}(1000132) / 255.0$ | $x_{train} = x_{train.astype}(troats2) / 255.0$ |
| y_train = tf.keras.utits.to_categoricat(y_train, 10) | y_train = tf.keras.utits.to_categoricat(y_train, 10) |
| ds_train = tr.data.bataset.trom_tensor_stices((x_train, y_train)).batch(64, drop_remainde | ds_train = tf.data.Dataset.from_tensor_stices((x_train, y_train)).batch(64, drop_rema |
| <pre>model = tf.keras.Sequential([</pre> | <pre>model = tf.keras.Sequential([</pre> |
| <pre>Conv2D(32, (3, 3), padding='same', input_shape=x_train.shape[1:]),</pre> | <pre>Conv2D(32, (3, 3), padding='same', input_shape=x_train.shape[1:]),</pre> |
| Activation('relu'), | Activation('relu'), |
| Conv2D(32, (3, 3)), | Conv2D(32, (3, 3)), |
| Activation('relu'), | Activation('relu'), |
| <pre>MaxPooling2D(pool_size=(2, 2)),</pre> | <pre>MaxPooling2D(pool_size=(2, 2)),</pre> |
| Dropout(0.25), | Dropout(0.25), |
| Conv2D(64, (3, 3), padding='same'), | Conv2D(64, (3, 3), padding='same'), |
| Activation('relu'), | Activation('relu'), |
| Conv2D(32, (3, 3)), | Conv2D(32, (3, 3)), |
| Activation('relu'), | Activation('relu'), |
| <pre>MaxPooling2D(pool_size=(2, 2)),</pre> | <pre>MaxPooling2D(pool_size=(2, 2)),</pre> |
| Dropout(0.25), | Dropout(0.25), |
| Flatten(), | Flatten(), |
| Dense(512), | Dense(512), |
| Activation('relu'), | Activation('relu'), |
| Dropout(0.5), | Dropout(0.5), |
| Dense(10), | Dense(10), |
| Activation('softmax') | Activation('softmax') |
| 1) | |
| model compile(loss='categorical crossentrony' | model compile(loss='categorical crossentropy' |
| ontimizer-tf ontimizers SCD(learning rate-0 016) | optimizer-tf optimizers SGD(learning rate-0 016) |
| metrics=['accuracy']) | metrics=['accuracy']) |
| meetics-[accuracy]/ | metrics-[accuracy]/ |
| <pre>model.fit(ds_train, epochs=40)</pre> | <pre>model.fit(ds_train, epochs=40)</pre> |

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TF2/KERAS TUTORIALS

github.com/graphcore/examples/tree/master/tutorials/tutorials/tensorflow2





INTRO TO POPTORCH

GRAPHCORE



WHAT IS POPTORCH?



| | _, ind = corch.max(predictions, i) | | _, ind = corch.max(predictions, i) |
|---|---|---|--|
| | <pre># provide labels only for samples, where prediction is available (during the training, no ions.size()[0]:]</pre> | | <pre># provide labels only for samples, where prediction is available (during the training, not labels = labels[-predictions.size()[0]:]</pre> |
| | PyTorch ^{ch.eq(ind, labels)).item() / labels.size(GPU} | | <pre>accuracy = torch.sum(torch.eq(ind, labels)).item() / labels.si IPU I00.0 return accuracy</pre> |
| | ifname == 'main': | i | fname == 'main': |
| | <pre>parser = argparse.ArgumentParser(description='MNIST training in PopTorch')</pre> | | <pre>parser = argparse.ArgumentParser(description='MNIST training in PopTorch')</pre> |
| | <pre>parser.add_argument('batch-size', type=int, default=8, help='batch size for training (default=1)</pre> | | <pre>parser.add_argument('batch-size', type=int, default=8, help='batch size for training (default=1)</pre> |
| | <pre>parser.add_argument('test-batch-size', type=int, default=8, help='batch size for testing</pre> | | <pre>parser.add_argument('test-batch-size', type=int, default=8, help='batch size for testing</pre> |
| | <pre>parser.add_argument('epochs', type=int, default=10, help='number of epochs to train (de</pre> | | <pre>parser.add_argument('epochs', type=int, default=10, help='number of epochs to train (definition)</pre> |
| | parser.add_argument('lr', type=float, default=0.05, help='learning rate (default: 0.05) | | parser.add_argument('lr', type=float, default=0.05, help='learning rate (default: 0.05)' |
| | | + | parser.add_argument('device-iterations', type=int, detault=50, help='device iterations') |
| | args = parser.parse_args() | | args – parser parse_args() |
| - | <pre>training_data = torch.utils.data.DataLoader(</pre> | + | <pre>opts = poptorch.Options().deviceIterations(args.device_iterations)</pre> |
| | /////////////////////////////////////// | + | <pre>training_data = poptorch.DataLoader(opts,</pre> |
| | torchvision.datasets.MNIST('mnist_data/', train=True, download=True, | | torchvision.datasets.MNIST('mnist_data/', train=True, download=True, trans |
| | batch_size=args.batch_size, shuffle=irue, drop_last=irue) | | batch_size=args.batch_size, shuffle=Irue, drop_last=Irue) |
| _ | torchyision_datasets_MNIST('mnist_data/'train=Falsedownload=True | Ŧ | torchyision_datasets_MNIST('mnist_data/'train=Ealsedownload=Truetrain |
| | | | |
| | <pre>model = Network()</pre> | | <pre>model = Network()</pre> |
| | <pre>training_model = TrainingModelWithLoss(model)</pre> | | <pre>training_model = TrainingModelWithLoss(model)</pre> |
| | <pre>optimizer=optim.SGD(model.parameters(), lr=args.lr)</pre> | | <pre>optimizer=optim.SGD(model.parameters(), lr=args.lr)</pre> |
| | | + | <pre>training_model = poptorch.trainingModel(training_model, opts, optimizer=optimizer)</pre> |
| | | + | <pre>inference_model = poptorch.inferenceModel(model)</pre> |
| | # Run training | | # Run training |
| | for in range(args.epochs): | | for in range(args.epochs): |
| | for data, labels in training data: | | for data, labels in training data: |
| | preds, losses = training_model(data, labels) | | <pre>preds, losses = training_model(data, labels)</pre> |
| — | <pre>optimizer.zero_grad()</pre> | + | |
| - | losses.backward() | + | # Detach the training model so that the same IPU could be used for validation |
| — | <pre>optimizer.step()</pre> | + | <pre>training_model.detachFromDevice()</pre> |
| | # Pup validation | | # Pup validation |
| | * run vacuation | | = 0.0 |
| | with torch.no grad(): | | with torch.no grad(): |
| | for data, labels in test_data: | | for data, labels in test_data: |
| — | <pre>output = model(data)</pre> | + | <pre>output = inference_model(data)</pre> |
| | <pre>sum_acc += accuracy(output, labels)</pre> | | <pre>sum_acc += accuracy(output, labels)</pre> |
| | <pre>print("Accuracy on test set: {:0.2f}%".format(sum_acc / len(test_data)))</pre> | | <pre>print("Accuracy on test set: {:0.2f}%".format(sum_acc / len(test_data)))</pre> |
| | | | |

POPTORCH TUTORIALS

github.com/graphcore/examples/tree/master/tutorials/tutorials/pytorch





UNDER THE HOOD: BSP



BULK SYNCHRONOUS PARALLEL (BSP)

BSP software bridging model – massively parallel computing with no concurrency hazards

3 phases: compute, sync, exchange

Easy to program - no live-locks or dead-locks

Widely-used in parallel computing - Google, FB, ...

First use of BSP inside a parallel processor



COMPUTATIONAL GRAPH



TIME

GRAPH EXECUTION MODEL COMMUNICATION COMPUTE SYNC COMMUNICATION COMPUTE SYNC COMMUNICATION

POPLAR FRAMEWORK



WHAT IS POPLAR?

- Parallel programming framework that targets the IPU
- Simple but powerful programming model
- Close to the metal
- General purpose, extensible



POPLAR FRAMEWORK

1. Graphs, Variables & Vertices

- 2. Compute Sets & Execution
- 3. Host IPU Execution Model



THE POPLAR GRAPH







Data is stored in the graph in fixed size multi-dimensional tensors.



VARIABLES





VERTICES



A **vertex** is a specific piece of work to be carried out.

The edges determine which variable elements are processed by the vertex. A vertex can connect to a single element or a range of elements.



3.13

0.23

VERTICES



Codelet A

Input<float> x; Input<Vector<float>> y; Output<float> z;

*z = x + sum(y);

Each vertex is associated with a **codelet**.



3.13

0.23





VERTICES



Many vertices are needed to fully utilize the device



POPLAR FRAMEWORK

- 1. Graphs, Variables & Vertices
- 2. Compute Sets & Execution
- 3. Host IPU Execution Model



COMPUTE SETS



Compute sets specify sets of vertices to execute in parallel

Poplar verifies the compute set is free of data races



A compute sets execute in 3 steps:

- **1. Exchange** Transfer inputs
- 2. Compute Run vertices in parallel
- **3. Exchange** Transfer outputs

Exchange code is generated by Poplar







Exchange is required when a vertex in a compute set needs to read or write data which is stored on another tile's memory.





For each compute set, each tile will have a number of vertices to execute.





All tiles start by **sync**ing.





The tiles then move to **exchange:** Required vertex input data is copied between memory.





Tiles will move to **compute** when they have finished exchange.

During compute vertices will read from and write to local tile memory.





Each tile processor has several independent hardware threads (workers) to execute code.

Once exchange is complete, a hardware scheduler (**supervisor**) dispatches vertices onto the workers to run.

The tiles will run all vertices and then sync.

SUMMARY

A graph is made up of:

- Data (variables in the graph)
- Compute tasks (vertices)
- Edges that connect them

Vertices:

- Are associated to a codelet (code)
- Run on a single tile

Compute sets:

- Specify sets of vertices to execute in parallel
- Are executed in 3 steps: Exchange inputs, Compute, Exchange outputs

Control program:

• Specifies the order of operations

The program resides on the chip:

• The host takes care of compilation and of the data stream preparation



THE HOST PROGRAM

Host programs use the poplar library. #include <poplar/Engine.hpp> Codelets are loaded into using namespace poplar; the graph. using namespace poplar::program; The **Graph** class is used to build up ... the computation Graph graph(target); graph. graph.addCodelets("my-codelets.cpp"); Control Program prog1, prog2; programs are The **Engine** class built up out of constructMyGraph(graph, &prog1, &prog2); represents a fully instances of compiled program Engine eng(device, graph, {prog1, prog2}); the **Program** ready to run on class. hardware. eng.run(0);



CODELET DEFINITIONS

The fields of the vertex specify its inputs, outputs and internal data.

| | Each codelet is defined as a C++ class that inherits from the |
|---|---|
| <pre>class AdderVertex : public Vertex { public: Input<float> x; Input<float> y; Output<float> z;</float></float></float></pre> | Vertex class. |
| <pre>float bias; bool compute() { *z = x + y + bias; return true; }</pre> | The compute method specifies the vertex execution behaviour. |

BUILDING THE COMPUTE GRAPH







CREATING CONTROL PROGRAMS

```
Graph g(device);
g.addCodelets("codelets.cpp");
```

```
•••
```

```
auto prog = Sequence();
prog.add(Execute(cs1));
prog.add(Execute(cs2));
```









CREATING THE ENGINE



SUMMARY

- Poplar lets you define your own operations by writing codelets
- Poplar generates "glue code" required to synchronize / exchange data
- Frees you to concentrate on parallel algorithm design





LIBRARIES = MODULAR GRAPH BUILDING





POPLIBSTM

| C / C++ and Python language bindings | | | | | | | |
|---------------------------------------|-----------------------------------|---|-----------------------------|--|--|--|--|
| poputil | popops | poplin | poprandom | popnn | | | |
| Utility functions for building graphs | Pointwise and reduction operators | Matrix multiply and convolution functions | Random number generation | Neural network functions (activation fns, pooling, loss) | | | |
| POPLAR® | | | | | | | |



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Director's Discretionary Allocation Program

The ALCF Director's Discretionary program provides "start up" awards to researchers working to achieve computational readiness for for a major allocation award.



Molecular dynamics simulations based on machine learning help scientists learn about the movement of the boundary between ice grains (yellow/green/cyan) and the stacking disorder that occurs when hexagonal (orange) and cubic (blue) pieces of ice freeze together. Image: Henry Chan and Subramanian Sankaranarayanan, Argonne National Laboratory

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general ~

charlieb 6:05 AM

🎉 Pleased to share with you all some new work from the Graphcore research team! 🎉

Our paper *Unit Scaling* introduces a new method for low-precision number formats, making FP16 We've managed to train BERT in these formats for the first time without loss scaling.

- You can find our blog post here: https://www.graphcore.ai/posts/simple-fp16-and-fp8-trainir
- Paperspace notebook (try it yourself!): https://ipu.dev/qXfm2a
- Arxiv paper: https://arxiv.org/abs/2303.11257

(& we were also featured on Davis Blalock's popular ML newsletter this week) (edited)

graphcore.ai

Simple FP16 and FP8 training with unit scaling

Unit Scaling is a new low-precision machine learning method able to train language models in FP16 and FP8 without loss scaling. (69 kB) \star



🗎 arXiv.org

Unit Scaling: Out-of-the-Box Low-Precision Training

We present unit scaling, a paradigm for designing deep learning models that simplifies the use of low-precision number formats. Training in FP16 or the recently proposed FP8 formats offers substantial efficiency gains, but can lack sufficient range for out-of-the-box training. Unit scaling addresses this by introducing a principled approach to model numerics: seeking unit variance of Show more

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Join at graphcore.ai/join-community

| | Tuesday, 11 June | |
|--|--|----------------------------------|
| 1:00 PM → 1:15 PM | Introduction | ③ 15m |
| 1:15 PM → 1:45 PM | Graphcore BowPod64 Hardware | () 30m |
| 1:45 PM → 2:30 PM | Software Stack: TensorFlow, PyTorch, and Poplar | () 45m |
| 2:30 PM → 2:45 PM | Break | () 15m |
| 2:45 PM → 3:15 PM | Porting applications with Poplar | ③ 30m |
| 3:15 PM → 4:00 PM | How to use Bow Pod64@ ALCF | ③ 45m |
| | Wednesday, 12 June | |
| 1:00 PM → 1:45 PM | Deep Dive on Graph neural networks and Large Language Models | Q.45m |
| 1:45 DM > 2:15 DM | | 0 +0111 |
| | Profiling with PonVision | () 20m |
| 2:15 PM | Profiling with PopVision | ③ 30m |
| 2:15 PM → 2:30 PM | Profiling with PopVision Break | (§ 30m) (§ 15m) |
| 2:15 PM → 2:30 PM 2:30 PM → 3:15 PM | Profiling with PopVision Break Hands-on session | (\$ 30m) (\$ 15m) (\$ 45m) |



THANK YOU

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