

Docker with GPU's in Batch

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Outline

- GPU Resources available in Batch
- How these are provisioned
- Monitoring
- Why Docker?
- Nvidia Docker Stack
- Docker Setup
- GPU Docker Setup
- Integrate GPU Docker with HTCondor
- Issues & Future Work



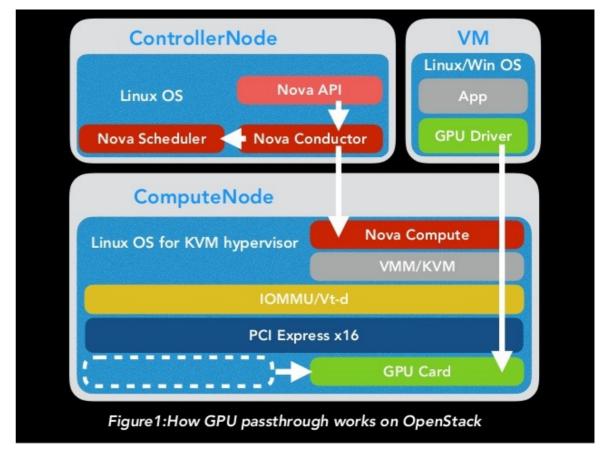
Resources available in Batch

- 16 Nvidia Tesla V100's
- Double Precision
- 32GB DDR3 Memory
- "With 640 Tensor Cores, V100 is the world's first GPU to break the 100 teraFLOPS (TFLOPS) barrier of deep learning performance." (Nvidia)





- 4 GPU's installed per server
 - 32Gb DDR3 Memory
 - 2 x Intel Xeon Silver 4114 CPU's
 - 384GB of DDR4 ECC Reg memory
 - 2x 960GB Micron 5200 Eco SSDs
- Provisioned via PCI passthrough
- Driver installed at the VM level
- 1 or 4 GPU's per VM currently configured via OpenStack Flavour – no change in driver requirements
- 1 VM per GPU but >1 GPU per VM



https://www.slideshare.net/LCChina/gpgpu-on-openstack-the-best-practice-for-gpggu-internal-cloud



- Batch GPU's → Scheduling with HTCondor
 - Machine ClassAd advertises GPU
 - Submit file 'Request_GPUs>0' adds this parameter to JobAd
 - Condor matches



```
[jfenech@lxplus723 newdocker]$ cat cpu_docker.py
#!/usr/bin/env python

import sys
import numpy as np
import tensorflow as tf

device_name = "/cpu:0"
with tf.device(device_name):
    random_matrix = tf.random.uniform(shape=[4,4], minval=0, maxval=1)
    dot_operation = tf.matmul(random_matrix, tf.transpose(random_matrix))
    sum_operation = tf.reduce_sum(dot_operation)

with tf.compat.v1.Session(config=tf.compat.v1.ConfigProto(log_device_placement=True)) as session:
        result = session.run(sum_operation)
        print(result)
```

```
[jfenech@lxplus723 newdocker]$ cat gpu.sh
#!/bin/bash
python -m virtualenv myvenv
source myvenv/bin/activate
pip install tensorflow
pip install tensorflow-gpu
python docker.py
```

```
[jfenech@lxplus723 newdocker]$ cat gpu.sub
                     = vanilla
universe
executable
                     = ./gpu.sh
arguments
                     = /etc/hosts
should_transfer_files = YES
when_to_transfer_output = ON_EXIT
                     = out/$(ClusterId).$(ProcId).out
output
error
                    = err/$(ClusterId).$(ProcId).err
log
                     = log/$(ClusterId).$(ProcId).log
+JobFlavour
                     = "espresso"
request_gpus
queue 1
```



```
[jfenech@lxplus723 newdocker]$ cat gpu_docker.py
#!/usr/bin/env python

import sys
import numpy as np
import tensorflow as tf

device_name = "/gpu:0"
with tf.device(device_name):
    random_matrix = tf.random.uniform(shape=[4,4], minval=0, maxval=1)
    dot_operation = tf.matmul(random_matrix, tf.transpose(random_matrix))
    sum_operation = tf.reduce_sum(dot_operation)

with tf.compat.v1.Session(config=tf.compat.v1.ConfigProto(log_device_placement=True)) as session:
        result = session.run(sum_operation)
        print(result)
```

```
[jfenech@lxplus723 newdocker]$ cat gpu_docker.sub
universe
                      = docker
docker_image
                     = tensorflow/tensorflow:1.15.0-gpu
executable
                      = ./gpu_docker.py
arguments
                      = /etc/hosts
should_transfer_files = YES
when_to_transfer_output = ON_EXIT
output
                      = out/$(ClusterId).$(ProcId).out
                      = err/$(ClusterId).$(ProcId).err
error
log
                      = log/$(ClusterId).$(ProcId).log
                      = "espresso"
+JobFlavour
request_gpus
                      = Machine == "b7g47n0009.cern.ch"
requirements
queue 1
```



- GPU's monitored at VM level via Collectd-cuda plugin now in Production
- Collects information via Nvidia Service Management Interface
- Current month's logs visible in Cuda Monitoring Grafana dashboard
- Logs stored in S3 available via Spark in SWAN
- Metrics available include temperature, utilisation, clocks, power
- Later HTCondor versions (>8.8) include enhanced per job GPU monitoring
 - Average Usage
 - Peak Memory Usage



[root@b7g47n0008 ~]# nvidia-smi Thu Jan 30 16:58:35 2020

Perf	Persis Pwr:Us	tence-M age/Cap	Bus-Id	Disp.A Memory-Usage	Volatile GPU-Util	Uncorr. ECC Compute M.
V100- P0	PCIE 27W	Off / 250W	0000000 12M	0:00:05.0 Off iB / 32510MiB	0%	0 Default
						+
PID	Туре	Process	name			GPU Memory Usage
	Perf ====== V100- P0	Persis Perf Pwr:Us: V100-PCIE P0 27W	Persistence-M Perf Pwr:Usage/Cap =======+ V100-PCIE Off P0 27W / 250W	Persistence-M Bus-Id Perf Pwr:Usage/Cap 	Persistence-M Bus-Id Disp.A Perf Pwr:Usage/Cap Memory-Usage V100-PCIE Off 00000000:00:05.0 Off P0 27W / 250W 12MiB / 32510MiB	



Percent of time over the past sample between 1 second and 1/6 second period during which one or more kernels was executing on the GPU



[root@b7g47n0008 ~]# nvidia-smi Thu Jan 30 16:58:35 2020



https://monit-grafana.cern.ch/d/0EfSgd4Zz/cuda-monitoring?orgId=5&from=now%2FM&to=now

31/01/2020



Utilisation

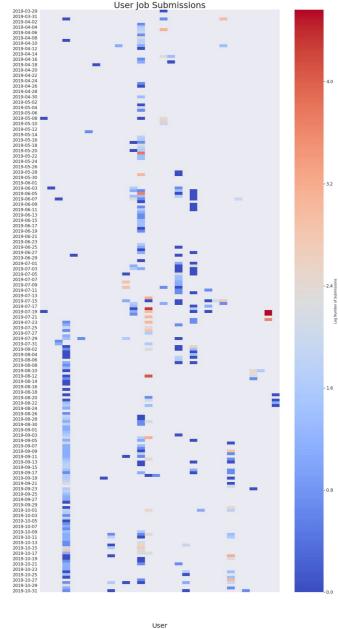
Percent of time over the past

sample between 1 second and 1/6 second period during which

[root@b7g47n0008 ~]# nvidia-smi Thu Jan 30 16:58:35 2020



https://monit-grafana.cern.ch/d/0EfSgd4Zz/cuda-monitor orgId=5&from=now%2FM&to=now





Why Docker?

- Reproducible environment independent of host
- Let user manage requirements and dependencies
- Variety of GPU use-cases and frameworks







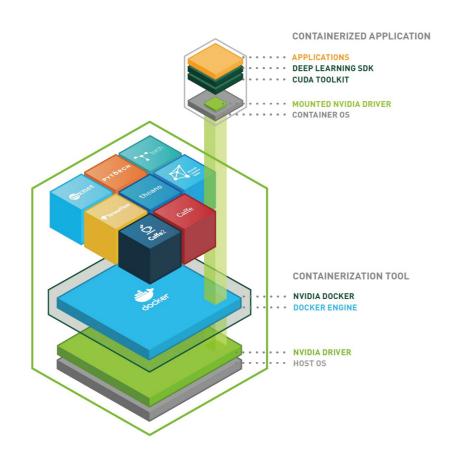
Why Docker?

Benchmarking!

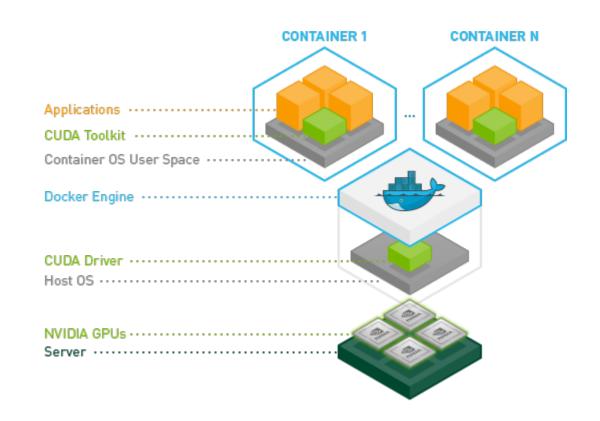


NVIDIA Docker Structure

Nvidia ≡ Cuda



https://dgx-wiki.readthedocs.io/en/latest/docs/docker/service.html#nvidia-docker



https://www.slideshare.net/LCChina/gpgpu-on-openstack-the-best-practice-for-gpggu-internal-cloud



Setup GPU Docker

- Install libnvidia-container, nvidia-container-runtime, nvidia-container-runtime-hook
- GPU integration is better in later versions of Docker
- CentOS repo ships with Docker version: 1.13
- Latest Docker Version: 19.03
- → Upgrade Docker
- Docker → Docker-ce
- Docker-ce requires containerd.io, which conflicts with earlier Docker versions
- → Remove Docker → Install containerd.io, docker-ce, docker-ce-cli
- Add Docker to Condor user group to give required permissions



Integrate GPU Docker with HTCondor

- Docker requires 2 flags/parameters in order to run on GPU, "--gpus all"
 - docker run –gpus all tensorflow/tensorflow:latest-gpu
- HTCondor 8.9.1 introduced greater GPU Docker integration:
- GPU's mounted automatically in (GPU + Docker) jobs
- DOCKER_EXTRA_ARGUMENTS flag allows the crucial '--gpus all' to be added to the docker create command that Condor utilises
- Improved GPU job monitoring with the additional metrics (average usage, peak memory usage)
- Upgrade HTCondor (from default 8.8.6) on GPU machines
- Requires dev repo
- Also need to ensure Machine ClassAd contains "HasDocker" flag



Issues & Future Work

- Varied workloads but single hardware
 - Persistent low utilisation
 - Some algorithms (e.g. LSTM's) struggle to make full use of GPU
 - Making full use of Tensor Cores depends on user providing precisely shaped inputs (well, divisible by 8 (for FP16 data) or 16 (INT8 data)) to match the GPU architecture
- Fairsharing limited number of resources
 - Many jobs are long (nextweek), blocking larger number of short jobs
- Can Docker manage GPU sharing??
 - Harness slack GPU resources
- Machine learning vs SixTrack workloads



Batchdocs GPU Tutorial

Grafana GPU Dashboard

