

# WHY CONTAINERS?

#### **Benefits of Containers:**

Simplify deployment of GPU-accelerated software, eliminating time-consuming software integration work

Isolate individual deep learning frameworks and applications

Share, collaborate, and test applications across different environments

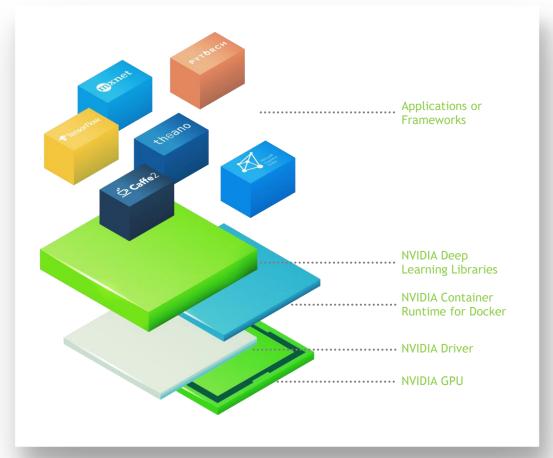


## CHALLENGES WITH COMPLEX SOFTWARE

Current DIY GPU-accelerated AI and HPC deployments are complex and time consuming to build, test and maintain

Development of software frameworks by the community is moving very fast

Requires high level of **expertise** to manage driver, library, framework dependencies



## **NVIDIA GPU CLOUD**

Simple access to a comprehensive catalog of GPU-accelerated software

#### **Discover 30 GPU-Accelerated Containers**

Deep learning, third-party managed HPC applications, NVIDIA HPC visualization tools, and partner applications

#### Innovate in Minutes, Not Weeks

Get up and running quickly and reduce complexity

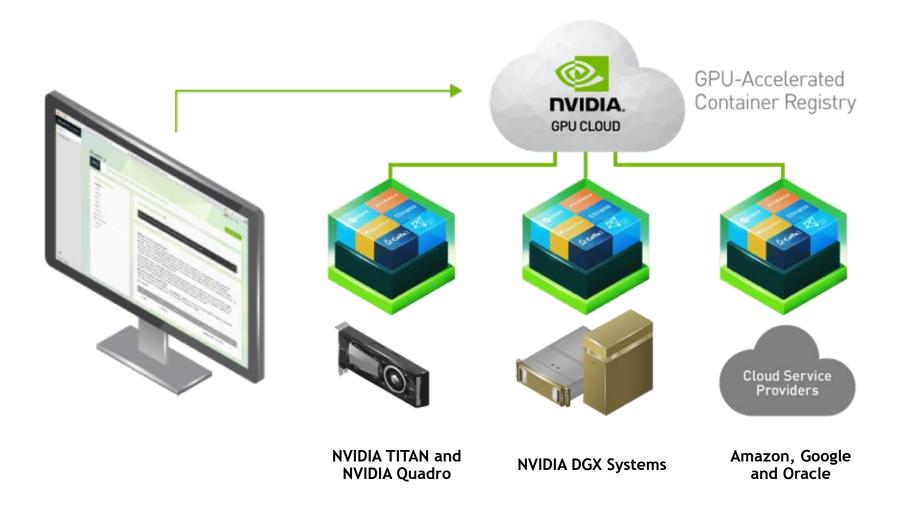
#### **Access from Anywhere**

Use on PCs with NVIDIA Volta or Pascal™ architecture GPUs, NVIDIA DGX Systems, and supported cloud providers





# FROM DESKTOP, TO DATA CENTER, TO CLOUD



# NGC GPU-OPTIMIZED DEEP LEARNING CONTAINERS

A comprehensive catalog of deep learning software

- NVCaffe
- Caffe2
- Microsoft Cognitive Toolkit (CNTK)
- ▶ DIGITS
- MXNet
- PyTorch

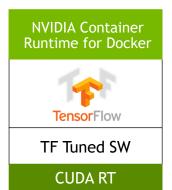
- TensorFlow
- Theano
- Torch
- CUDA (base level container for developers)
- NVIDIA TensorRT inference accelerator with ONNX support



## **ALWAYS UP-TO-DATE**

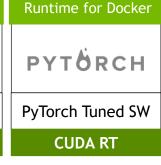
#### Monthly updates from NVIDIA to deep learning containers

#### **Containerized Applications**

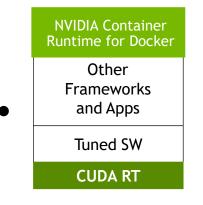








**NVIDIA Container** 



Linux Kernel + CUDA Driver

# **END-TO-END PRODUCT FAMILY**

#### **TRAINING**





#### **INFERENCE**







## GET STARTED TODAY WITH NGC

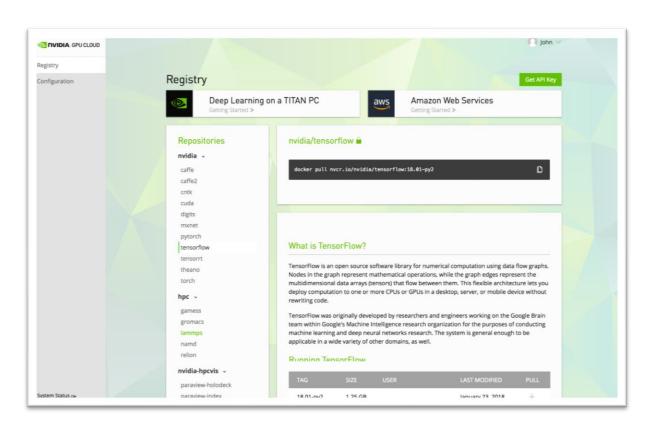
Sign up for no cost access

To learn more about all of the GPU-accelerated software on NVIDIA GPU Cloud, visit:

nvidia.com/cloud

To sign up, go to:

nvidia.com/ngcsignup



## NGC CONTAINERS ON CSCS

In just 1,2,3.....

CSCS uses shifter to manage containers. Shifters essentially does not allow root access and maintains user privileges.

Run your first NGC Tensorflow container in 3 simple steps

- 1) shifter --login docker pull nvcr.io/nvidia/tensorflow:18.03-py3
- 2) Username: \$oauthtoken Password: <API\_KEY>
- 3) srun -N1 -C gpu --pty shifter run --mount=type=bind,source=\$HOME,destination=\$HOME nvcr.io/nvidia/tensorflow:18.03-py3 bash

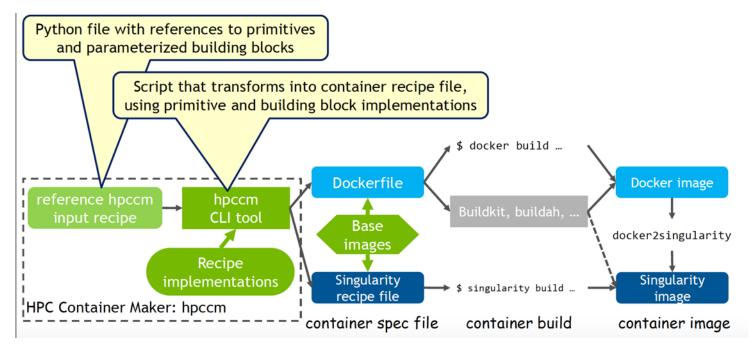




# **HPC CONTAINER MAKER - HPCCM**

"h-p-see-um"

- HPC Container Maker (HPCCM) generates container specification files (Dockerfiles or Singularity recipe) based on recipes
- A recipe specifies the series of steps to be performed when building a container



## **HPC CONTAINER MAKER - HPCCM**

- Container implementation abstraction
  - The same recipe file generates specification files for Docker or Singularity
- Availability of full programming languages
  - A recipe is Python code. This means that you can use the full power of Python in a recipe for conditional branching, input validation, searching the web for the latest version of a component, etc.
- Higher level abstraction
  - Provides building blocks to simplify recipes and encapsulate best practices
- Container Maker generates human readable Dockerfiles and Singularity recipe files

## **FEATURES**

#### **BASE IMAGE & PACKAGES**

- Ubuntu 16.04
- Cent OS

#### **COMPILER**

- PGI COMMUNITY EDITION
- GNU
- NVCC (CUDA TOOLKIT)

#### **SOFTWARE - MORE ON REQUEST**

- CMAKE
- HDF5
- NETCDF
- MLNX OFED
- OpenMPI
- MVAPICH2

- Python
- FFTW
- MKL

#### **FULL SCALE APPLICATIONS**

- MILC
- GROMACS

```
# Choose a base image
Stage0.baseimage('ubuntu:16.04')
# Install GNU Compilers
Stage0 += apt_get(ospackages=['gcc', 'g++', 'gfortran'])
                                                        hpccm.py --recipe recipes/basic.py --format singularity
hpccm.py --recipe recipes/basic.py --format docker
FROM ubuntu: 16.04
                                                          BootStrap: docker
RUN apt-get update -y && \
                                                          From: ubuntu:16.04
apt-get install -y --no-install-recommends \
                                                          %post
gcc \
                                                          apt-get update -y
g++ \
                                                          apt-get install -y --no-install-recommends \
gfortran && \
                                                          gcc \
rm -rf /var/lib/apt/lists/*
                                                          g++ \
                                                          gfortran
                                                          rm -rf /var/lib/apt/lists/*
```

# AVAILABILITY OF A FULL PROGRAMMING LANGUAGE

 Full power of Python in a recipe for conditional branching, input validation, searching the web for the latest version of a component, etc.

For example, the LAMMPS application may be built in single, double, or mixed precision mode. (hpccm.py --userarg LAMMPS\_PRECISION=...)

```
# get and validate precision
VALID_PRECISION = ['single', 'double', 'mixed']
precision = USERARG.get('LAMMPS_PRECISION', 'single')
if precision not in VALID_PRECISION:
    raise ValueError('Invalid precision')
...

Stage0 += shell(commands=['make -f Makefile.linux.{}'.format(precision), ...])
...
```

## HIGHER LEVEL ABSTRACTION

Building blocks to simplify recipes and encapsulate best practices

```
Stage0 += openmpi(cuda=True, infiniband=True, prefix='/usr/local/openmpi', version='3.0.0')
```

```
# OpenMPI version 3.0.0 RUN
apt-get update -y && \
apt-get install -y --no-install-recommends \
 file \
 hwloc \
 openssh-client \
 wget && \
rm -rf /var/lib/apt/lists/*
RUN mkdir -p /tmp && wget -q --no-check-certificate -P /tmp https://www.open-
mpi.org/software/ompi/v3.0/downloads/openmpi-3.0.0.tar.bz2 && \
 tar -x -f /tmp/openmpi-3.0.0.tar.bz2 -C /tmp -j && \
 cd /tmp/openmpi-3.0.0 && ./configure --prefix=/usr/local/openmpi --disable-
getpwuid --enable-orterun-prefix-by-default --with-cuda --with-verbs && \
 make -i4 && \
 make -j4 install && \
 rm -rf /tmp/openmpi-3.0.0.tar.bz2 /tmp/openmpi-3.0.0
ENV PATH=/usr/local/openmpi/bin:$PATH \
LD_LIBRARY_PATH=/usr/local/openmpi/lib:$LD_LIBRARY_PATH
```

```
# OpenMPI version 3.0.0
%post
 apt-get update -y
 apt-get install -y --no-install-recommends \
   file \
   hwloc \
   openssh-client \
   wget
 rm -rf /var/lib/apt/lists/*
%post
 mkdir -p /tmp && wget -q --no-check-certificate -P /tmp https://www.open-
mpi.org/software/ompi/v3.0/downloads/openmpi-3.0.0.tar.bz2
tar -x -f /tmp/openmpi-3.0.0.tar.bz2 -C /tmp -j
 cd /tmp/openmpi-3.0.0 && ./configure --prefix=/usr/local/openmpi --disable-getpwuid
--enable-orterun-prefix-by-default --with-cuda --with-verbs
 make -i4
make -j4 install
 rm -rf /tmp/openmpi-3.0.0.tar.bz2 /tmp/openmpi-3.0.0
%environment
export PATH=/usr/local/openmpi/bin:$PATH
 export LD_LIBRARY_PATH=/usr/local/openmpi/lib:$LD_LIBRARY_PATH
```

# **TOOLCHAIN SUPPORT**

#### Using toolchain for building blocks

```
GNU = gnu()
Stage0 += GNU
TC = GNU.toolchain
TC.CUDA_HOME = /usr/local/cuda

OMPI = openmpi(version='3.0.0',toolchain=TC)
Stage0 += OMPI

FFTW = fftw(version='3.3.7',toolchain=TC)
Stage0 += FFTW

HDF5 = hdf5(version='1.10.1',toolchain=TC)
Stage0 += HDF5

NETCDF =netcdf(version='4.6.1',toolchain=TC)
Stage0 += NETCDF
```

## USING HPCCM FOR CSCS

#### ON LOCAL LINUX BOX:

- 1) Git clone <a href="https://github.com/NVIDIA/hpc-container-maker">https://github.com/NVIDIA/hpc-container-maker</a>
- 2) ./hpccm -format docker -recipe hpcbase-gnu-mvapich2.py > Dockerfile
- 3) Docker build -t container < Dockerfile
- 4) Docker save container -o container.tar.gz

ON CSCS: (https://user.cscs.ch/user\_services/containers/advanced\_shifter/)

- 1) Srun -N 1 -C gpu shifter load container.tar.gz container
- 2) srun -N1 -C gpu --pty shifter run --mount=type=bind,source=\$HOME,destination=\$HOME load/library/container:latest bash

