

- Go to the NVIDIA website and find the latest driver version compatible with your graphics card, if you have a NVIDIA GeForce, go to: http://www.geforce.com/drivers
- Check the latest version compiled in "Proprietary GPU Drivers" PPA which also matches your graphics card in: https://launchpad.net/~graphics-drivers/+archive/ubuntu/ppa. For stability, you may want to use the recommended driver in that page.
- Add the "Proprietary GPU Drivers" PPA repository:

```
sudo add-apt-repository ppa:graphics-drivers/ppa
```

· Update the repositories database:

```
sudo apt-get update
```

• Install your selected version of the graphics card (this could break your graphic system, be sure to know how to use a shell to be able to repair it), for example:

```
sudo apt-get install nvidia-352
```

· Restart your system:

```
sudo shutdown -r now
```

• Check the NVIDIA drivers installed:

```
cat /proc/driver/nvidia/version
```

Fix flickering

If your screen is flickering, you can solve it with the following steps.

Note: the following may not apply to you.

• Install Compiz Config:

```
sudo apt-get install compizconfig-settings-manager
```

- From the launcher, execute CompizConfig Settings Manager
- Check the checkbox in "Utility -> Workarounds -> Force full screen redraws (buffer swap) on repaint"

CUDA

• Check that your computer is CUDA enabled (you should have a compatible NVIDIA graphics card):

```
lspci | grep -i nvidia
```

- You need to have an account or register in the NVIDIA developer program (it's free): https://developer.nvidia.com/user
- Download CUDA 7 from: https://developer.nvidia.com/cuda-downloads, you can download the "deb (network)" package (cuda repo).
- · Go to the downloads directory:

```
cd ~/Downloads
```

• Install the package (the repository):

```
sudo dpkg -i cuda-repo-ubuntu1404*amd64.deb
```

· Update the repositories DB:

```
sudo apt-get update
```

• Install CUDA (it should list cuda packages):

```
sudo apt-get install cuda
```

• Add CUDA to the PATH variable (and make it permanent for the current user):

```
echo 'export PATH=/usr/local/cuda/bin:$PATH' >> ~/.bashrc
```

• Add CUDA to the LD LIBRARY PATH variable (and make it permanent for the current user):

```
echo 'export LD_LIBRARY_PATH=/usr/local/cuda/lib64:$LD_LIBRARY_PATH' >> ~/.bashrc
```

• Log out and log in again to activate the new variables (close and open the terminal).

Check CUDA installation

• To check the correct installation running the samples, install the samples in a directory under your home:

```
cuda-install-samples-7.0.sh ~/cuda-samples
```

• You can check the version of the NVIDIA drivers installed with:

```
cat /proc/driver/nvidia/version
```

· You can check the version of CUDA installed with:

```
nvcc -V
```

• Go to the samples directory:

```
cd ~/cuda-samples/NVIDIA*Samples
```

• Build the samples:

```
make -j $(($(nproc) + 1))
```

Note: you could just run make, but that last part (-j (((nproc) + 1))) executes the make command in parallel, using the number of cores in your machine (plus one), so the compilation would finish faster.

• Restart your computer:

```
sudo shutdown -r now
```

• Check the CUDA installation, run:

```
bin/x86_64/linux/release/deviceQuery
```

- Check that it detects a CUDA device.
- Check that it detects your graphics card.
- Check at the end that it says that the test passed.
- Run the bandwidth test to check that the system and the CUDA device are capable to communicate:

bin/x86_64/linux/release/bandwidthTest

Install OpenBLAS

Note: you could just use "ATLAS" with: sudo apt-get install libatlas-base-dev. But you would probably get a better performance from OpenBLAS following the next instructions, and it's quite straightforward.

• Install Fortran compiler:

```
sudo apt-get install gfortran
```

• Create a "code" directory to store your code:

Note: you can put your code wherever you want, so, modify this part as you wish, but I will assume that your are doing it.

```
mkdir ~/code
```

• Enter that directory:

```
cd ~/code
```

• Clone OpenBLAS:

```
git clone https://github.com/xianyi/OpenBLAS.git
```

• Go to the OpenBLAS directory:

```
cd OpenBLAS
```

· Compile OpenBLAS:

```
make -j $(($(nproc) + 1))
```

Note: you could just run make, but that last part (-j (((nproc) + 1))) executes the make command in parallel, using the number of cores in your machine (plus one), so the compilation would finish faster.

· Install OpenBLAS:

```
sudo make PREFIX=/usr/local install
```

Install Boost

```
sudo apt-get install libboost-all-dev
```

Install OpenCV

sudo apt-get install libopencv-dev

Install: protobuf, glog, gflags

sudo apt-get install libprotobuf-dev libgoogle-glog-dev libgflags-dev protobuf-compiler

Install IO libraries: hdf5, leveldb, snappy, lmdb

sudo apt-get install libhdf5-serial-dev libleveldb-dev libsnappy-dev liblmdb-dev

Install Anaconda Python

- Go to the Anaconda website and download Anaconda for Linux (you would probably want the 64x version): http://continuum.io/downloads
- · Go to the downloads directory:

cd ~/Downloads

• Execute the installer:

bash Anaconda*.sh

- When asked if add Anaconda to the PATH for the current user, type "yes".
- Log out and log in again to activate the new variables (close and open the terminal).

Install HDF5

• Install HDF5 with conda, although it will only be used by Anaconda:

conda install hdf5

Configure the HDF5 version

Note: this is a quick fix for a minor bug / issue with the version of libhdf5. If you know a better / proper way to solve it, let me know. If this section doesn't apply to you, omit it.

• Go to the libraries directory:

cd /usr/lib/x86_64-linux-gnu

• Link the system version of HDF5 from 7 to 9:

```
sudo ln -s libhdf5.so.7 libhdf5.so.9
sudo ln -s libhdf5_hl.so.7 libhdf5_hl.so.9
```

• Update the "Dynamic Linker":

sudo ldconfig

Install CuDNN

Note: for this to work, you need a NVIDIA GPU based on the "Kepler" architeture (or the newer, "Maxwell" architecture).

Note: although this guide includes all the build, installation and configuration of CuDNN, only the build and installation was tested, the usage hasn't been tested. (It was only tested on a NVIDIA GT 525m and a NVIDIA GTX 580 which have a "Fermi" (older) architecture, and is not supported by CuDNN.

- You need to have an account or register in the NVIDIA developer program (it's free): https://developer.nvidia.com/user
- Using your account, download CuDNN: https://developer.nvidia.com/cuDNN
- Go to the downloads directory:

```
cd ~/Downloads/
```

· Uncompress the CuDNN download:

```
tar xvf cudnn*.tgz
```

• Enter the uncompressed directory:

```
cd cuda
```

· Copy the *.h files to your CUDA installation:

```
sudo cp */*.h /usr/local/cuda/include/
```

• Copy the .so files to your CUDA installation:

```
sudo cp */*.so* /usr/local/cuda/lib64/
```

Configure and compile Caffe

• Create a "code" directory to store your code:

Note: you can put your code wherever you want, so, modify this part as you wish, but I will assume that your are following the tutorial.

```
mkdir ~/code
```

• Enter that directory:

cd ~/code

• Clone Caffe

```
git clone https://github.com/BVLC/caffe.git
```

• Enter the new Caffe directory

```
cd caffe
```

· Copy the Makefile.config

```
cp Makefile.config.example Makefile.config
```

• Enable CuDNN, set the USE_CUDNN := 1 flag in Makefile.config:

Note: this will only work with a GPU with Kepler or newer architecture (Maxwell). This tutorial was tested on an older GPU, so this part was omitted.

```
sed -i 's/# USE_CUDNN := 1/USE_CUDNN := 1/' Makefile.config
```

• Enable OpenBLAS:

Note: if you didn't install OpenBLAS and installed ATLAS, omit this part.

```
sed -i 's/BLAS := atlas/BLAS := open/' Makefile.config
```

• Enable Anaconda:

```
sed -i 's|# ANACONDA_HOME := $(HOME)/anaconda|ANACONDA_HOME := $(HOME)/anaconda|' Makefile.config
sed -i 's|# PYTHON_INCLUDE := $(ANACONDA_HOME)|PYTHON_INCLUDE := $(ANACONDA_HOME)|' Makefile.config
sed -i 's|# $(ANACONDA_HOME)|$(ANACONDA_HOME)|' Makefile.config
sed -i 's|# PYTHON_LIB := $(ANACONDA_HOME)|PYTHON_LIB := $(ANACONDA_HOME)|' Makefile.config
sed -i 's|# WITH_PYTHON_LAYER := 1|WITH_PYTHON_LAYER := 1|' Makefile.config
```

• Install Python dependencies using conda:

```
for req in $(cat python/requirements.txt); do conda install $req; done
```

• Install Python dependencies using pip (to install what couldn't be installed by conda):

```
pip install -r python/requirements.txt
```

Remove had Anaconda libm:

Note: this is a fix for a common bug with Anaconda. It may not apply to you.

```
mkdir ~/anaconda/lib/libm.orig.d
mv ~/anaconda/lib/libm.* ~/anaconda/lib/libm.orig.d
```

• Build Caffe:

```
make all -j $(($(nproc) + 1))
```

· Build the tests:

```
make test -j $(($(nproc) + 1))
```

• Run the tests:

```
make runtest -j $(($(nproc) + 1))
```

- After running make runtest , check that all the tests passed.
- Build PyCaffe (Python interface):

```
make pycaffe -j $(($(nproc) + 1))
```

Note: you could just run make all or make test or make runtest or make pycaffe respectively, but that last part (-j \$((\$(nproc) + 1))) executes the make command in parallel, using the number of cores in your machine (plus one), so the compilation would finish faster.

• Export a variable CAFFE_ROOT with your Caffe root directory and make it permanent:

```
echo "export CAFFE_ROOT=$(pwd)" >> ~/.bashrc
```

• Add Caffe to your PYTHONPATH variable using your CAFFE_ROOT:

Technical Note: you could add Caffe to your PYTHONPATH directly, but by using a CAFFE_ROOT variable, you could move your Caffe directory and would only have to modify your CAFFE_ROOT variable. And you could use the same to export Caffe binaries to your PATH variable.

echo 'export PYTHONPATH=\$CAFFE_ROOT/python:\$PYTHONPATH' >> ~/.bashrc

- Log out and log in again to activate the new variables (close and open the terminal).
- To test that PyCaffe is working open IPython:

ipython

• Import "caffe":

import caffe

• Check that there are no errors.

Congratulations, you have an up and running Caffe installation.

Now you may want to follow one of the tutorials:

- LeNet MNIST Tutorial
- ImageNet tutorial

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