

Magic Tools to Install / Manage Software

Part 1: CONDA Virtual Environment

Jason Li

HPC User Services

LSU HPC / LONI

sys-help@loni.org

Louisiana State University, Baton Rouge

Oct 25, 2023

Magic Tools to Install / Manage Software

Part 1:  **CONDA** Virtual Environment

Part 2:  **ingularity Container**

1. Why Conda?

- 1) Problems
- 2) Virtual environment & Conda

2. Basic Usage

- 1) Get Conda
- 2) Typical workflow
- 3) Creating a virtual environment
- 4) Installing software packages

3. Advanced Tips

1. Why Conda?

- 1) Problems
- 2) Virtual environment & Conda

2. Basic Usage

- 1) Get Conda
- 2) Typical workflow
- 3) Creating a virtual environment
- 4) Installing software packages



3. Advanced Tips

- **Core problem:**


Installing software on an HPC system

- **Traditional Linux solution:**
 - Compiling from source code

a) Dependencies (Welcome to Linux!)



UNIVERSITÉ
DE GENÈVE
FACULTÉ DE MÉDECINE



SIB
Swiss Institute of
Bioinformatics

BUSCO

from QC to gene prediction and phylogenomics

BUSCO v5.4.7 is the current stable version!

[Gitlab](#), a [Conda package](#) and [Docker container](#) are also available.

Based on evolutionarily-informed expectations of gene content of near-universal single-copy orthologs, BUSCO metric is complementary to technical metrics like N50.

a) Dependencies (Welcome to Linux!)

Third-party components

A full installation of BUSCO requires *Python 3.3+* (2.7 is not supported from v4 onwards), *BioPython*, *pandas*, *BBMap*, *tBLASTn 2.2+*, *Augustus 3.2+*, *Prodigal*, *Metaeuk*, *HMMER3.1+*, *SEPP*, and *R + ggplot2* for the plotting companion script. Some of these tools are necessary only for analysing certain type of organisms and input data, or for specific run modes.

- <https://biopython.org/>
- <https://pandas.pydata.org/>
- <https://jgi.doe.gov/data-and-tools/software-tools/bbtools/>
- <https://ftp.ncbi.nlm.nih.gov/blast/executables/blast+/LATEST>
- <http://bioinf.uni-greifswald.de/augustus/>
- <https://github.com/soedinglab/metaeuk>
- <https://github.com/hyattpd/Prodigal>
- <http://hmmer.org/>
- <https://github.com/smirarab/sepp/>
- <https://www.r-project.org/>

Please make sure that each software package listed above works INDEPENDENTLY of BUSCO before attempting to run any BUSCO assessments.

a) Dependencies (Welcome to Linux!)

Third-party components

A full installation of BUSCO requires *Python 3.3+* (2.7 is not supported from v4 onwards), *BioPython*, *pandas*, *BBMap*, *tBLASTn 2.2+*, *Augustus 3.2+*, *Prodigal*, *Metaeuk*, *HMMER3.1+*, *SEPP*, and *R + ggplot2* for the plotting companion script. Some of these tools are necessary only for analysing certain type of organisms and input data, or for specific run modes.

- <https://biopython.org/>
- <https://pandas.pydata.org/>
- <https://jgi.doe.gov/data-and-tools/software-tools/bbtools/>
- <https://ftp.ncbi.nlm.nih.gov/blast/executables/blast+/LATEST>
- <http://bioinf.uni-greifswald.de/augustus/>
- <https://github.com/soedinglab/metaeuk>
- <https://github.com/hyattpd/Prodigal>
- <http://hmmer.org/>
- <https://github.com/smirarab/sepp/>
- <https://www.r-project.org/>

Please make sure that each software package listed above works INDEPENDENTLY of BUSCO before attempting to run any BUSCO assessments.

a) Dependencies (Welcome to Linux!)

Third-party components

A full installation of BUSCO requires *Python 3.3+* (2.7 is not supported from v4 onwards), *BioPython*, *pandas*, *BBMap*, *tBLASTn* 2.2+, *Augustus* 3.2+, *Prodigal*, *Metaeuk*, *HMMER3.1+*, *SEPP*, and *R + ggplot2* for the plotting companion script. Some of these tools are necessary only for analysing certain type of organisms and input data, or for specific run modes.

- <https://biopython.org/>
- <https://pandas.pydata.org/>
- <https://jgi.doe.gov/data-and-tools/software-tools/bbtools/>
- <https://ftp.ncbi.nlm.nih.gov/blast/executables/blast+/LATEST>
- <http://bioinf.uni-greifswald.de/augustus/>
- <https://github.com/soedinglab/metaeuk>
- <https://github.com/hyattpd/Prodigal>
- <http://hmmer.org/>
- <https://github.com/smirarab/sepp/>
- <https://www.r-project.org/>

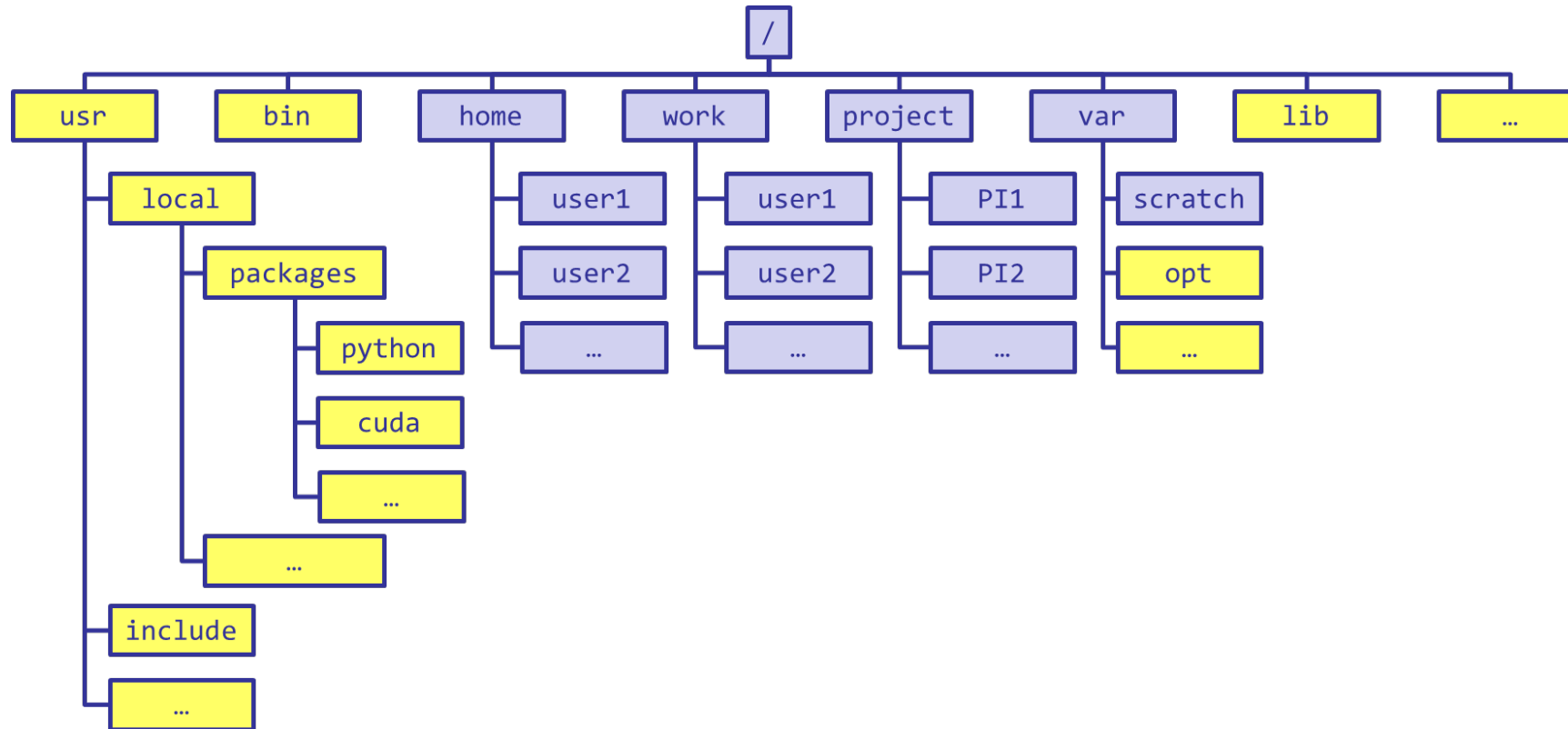
Please make sure that each software package listed above works INDEPENDENTLY of BUSCO before attempting to run any BUSCO assessments.



- Dependencies
- The following dependencies are required for AUGUSTUS:
- for gzip compressed input: (set ZIPINPUT = false in `common.mk` if this feature is not available)
 - libboost-iostreams-dev
 - zlib1g-dev
 - for comparative AUGUSTUS (multi-species, CGP): (set COMPGENEPRED = false in `common.mk` if this feature is not available. Augustus can then only be run in single-genome mode, which is what most users need.)
 - libgsl-dev
 - libboost-all-dev
 - libsuitesparse-dev
 - liblsolve55-dev
 - libsqlite3-dev (add SQLITE = false to `common.mk` if this feature is not required or the required library is not available)
 - libmysql++-dev (add MYSQL = false to `common.mk` if this feature is not required or the required library is not available)
 - for compiling utilities bam2hints and filterBam:
 - libbamtools-dev zlib1g-dev
 - for compiling utility utrnanseq:
 - libboost-all-dev (version must be > Boost_1_49_0)
 - for compiling utility bam2wig:
 - Follow [these instructions](#). Note that it shouldn't be a problem to compile AUGUSTUS without bam2wig. In practice, you can simply use `bamToWig.py` to accomplish the same task.
 - For compiling homgenemapping (set BOOST = FALSE in `auxprogs/homgenemapping/src/Makefile` if the option --printHomologs is not required or the required libraries are not available)
 - libboost-all-dev
 - for scripts:
 - Perl
 - Python3
 - for the python3 script `bamToWig.py`:
 - twoBitInfo and faToTwoBit from <http://hgdownload.soe.ucsc.edu/admin/exe> . `bamToWig.py` will automatically download these tools to the working directory during execution if they are not in your \$PATH.
 - SAMtools (available e.g. via package managers or [here](#) - see notes below)

1) Problems

b) Permission denied (Welcome to HPC!)



b) Permission denied (Welcome to HPC!)

```
[jasonli3@smic2 ~]$ module load python/3.6.2-anaconda-tensorflow
[jasonli3@smic2 ~]$ module li
Currently Loaded Modulefiles:
 1) python/3.6.2-anaconda-tensorflow
```

b) Permission denied (Welcome to HPC!)

```
[jasonli3@smic2 ~]$ module load python/3.6.2-anaconda-tensorflow
[jasonli3@smic2 ~]$ module li
Currently Loaded Modulefiles:
 1) python/3.6.2-anaconda-tensorflow
[jasonli3@smic2 ~]$ pip install geos
Collecting geos
  Downloading https://files.pythonhosted.org/packages/49/5b/b8acf74c01187a36aa41b6523deb9baa59c
100% |████████████████████████████████████████████████████████████████████████████████| 409kB 3.0MB/s
```

1) Problems

b) Permission denied (Welcome to HPC!)

```
[jasonli3@smic2 ~]$ module load python/3.6.2-anaconda-tensorflow
[jasonli3@smic2 ~]$ module li
Currently Loaded Modulefiles:
  1) python/3.6.2-anaconda-tensorflow
[jasonli3@smic2 ~]$ pip install geos
Collecting geos
  Downloading https://files.pythonhosted.org/packages/49/5b/b8acf74c01187a36aa41b6523deb9baa59c
100% |████████████████████████████████████████████████████████████████████████████████| 409kB 3.0MB/s
File "/usr/local/packages/python/3.6.2-anaconda/lib/python3.6/site-packages/pip/_utils/_
os.makedirs(path)
File "/usr/local/packages/python/3.6.2-anaconda/lib/python3.6/os.py", line 220, in makedirs
  mkdir(name, mode)
PermissionError: [Errno 13] Permission denied: '/usr/local/packages/python/3.6.2-anaconda/lib/p
You are using pip version 9.0.1, however version 23.0.1 is available.
You should consider upgrading via the 'pip install --upgrade pip' command.
[jasonli3@smic2 ~]$
```

b) Permission denied (Welcome to HPC!)

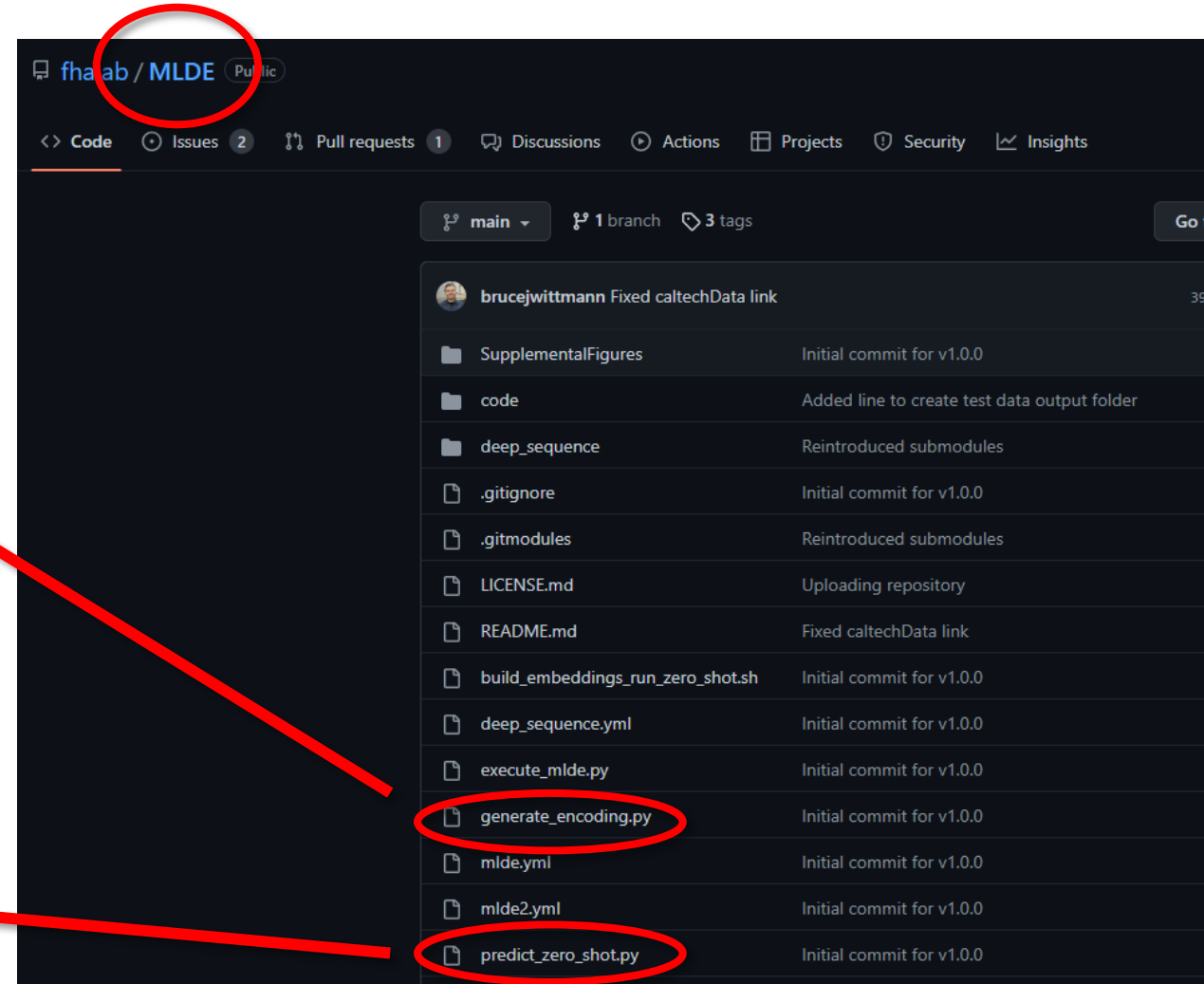
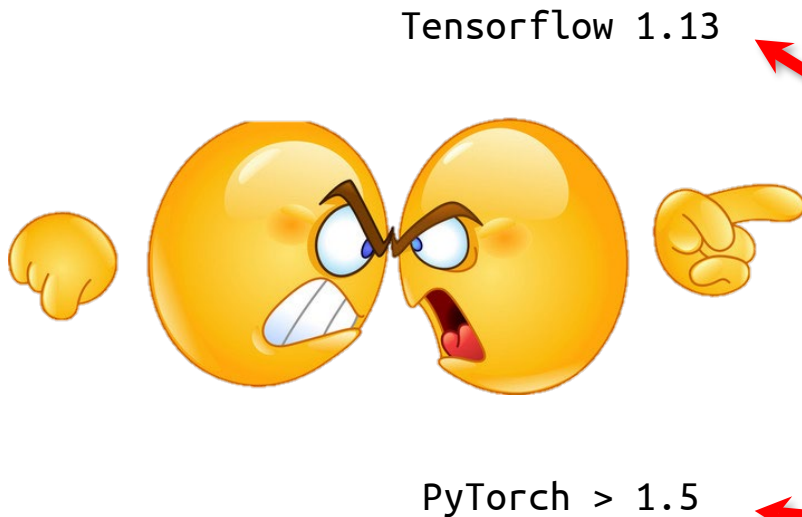
- If you ask Google / ChatGPT...

```
$ sudo yum install ...  
$ sudo apt-get install ...  
$ sudo make install
```



c) Conflicted packages

- What if I need two packages w/ conflicted dependencies?



d) Sharing / Migrating your environment

- Huge effort & large disk quota to install
 - What if my colleagues want to use?
 - What if I want to migrate a different cluster?

Any of those apply to you?

Magic Tools to Install / Manage Software

Part 1:  **CONDA** Virtual Environment

Part 2:  **singularity** Container

1. Why Conda?

- 1) Problems
- 2) Virtual environment & Conda

2. Basic Usage

- 1) Get Conda
- 2) Typical workflow
- 3) Creating a virtual environment
- 4) Installing software packages

3. Advanced Tips

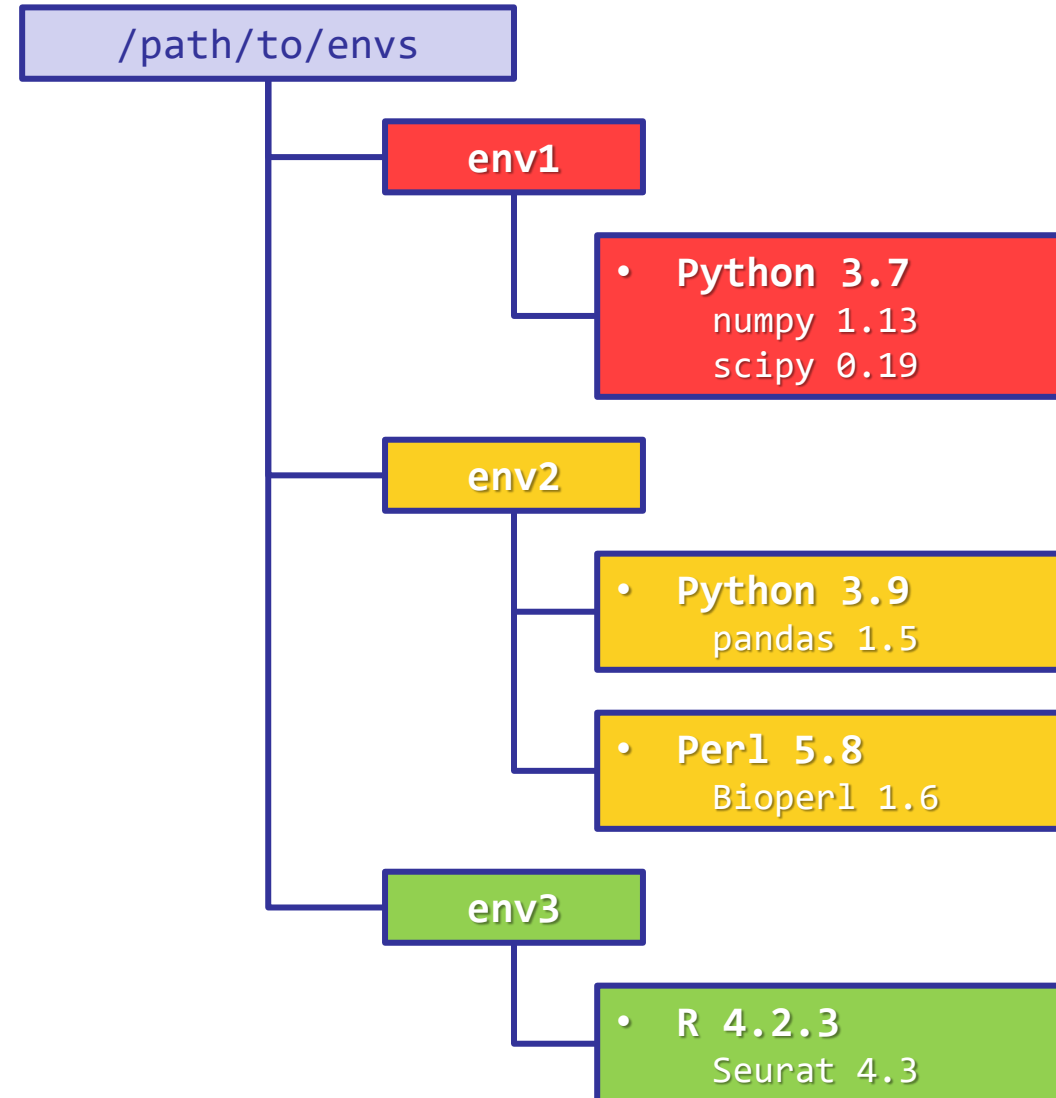


Virtual Environment

2) Virtual environment & Conda

a) What is a **virtual environment**?

- A comprehensive **software framework**, (usually) consists of:
 - A **single directory** contains all files (e.g., executables, dependencies, ...)
 - Proper **configurations** (e.g., environmental variables)



2) Virtual environment & Conda

a) What is a **virtual environment**?

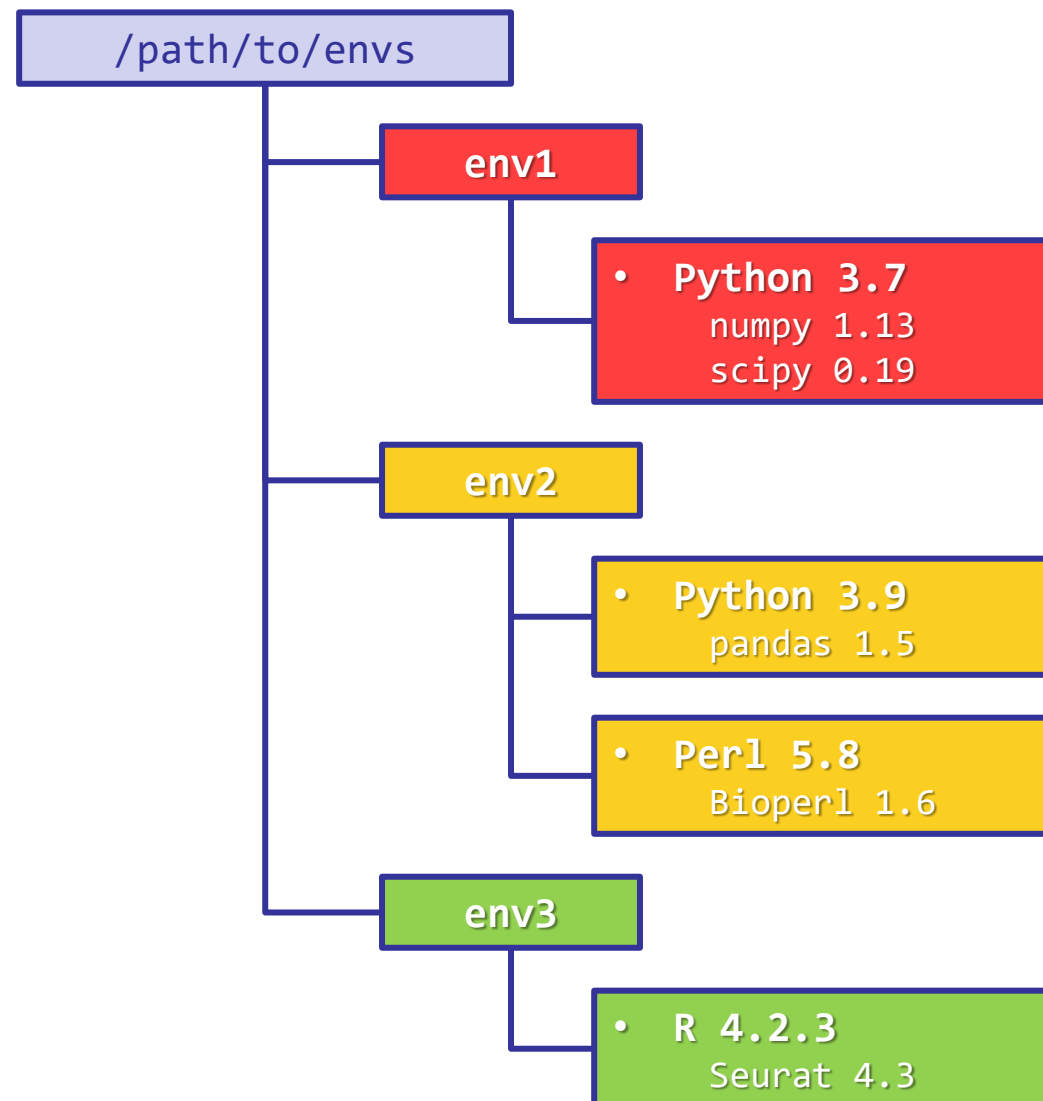
- **Properties**

- **Contained**

All dependencies are installed within the VE

- **Isolated**

Whatever happens in a VE stays in that VE...



b) What is **Conda**



- A **software** tool
- Creates / loads / switches between virtual environments
- Installs / updates / manages packages & dependencies in virtual environments
- Initially for Python → General purposes
- Advantage: Does **NOT** need sudo permission!

c) Relation

CONDA

- is a tool to create / manage
- is not the only tool to create / manage
- usually works with

Virtual
Environment

- In general:

**If a software package you need is managed by Conda,
you (most likely) can install / manage it without needing our help**

1. Why Conda?

- 1) Problems
- 2) Virtual environment & Conda

2. Basic Usage

- 1) Get Conda
- 2) Typical workflow
- 3) Creating a virtual environment
- 4) Installing software packages

3. Advanced Tips

1. Why Conda?

- 1) Problems
- 2) Virtual environment & Conda

2. Basic Usage

- 1) Get Conda
- 2) Typical workflow
- 3) Creating a virtual environment
- 4) Installing software packages

3. Advanced Tips

- ~~• First install a conda in your user environment ... ?~~

1) Get Conda

a) Use conda that comes with system-wide python module

- No installation / disk quota required.
- Sufficient for most user cases.

```
$ module load python
```

Step 1: Can use Conda

```
$ conda init
```

**Step 2: Can use Conda without loading python module
(recommended)**

1) Get Conda

b) Install miniconda

- Latest version: https://repo.anaconda.com/miniconda/Miniconda3-latest-Linux-x86_64.sh

```
$ wget https://repo.anaconda.com/miniconda/Miniconda3-latest-Linux-x86_64.sh
```

Step 1: Download miniconda

```
$ chmod u+x Miniconda3-latest-Linux-x86_64.sh
```

Step 2: Allow execution

```
$ ./Miniconda3-latest-Linux-x86_64.sh
```

Step 3: Run and follow prompts

1. Why Conda?

- 1) Problems
- 2) Virtual environment & Conda

2. Basic Usage

- 1) Get Conda
- 2) Typical workflow
- 3) Creating a virtual environment
- 4) Installing software packages

3. Advanced Tips

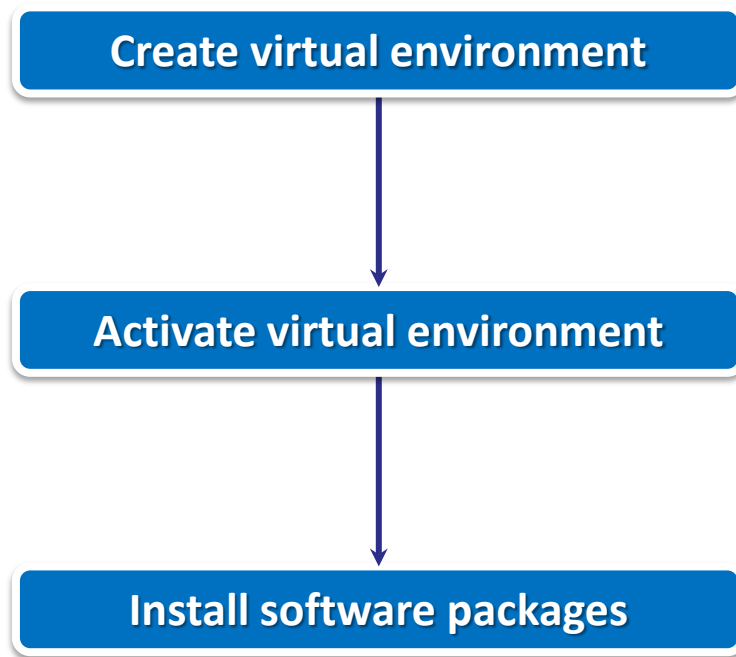
2) Typical workflow

- Key:

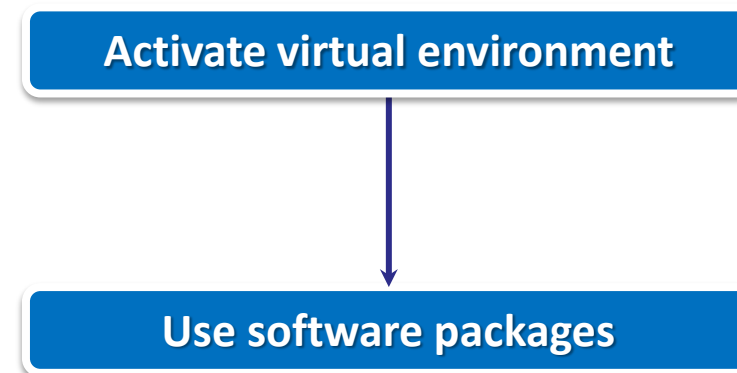
Always use a virtual environment!

2) Typical workflow

To install ...



To use ...



1. Why Conda?

- 1) Problems
- 2) Virtual environment & Conda

2. Basic Usage

- 1) Get Conda
- 2) Typical workflow
- 3) Creating a virtual environment
- 4) Installing software packages

3. Advanced Tips

3) Creating a virtual environment

- Most frequently used commands

Command	
conda create -n ENVIRONMENT	Cre

```
(base) [jasonli3@smic2 ~]$ conda create -n myenv
Collecting package metadata (current_repodata.json): done
Solving environment: done

==> WARNING: A newer version of conda exists. <==
current version: 4.12.0
latest version: 23.1.0

Please update conda by running

$ conda update -n base -c defaults conda

## Package Plan ##

environment location: /home/jasonli3/.conda/envs/myenv

Proceed ([y]/n)?

Preparing transaction: done
Verifying transaction: done
Executing transaction: done
#
# To activate this environment, use
#
#     $ conda activate myenv
#
# To deactivate an active environment, use
#
#     $ conda deactivate
```

[1] <https://docs.conda.io/projects/conda/en/latest/commands.html>



3) Creating a virtual environment

- Most frequently used commands

Command	Description
<code>conda create -n ENVIRONMENT</code>	Create a virtual environment
<code>source activate ENVIRONMENT</code>	Activate a virtual environment

```
(base) [jasonli3@smic2 ~]$ source activate myenv  
(myenv) [jasonli3@smic2 ~]$
```

3) Creating a virtual environment

- Most frequently used commands

Command	Description
<code>conda create -n ENVIRONMENT</code>	Create a virtual environment
<code>source activate ENVIRONMENT</code>	Activate a virtual environment
<code>conda deactivate</code>	Deactivate a virtual environment

```
(myenv) [jasonli3@smic2 ~]$ conda deactivate  
(base) [jasonli3@smic2 ~]$
```

3) Creating a virtual environment

- Most frequently used commands

Command	Description
<code>conda create -n ENVIRONMENT</code>	Create a virtual environment
<code>source activate ENVIRONMENT</code>	Activate a virtual environment
<code>conda deactivate</code>	Deactivate a virtual environment
<code>conda env list</code>	List all virtual environments

```
(base) [jasonli3@smic2 ~]$ conda env list
# conda environments:
#
myenv                /home/jasonli3/.conda/envs/myenv
base                  *  /usr/local/packages/python/3.8.5-anaconda
```

3) Creating a virtual environment

- Most frequently used commands

Command	Description
<code>conda create -n ENVIRONMENT</code>	Create a virtual environment
<code>source activate ENVIRONMENT</code>	Activate a virtual environment
<code>conda deactivate</code>	Deactivate a virtual environment
<code>conda env list</code>	List all virtual environments
<code>conda env remove -n ENVIRONMENT</code>	Remove a virtual environment and all installed packages

CAUTION! NO CONFIRMATION! IRREVOCABLE!

3) Creating a virtual environment

- Most frequently used commands

Command	Description
conda create -n ENVIRONMENT	Create a virtual environment
source activate ENVIRONMENT	Activate a virtual environment
conda deactivate	Deactivate a virtual environment
conda env list	List all virtual environments
conda env remove -n ENVIRONMENT	Remove a virtual environment and all installed packages

1. Why Conda?

- 1) Problems
- 2) Virtual environment & Conda

2. Basic Usage

- 1) Get Conda
- 2) Creating a virtual environment
- 3) Typical workflow
- 4) Installing software packages

3. Advanced Tips

4) Installing software packages

- Before installation...

Make sure a virtual environment is activated!

4) Installing software packages

a) Most frequently used commands

Command	Description
<code>conda install NAME</code>	Install a software package

4) Installing software packages

a) Most frequently used commands

Command	Description
<code>conda install NAME</code>	Install a software package
<code>conda install NAME=VERSION</code>	Install a specific version

4) Installing software packages

a) Most frequently used commands

Command	Description
<code>conda install NAME</code>	Install a software package
<code>conda install NAME=VERSION</code>	Install a specific version
<code>conda install NAME -c CHANNEL</code>	Install from a specific channel (e.g., conda-forge, bioconda, nvidia, ...)

4) Installing software packages

a) Most frequently used commands

Command	Description
<code>conda install NAME</code>	Install a software package
<code>conda install NAME=VERSION</code>	Install a specific version
<code>conda install NAME -c CHANNEL</code>	Install from a specific channel (e.g., conda-forge, bioconda, nvidia, ...)
<code>conda install NAME1 NAME2 ...</code>	Install multiple packages at once (let conda work out dependencies)

4) Installing software packages

a) Most frequently used commands

Command	Description
<code>conda install NAME</code>	Install a software package
<code>conda install NAME=VERSION</code>	Install a specific version
<code>conda install NAME -c CHANNEL</code>	Install from a specific channel (e.g., conda-forge, bioconda, nvidia, ...)
<code>conda install NAME1 NAME2 ...</code>	Install multiple packages at once (let conda work out dependencies)
<code>conda list</code>	List all installed software package

4) Installing software packages

b) Other useful commands

Command	Description
<code>conda search NAME</code>	Search available package versions

4) Installing software packages

b) Other useful commands

Command	Description
conda search NAME	Search available package versions
conda search NAME -c CHANNEL	Search available package versions in a specific channel

4) Installing software packages

b) Other useful commands

Command	Description
<code>conda search NAME</code>	Search available package versions
<code>conda search NAME -c CHANNEL</code>	Search available package versions in a specific channel
<code>conda search NAME --info</code>	Search available package versions with details

4) Installing software packages

b) Other useful commands

Command	Description
conda search NAME	Search available package versions
conda search NAME -c CHANNEL	Search available package versions in a specific channel
conda search NAME --info	Search available package versions with details
conda update/upgrade NAME	Update a package to the latest available version

4) Installing software packages

b) Other useful commands

Command	Description
<code>conda search NAME</code>	Search available package versions
<code>conda search NAME -c CHANNEL</code>	Search available package versions in a specific channel
<code>conda search NAME --info</code>	Search available package versions with details
<code>conda update/upgrade NAME</code>	Update a package to the latest available version
<code>conda uninstall/remove NAME</code>	Remove a package

[1] <https://docs.conda.io/projects/conda/en/latest/commands.html>

4) Installing software packages

c) Bonus: Hot packages!

i. PyTorch (w/ GPU support)

```
$ # Start an interactive job (using qsub or srun)
```

Step 1: Start an interactive job (otherwise will timeout)

```
$ conda create -n torch
```

```
$ source activate torch
```

```
$ conda install pytorch==2.0.1 torchvision==0.15.2 torchaudio==2.0.2 pytorch-cuda=11.8 -c pytorch -c nvidia
```

Step 2: Create a VE and install PyTorch & dependencies

4) Installing software packages

c) Bonus: Hot packages!

ii. Tensorflow (w/ GPU support)

```
$ # Start an interactive job on *GPU* nodes (using qsub or srun)
```

Step 1: Start an interactive job on *GPU* nodes (otherwise will fail)

```
$ conda create -n tf
```

```
$ source activate tf
```

```
$ conda install -c conda-forge -c nvidia tensorflow-gpu=2.12 cudatoolkit=11.8 cuda-nvcc=11.8
```

Step 2: Create a VE and install tensorflow & dependencies

```
$ mkdir -p $CONDA_PREFIX/etc/conda/activate.d
```

```
$ echo 'export XLA_FLAGS=--xla_gpu_cuda_data_dir=$CONDA_PREFIX' >> $CONDA_PREFIX/etc/conda/activate.d/env_vars.sh
```

Step 3: Run these commands to set up environment

- Your workflow should mostly look like...

To install ...

```
$ conda create ...  
  
$ source activate ...  
  
$ conda install ...
```

To use ...

```
$ source activate ...  
  
$ # Do whatever you need  
   to do with the packages
```


- Create a virtual environment
- Search for **SciPy** version and install the second-latest version (as well as dependencies)
- After you are done, type in chat the installed **SciPy** and **Python** version

1. Why Conda?

- 1) Problems
- 2) Virtual environment & Conda

2. Basic Usage

- 1) Get Conda
- 2) Typical workflow
- 3) Creating a virtual environment
- 4) Installing software packages

3. Advanced Tips

A little more than the basics...

1) Where to get software?

- You can get software from a lot of places using Conda
 - **Not that you should!**
- Concerns?
 - **Reliability** (some third-party or untested images may not work)
 - **Security risk** (some untrustworthy publishers may pack something you don't know about)
- Solution
 - Always get from a source that **you can trust**

1) Where to get software?

- **Tier 1: Developer release (official release)**

- On software's official website, look for “**Conda**”.
- E.g., [PyTorch](#), [Spyder](#), [CudaToolKit](#)

- **Tier 2: Trustworthy channels**

Name	Notes
default / anaconda	• Default channel, officially managed by Conda
conda-forge	• Community supported, but rule-enforced and generally trustworthy
bioconda	• Community supported for bioinformatics
nvidia / cuda	• Nvidia official channel
pytorch	• Pytorch official channel
intel	• Intel official channel
...	

2) Change Conda path

- Default Conda path

	System-wide Python modules	Customized Conda
Environments	<code>/home/\$USER/.conda/envs/</code>	<code>/path/to/conda/envs/</code>
Cache	<code>/work/\$USER/.conda/pkgs/</code>	<code>/path/to/conda/pkgs/</code>

```
[jasonli3@smic1 ~]$ balance
User filesystem quotas for jasonli3 (uid 15827):
  Filesystem      MB used   MB quota
  /home           950      10000
  /work /project  329639    0        6
Storage allocation  MB used   MB quota
```

2) Change Conda path

a) Method 1: Command lines

```
$ conda config --add envs_dirs /path/to/envs
```

Add path to environments

```
$ conda config --add pkgs_dirs /path/to/pkgs
```

Add path to downloaded packages

2) Change Conda path

b) Method 2: Configuration file

- Use any text editor to open: `~/.condarc`

```
$ vi ~/.condarc
```

```
envs_dirs:  
  - /project/jasonli3/.conda/envs  
pkgs_dirs:  
  - /project/jasonli3/.conda/pkgs
```

```
~
```


2) Change Conda path

c) Places to store your virtual environments:

Location	Pros	Cons
/home	<ul style="list-style-type: none">• All users have access• No expiration date• Backed up	<ul style="list-style-type: none">• Limited quota (10 GB)
/project	<ul style="list-style-type: none">• Larger quota (x 100 GB)• Valid for one year & renewable• Can be shared among group	<ul style="list-style-type: none">• Not all users have access (PI must apply for /project drive)
/work	<ul style="list-style-type: none">• All users have access• No quota limit	<ul style="list-style-type: none">• Files are subject to purge!

2) Change Conda path

c) Places to store your virtual environments:

Location	Pros	Cons
/home	<ul style="list-style-type: none">• All users have access• No expiration date• Backed up	<ul style="list-style-type: none">• Limited quota (10 GB)
/project	<ul style="list-style-type: none">• Larger quota (x 100 GB)• Valid for one year & renewable• Can be shared among group	<ul style="list-style-type: none">• Not all users have access (PI must apply for /project drive)
/work	<ul style="list-style-type: none">• All users have access• No quota limit	<ul style="list-style-type: none">• Files are subject to purge!

3) Share virtual environment

- **Scenario:**

- I made a huge effort to install an extensive collection of software packages for our group's research needs. I don't want to do it all over again for everyone in our group. Is it possible to just share the virtual environment with them?

3) Share virtual environment

- **Solution:**

- a) **PI:**

- Apply for a storage allocation (a.k.a. /project, if hasn't)
 - Add User 1 (sharing) and User 2 (shared) to /project

- b) **User 1 (sharing):**

- Set up **envs_dirs** to create a virtual environment in **a /project location**
 - Install software in the virtual environment

- c) **User 2 (shared):**

- Set up **envs_dirs** to create a virtual environment in **the same /project location**

4) Migrate / clone virtual environment

- **Scenario:**

- I have been using LSU HPC cluster. But now I want to switch to LONI and run the exactly same software. How can I do that?
- I am leaving my current position. But I may continue doing similar research. How can I replicate my environment to a different HPC system in a different institute?

4) Migrate / clone virtual environment

- **Solution**

To ...

Run command

4) Migrate / clone virtual environment

- Solution

To ...

Export virtual environment recipe to file

```
name: spyder
channels:
  - defaults
dependencies:
  - _libgcc_mutex=0.1=main
  - _openmp_mutex=5.1=1_gnu
  - arrow=1.2.3=py310h06a4308_1
  - astroid=2.14.2=py310h06a4308_0
  - attrs=22.1.0=py310h06a4308_0
  - babel=2.11.0=py310h06a4308_0
  - beautifulsoup4=4.11.1=py310h06a4308_0
  - black=22.6.0=py310h06a4308_0
  - blas=1.0=mkl
  - bottleneck=1.3.5=py310ha9d4c09_0
  - brotli=1.0.9=h5eee18b_7
```

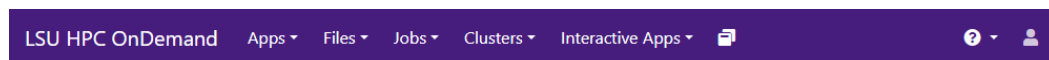
4) Migrate / clone virtual environment

- **Solution**

To ...	Run command
Export virtual environment recipe to file	<code>conda env export > myenv.yml</code>
Create a virtual environment from file	<code>conda env create -f myenv.yml</code>

5) Use virtual environment on Open OnDemand

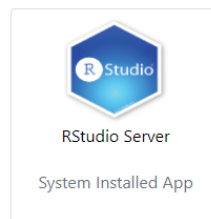
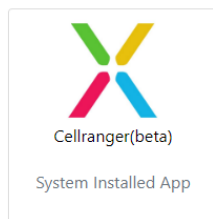
LSU HPC (SMIC / SuperMike 3)



OnDemand provides an integrated, single access point for all of your HPC resources.

Pinned Apps A featured subset of all available apps

Interactive Apps

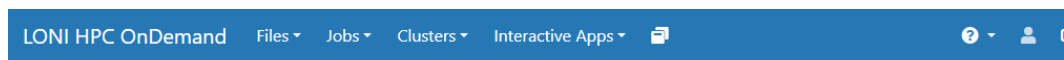


Message of the Day

Welcome to the LSU HPC OnDemand portal!

With the OnDemand web portal, you can:

LONI (QB2)



OnDemand provides an integrated, single access point for all of your HPC resources.

Message of the Day

Welcome to the LONI HPC OnDemand portal!

With the OnDemand web portal, you can:

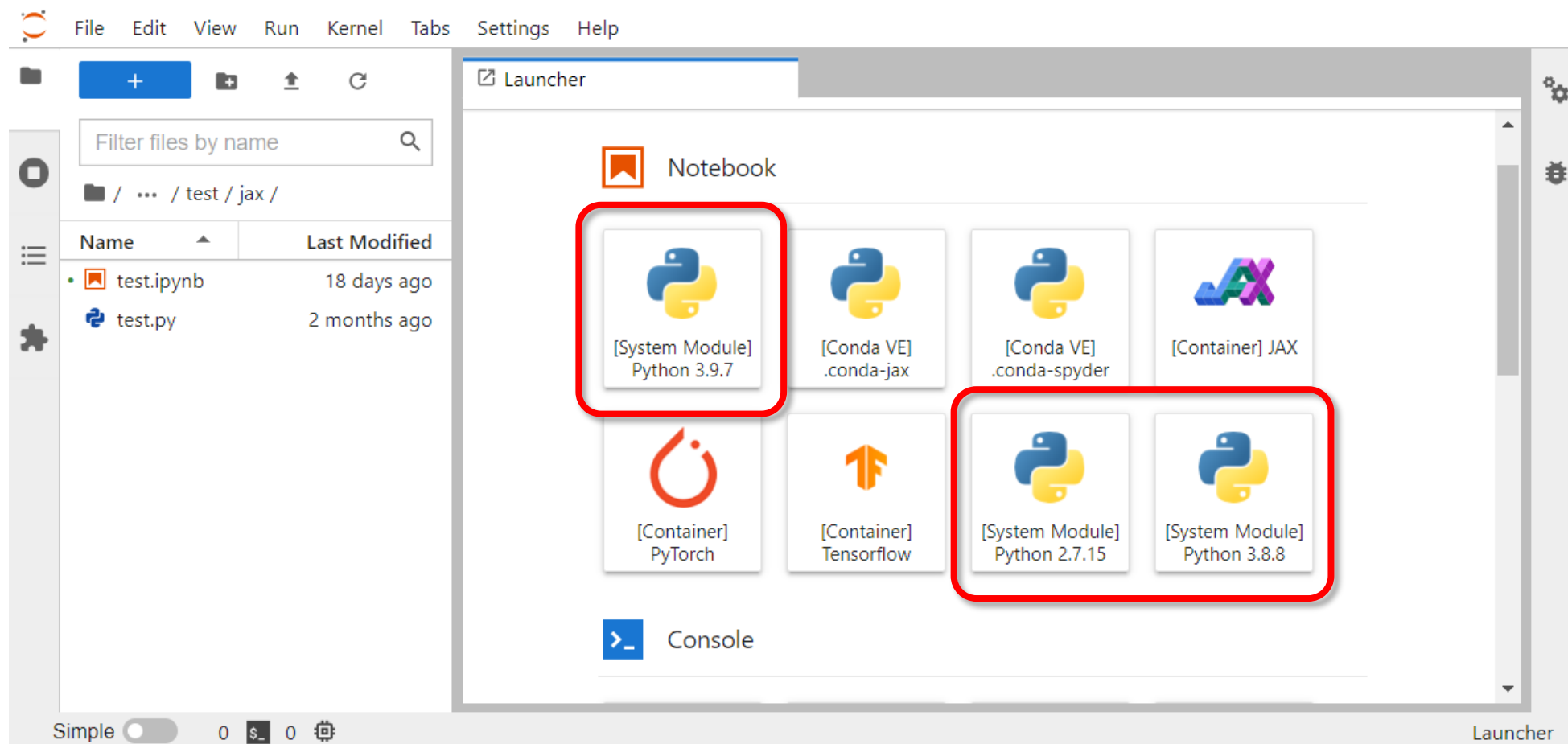
- - Manage, download and upload files to the HPC systems (click links in the "Files" on the top of this page)
- - Check allocation balances
- - Check disk usage and quotas
- - Check job status
- - Submit jobs using templates
- - Access HPC systems via a terminal
- - Run interactive apps such as Jupyter Notebook/Lab and Rstudio (click links in the "Interactive Apps" on the top of this page)

Getting started

[1] <https://youtu.be/JLXN0AZgaqc>

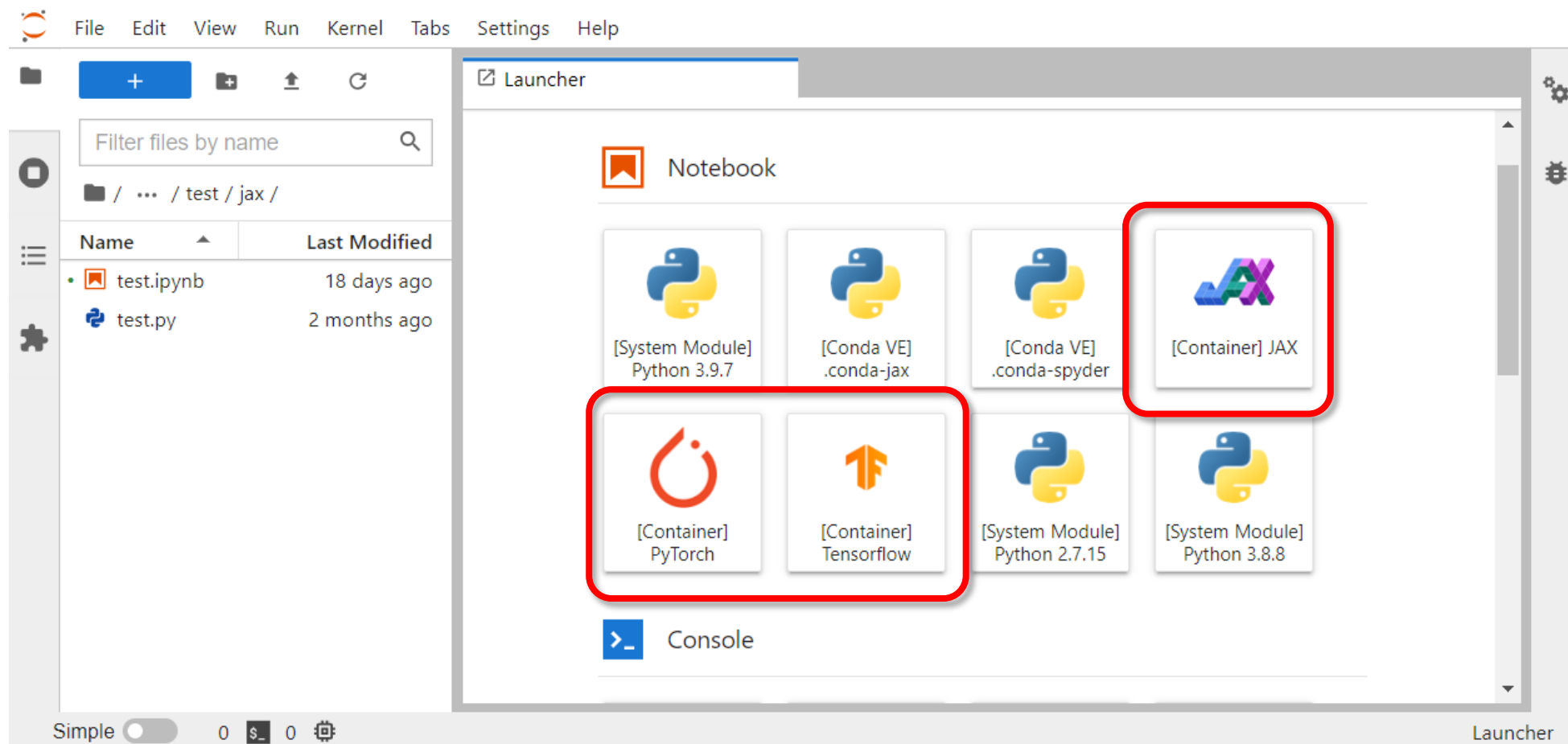


5) Use virtual environment on Open OnDemand



[1] <https://youtu.be/JLXN0AZgaqc>

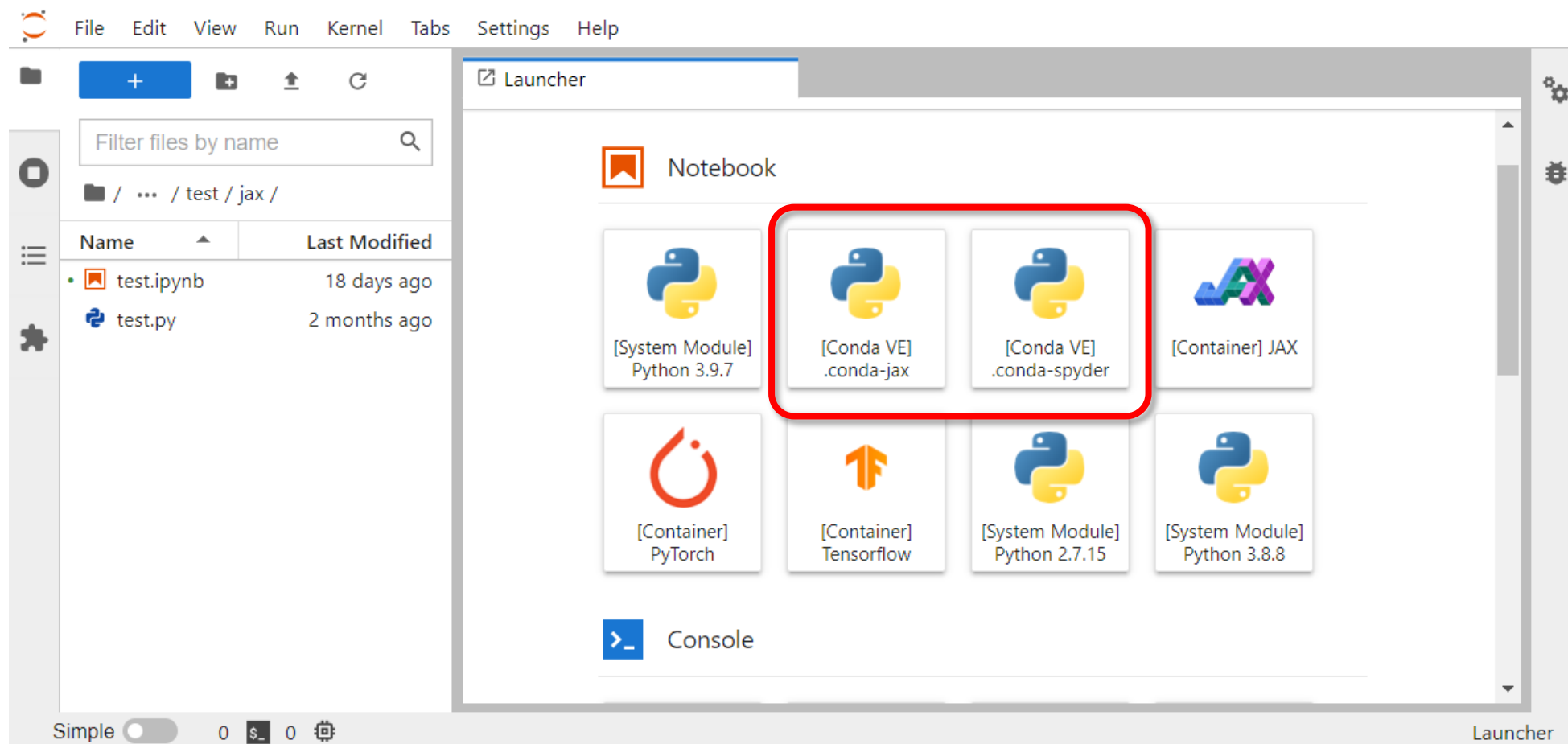
5) Use virtual environment on Open OnDemand



[1] <https://youtu.be/JLXN0AZgaqc>



5) Use virtual environment on Open OnDemand



[1] <https://youtu.be/JLXN0AZgaqc>



5) Use virtual environment on Open OnDemand

- **How to:**

Step 1: ssh to the cluster you want to use

LSU HPC	LONI
SMIC / Super Mike 3	QB2

Step 2: Activate the virtual environment you want to use in Jupyter

```
$ source activate ENVIRONMENT
```

Step 3: Install ipykernel

```
$ conda install ipykernel          # Either use this
$ pip install ipykernel            # Or use this
```

Step 4: Start a Jupyter session in Open OnDemand, and choose the environment in **kernel**

- **Scenario**
 - I need software packages other than Python (R / Perl / Lua / ...)
 - I need a different version than the system modules
 - I am using the system's R module, but having trouble installing some packages (e.g., rgdal)

```
> install.packages("rgdal")
Warning in install.packages("rgdal") :
  'lib = "/home/packages/r/4.1.2/5k5jengl/rlib/R/library"' is not writable
would you like to use a personal library instead? (yes/no) yes
no
configure: error: gdal-config not found or not executable.
ERROR: configuration failed for package 'rgdal'
* removing '/home/jasonli3/R/x86_64-pc-linux-gnu-library/4.1/rgdal'

The downloaded source packages are in
  '/tmp/Rtmpd2csho/downloaded_packages'
Warning message:
In install.packages("rgdal") :
  installation of package 'rgdal' had non-zero exit status
```

- Solutions

Many non-python packages are managed by Conda too!

To install ...		Run command ...
Languages	R	<code>conda install R</code>
	Perl	<code>conda install perl</code>
	Julia	<code>conda install julia -c conda-forge</code>
Dependencies	hdf5	<code>conda install hdf5</code>
	netcdf	<code>conda install libnetcdf -c conda-forge</code>
	FFTW	<code>conda install fftw</code>
		...

6) More than Python

- One more cool thing...
 - You can use language specific package management tools

Language	Tool
Python	pip
R	install.packages
Perl	cpan
Julia	Pkg

- Packages will be **isolated** in the virtual environment

6) More than Python

- E.g., Use Conda to solve your R issue

- Use system's R module:

```
$ module load r  
$ R  
> install.packages("rgdal")
```

→ Will fail!

- Use Conda:

```
$ conda create -n rgdal  
$ source activate rgdal  
$ conda install r-rgdal
```

→ Will succeed!

7) Troubleshooting

a) Conflict with system module

```
(spyder) [jasonli3@smic2 ~]$ module li
Currently Loaded Modulefiles:
1) python/3.8.5-anaconda
(spyder) [jasonli3@smic2 ~]$ python
Python 3.10.9 (main, Mar  8 2023, 10:47:38) [GCC 11.2.0] on linux
Type "help", "copyright", "credits" or "license" for more information.
>>> import numpy as np
/usr/local/packages/python/3.8.5-anaconda/lib/python3.8/site-packages/numpy/__init__.p
fore Intel(R) MKL initialization ensuring correct out-of-the box operation under c
ess is not assured. Please install mkl-se see http://github.com/IntelPyt
from . import _distributor_init
Traceback (most recent call last):
  File "/usr/local/packages/python/3.8.5-anaconda/lib/python3.8/site-packages/numpy/co
from . import multiarray
File "/usr/local/packages/python/3.8.5-anaconda/lib/python3.8/site-packages/numpy/co
```

Fail!

a) Conflict with system module

- Rule of thumb:

Do **NOT** load system module if you are using your own installation!

b) What if I made a mess?

- I mixed conda / pip back and forth, and broke the environment...
- It may be easier to create a new virtual environment and start fresh...

Conclusion

To install ...

```
$ conda create ...  
  
$ source activate ...  
  
$ conda install ...
```

To use ...

```
$ source activate ...  
  
$ # Do whatever you need  
   to do with the packages
```

Next week in our miniseries

Magic Tools to Install / Manage Software

Part 2: ingularity Container

Jason Li

HPC User Services

LSU HPC / LONI

sys-help@loni.org

Louisiana State University, Baton Rouge

Nov 1, 2023

Conda vs Singularity

	Conda / Virtual Environments	Singularity / Containers
Availability	All users	All users, but may need additional things
Functionality	Good (usually)	Better (more likely to work)
Self-contained	Yes	Yes
Isolation	Yes (but still accessible from outside)	Perfect
Editability	Yes	No (image file itself) / Yes (a detour with sandbox)
Disk usage	Large	Smaller
Portability	Possible (with .yaml file)	Great (copy-paste one file)
Security	Fair	Good
Ease of use	Good	May require more effort

	Conda / Virtual Environments	Singularity / Containers
Good for	<ul style="list-style-type: none">• Less hassle (usually?) to create and install software from scratch• If you need to frequently make modifications• If you need to access files from outside of the environment (e.g., compiling a code that uses some files in the virtual environment as dependencies)	<ul style="list-style-type: none">• Less hassle if the developer releases a working container• If you don't or don't want to make changes after it is created• Portability• Reduce disk usage• Your system admins yelled at you about security risk

- **Contact user services**

- Email Help Ticket: sys-help@loni.org
- Telephone Help Desk: +1 (225) 578-0900