Running minAone optimization code

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Usage

```
equations.txt → python_scripts directory
```

1. Setup python dependencies (sympy)

```
conda create --name anaconda_env anaconda conda activate anaconda-env
```

- 2. Collect all your data files based on specs.txt
- 3. python original_minaone.py → Makefile
- 4. make → neutrinos_cpp executable
- 5. ./neutrinos.cpp OR submit.sbatch → .dat files (a LOT for 8-beam; ~256MB), ipopt.out

Troubleshooting a Segfault:

- Are all the data files correctly labeled in specs.txt? Are they all in the directory?
- Are all data files the right/same length? (Use wc)
- Special characters can be viewed with vim -b FILENAME
 - trailing whitespace " " matters
 - linebreak-likes "^M" might matter
- If minAone.py runs, equations.txt is OK
- If neutrinos cpp runs, specs.txt is OK

Troubleshooting other issues/dependencies

```
==> WARNING: A newer version of conda exists. <==
current version: 4.6.12
latest version: 4.13.0

Please update conda by running
$ conda update -n base -c defaults conda
```

Jupyter notebook tunnel:

```
On remote, run jupyter notebook --no-browser --port=NNNN

Prerequisite: module add anaconda

On local, run ssh -t -t USERNAME @hpc-logon.nyit.edu -L NNNN:localhost:NNNN ssh node002 -L

NNNN:localhost:NNNN

You ran ssh -t -t swong25@10.10.32.70 -L 6969:localhost:6969 ssh node002 -L

8892:localhost:6969
```

GitHub on the cluster

prequisites:

- SSH key (haven't figured out how that works yet)
- OR Personal Access Token

```
git clone <URL>
git add FILES; git commit -m "MESSAGE"
git push origin master
```