# a guide to environment preparation

[ python environment for simple numerical quantum mechanics ]

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## 2 Steps

```
Install Miniconda (python environment)
Install python modules
Test Jupyter Notebook
```

#### **Download Miniconda**

#### From here: URL

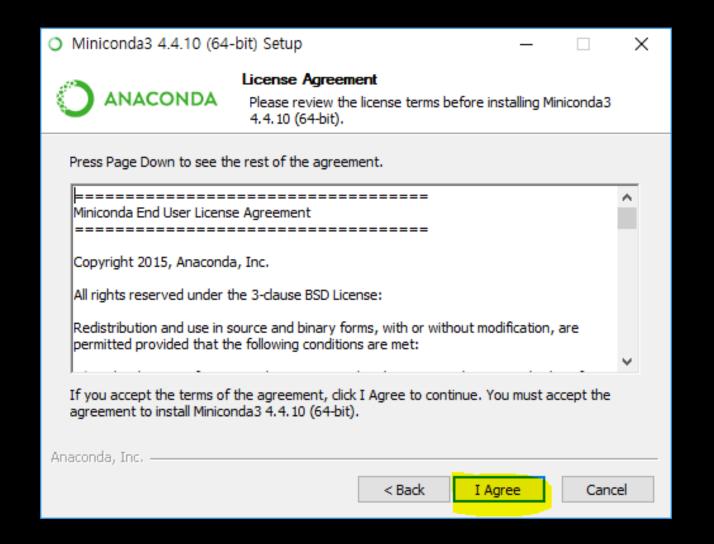
#### Windows Mac OS X Linux 64-bit (bash installer) Python 3.6 64-bit (exe installer) 64-bit (bash installer) 32-bit (exe installer) 32-bit (bash installer) Python 2.7 64-bit (exe installer) 64-bit (bash installer) 64-bit (bash installer) 32-bit (exe installer) 32-bit (bash installer)

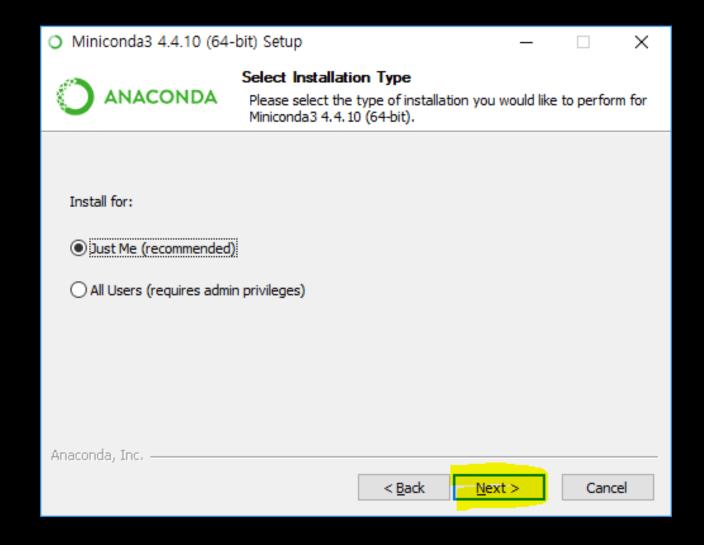


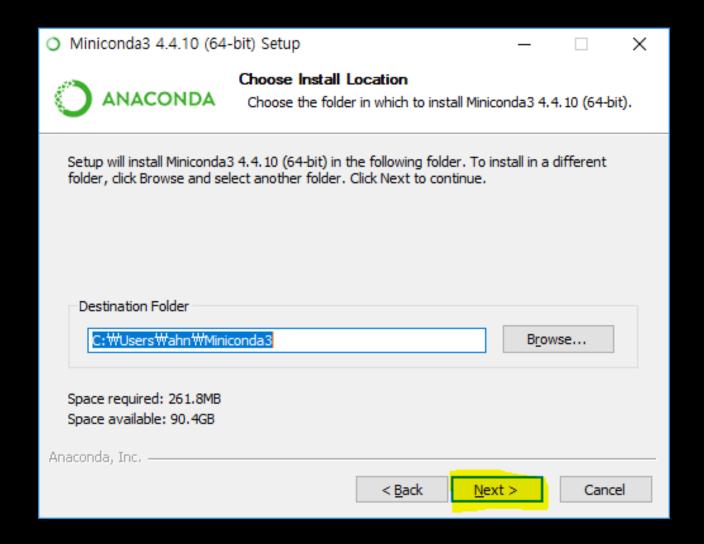
This is for Windows.

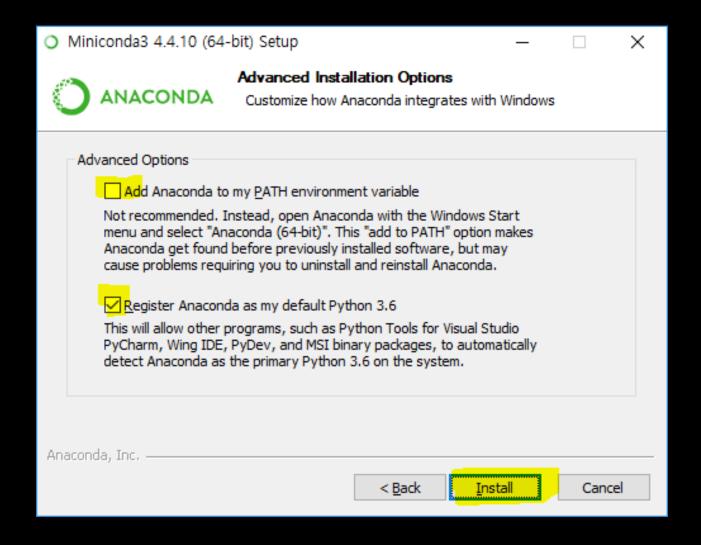
For other operating system,

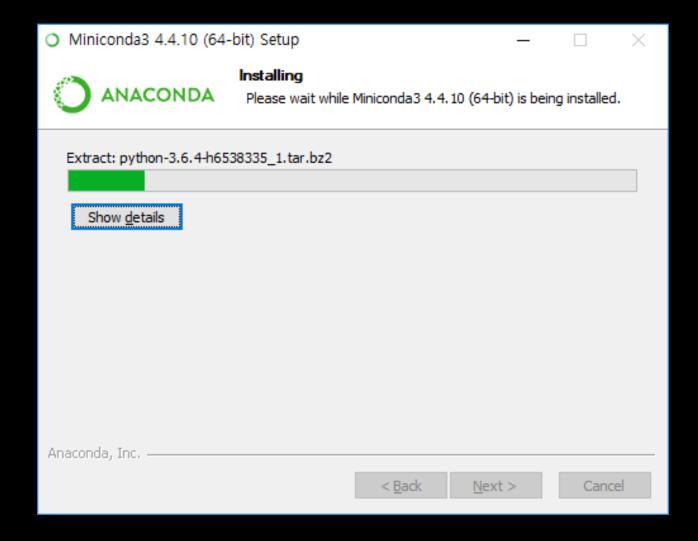
refer to HERE

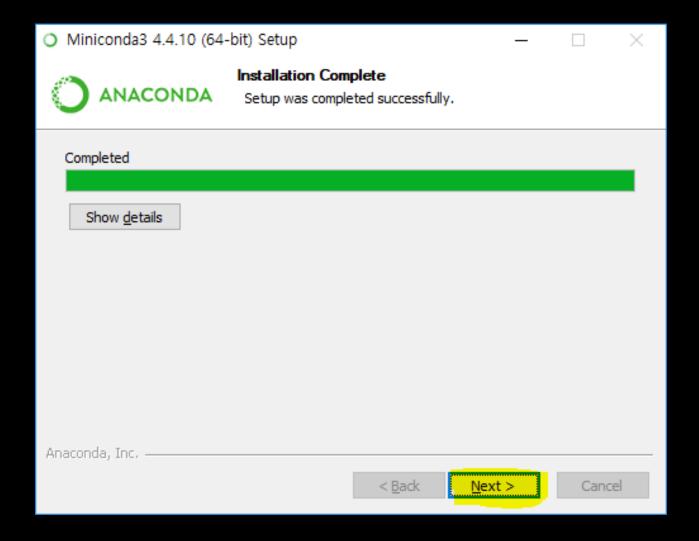








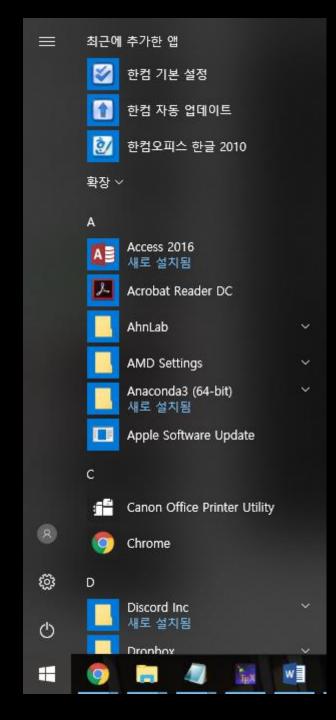






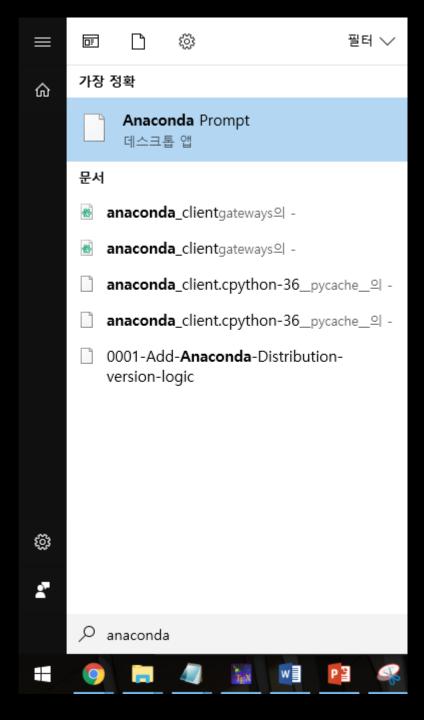
#### On Windows:

- 1. Press Windows Key
- → You will see a 'Start Menu'



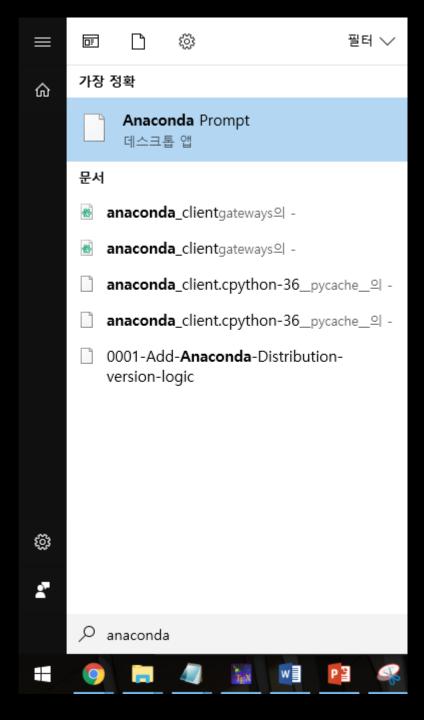
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- 3. Hit 'Enter' Key to open it



On Mac OS: Assuming your miniconda installation directory is "\$HOME/miniconda"

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- → PATH should include installation binary directory
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For more info: **LINK** 

# Install python modules

#### **Enter:**

pip install numpy matplotlib jupyter vis

**Enter:** 

Anaconda Prompt

pip install numpy matplotlib jupyter vis

(base) C:₩Users₩ahn>pip install numpy matplotlib jupyter vis

Same for Mac OS

**Enter:** 

Anaconda Prompt

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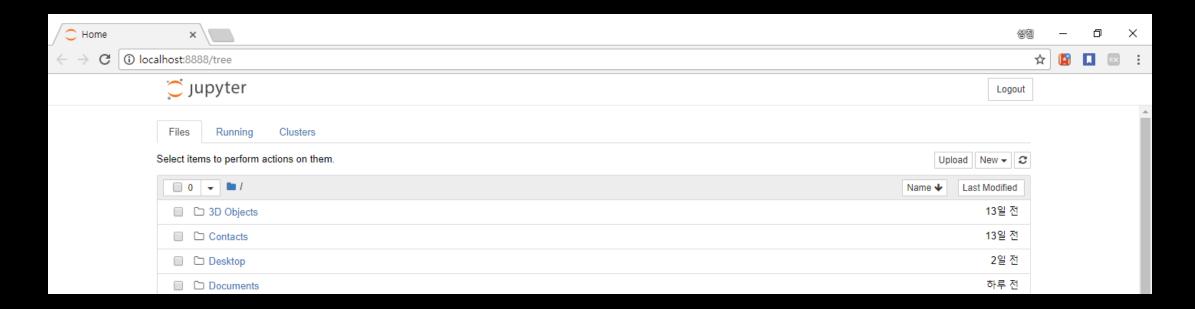
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# Test Jupyter Notebook

#### **Enter:**

jupyter notebook

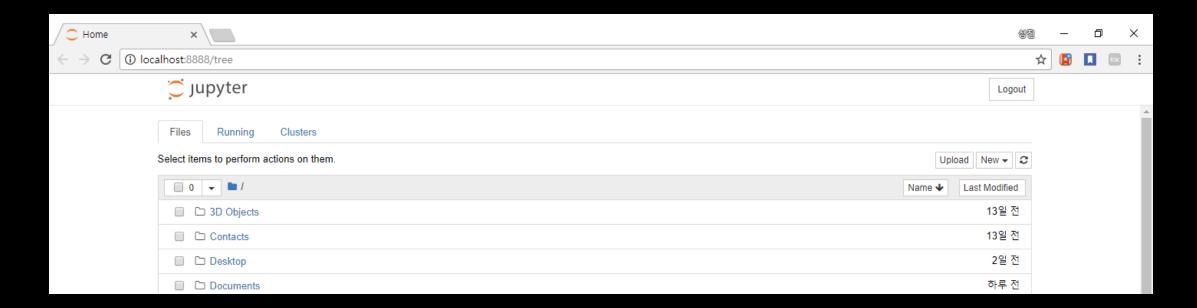


**Enter:** 

You can get some basics from **HERE** 



jupyter notebook



fin.