

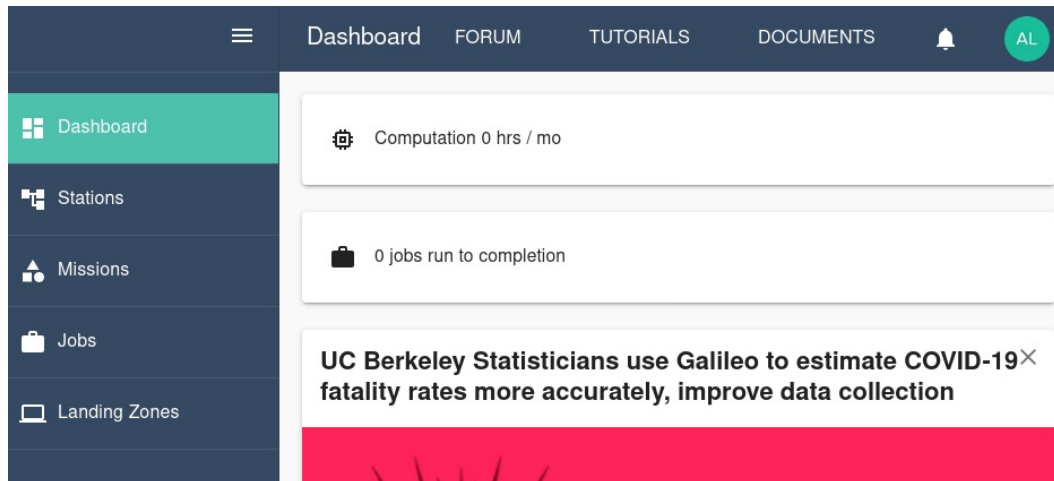
# Tutorial: Running Quantum Espresso in Galileo

## Getting started with Quantum Espresso in Galileo

To get started with Galileo, [log into your account](#) using Firefox or Chrome.

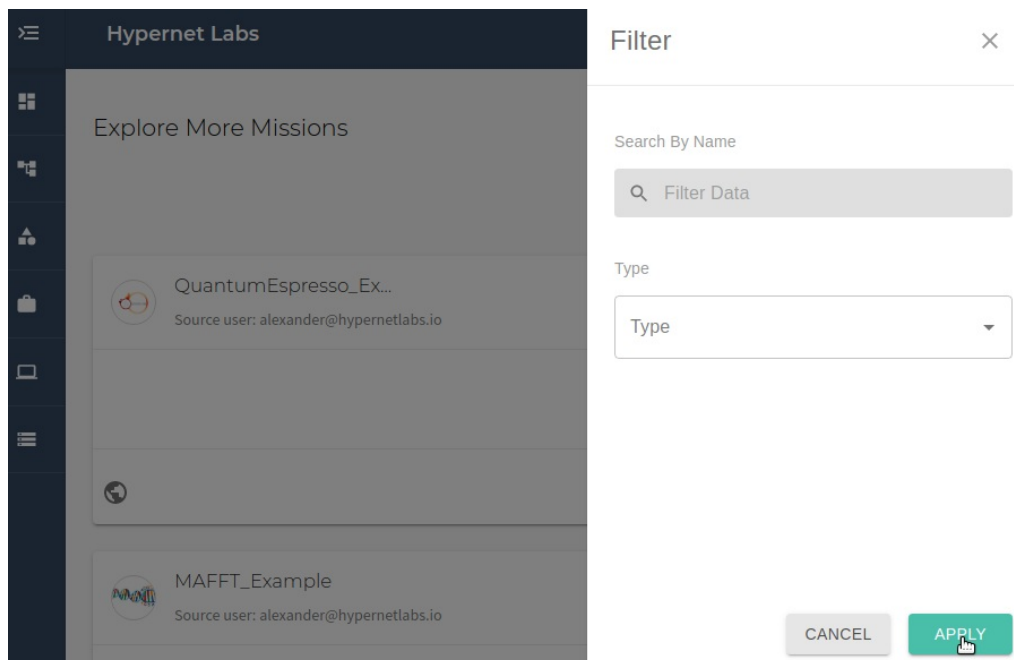
## Understanding the user interface and cloning a Mission

When you log into Galileo, the first thing you'll see is your Dashboard:



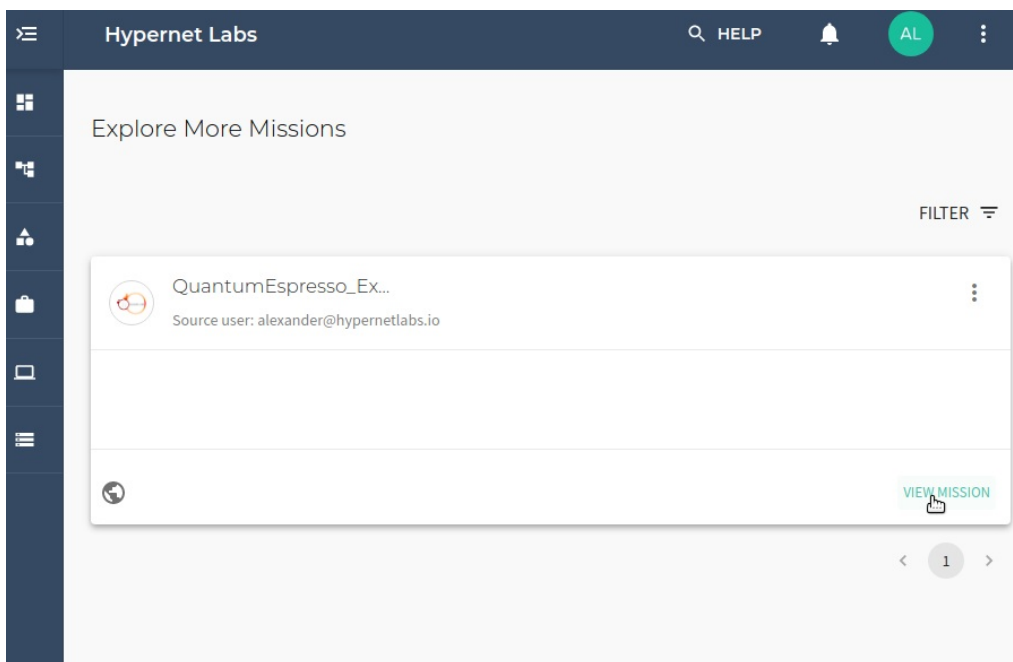
View of the Galileo Dashboard

To run the Quantum Espresso example, start by navigating to the Missions tab using the side menu. Clone the Quantum Espresso example Mission from the Explore Missions tab. Use the filter to search for the mission by name and click “Apply”.



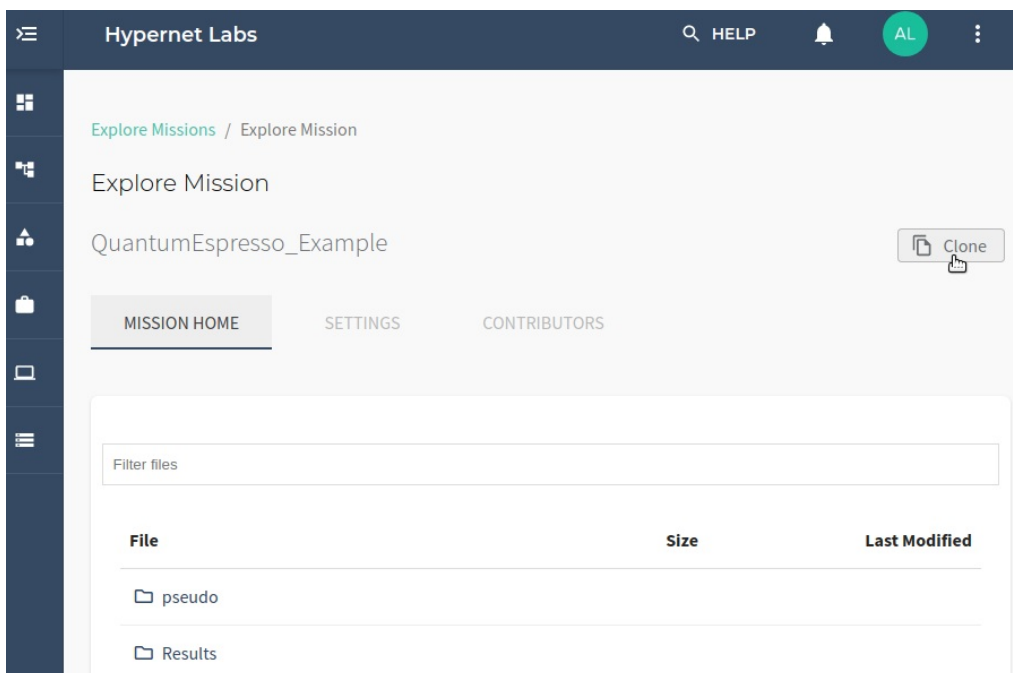
Find the public example mission by name

Once you have found the correct Mission, click “View Mission”.



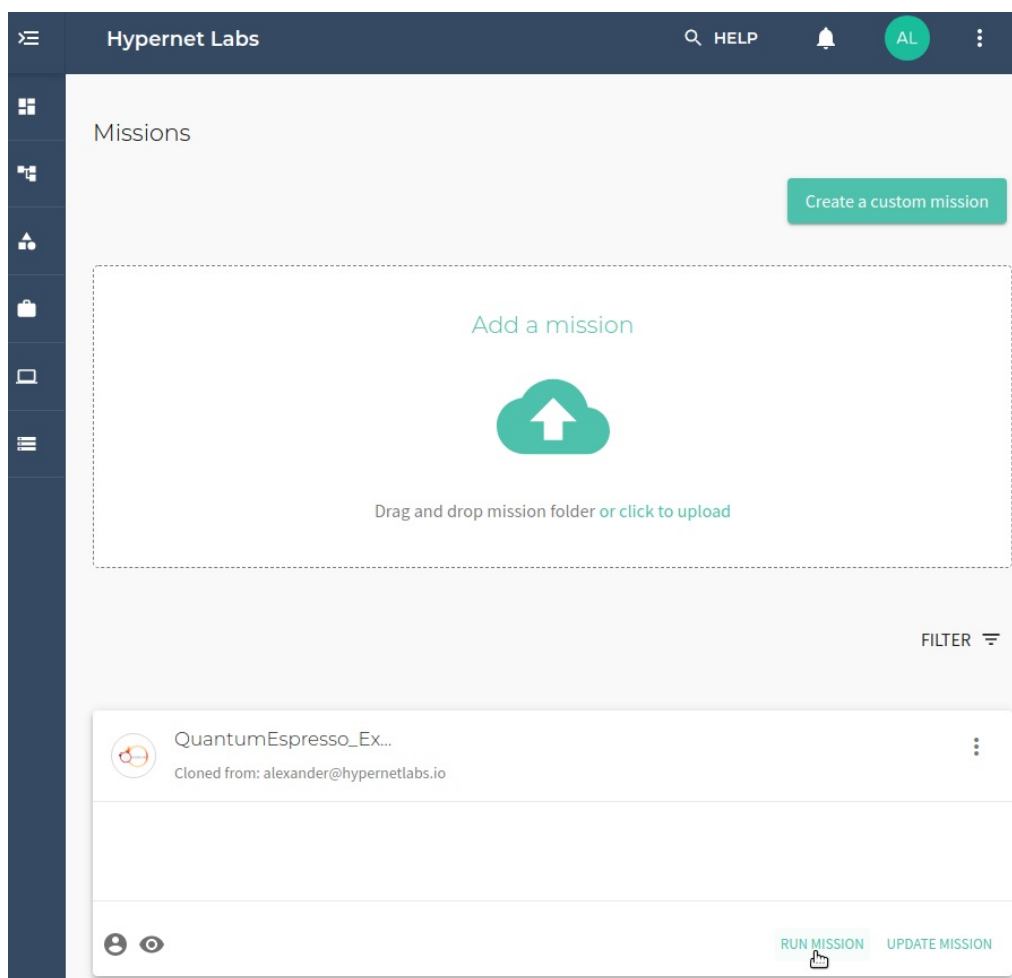
Click View Mission

To clone the public Mission to your account, click the “Clone” button in the upper right corner of the interface. Choose between creating a public or private clone and also choose which Cargo Bay to use.



Clone the mission

You will now see a cloned copy of the Mission in your Missions.



The cloned copy

### Let's take a look at our files


First, we have a series of files to run our computations: `pw.methane.in` is an input file that conducts the SCF ground-state calculation, `turbo_davidson.methane.in` runs a Davidson calculation of our eigenvalues, and `turbo_spectrum.methane.in` completes a post-processing analysis of our spectrum.


We also have the files `plot_spectrum_nohyb.gp` and `plot_spectrum_hyb.gp`, which plot our spectrum using gnuplot with and without B3LYP pseudo-potential, respectively.

Finally, we have a shell script called `runqe.sh` that contains our QE commands.

### Running a job and collecting results

Now we are ready to run a job using the Mission. Click the **Run** button in the upper right corner of the Mission tab. You will see a "Mission run successfully!" message. At the bottom of the Mission tab, you can track the progress of the job.

Jobs 

QuantumEspresso\_Example [Show More](#) 


**Queued**


UPLOADED	: Thu Mar 11 2021 15:47:54	SENT TO	: Landing Zone Pending
SENT BY	: alexander_berry@brown.edu	TIME TAKEN	: 00:00:00

Rows per page: 20 ▾ 1-1 of 1 |< < > >|

Track job progress

Once the computation is completed, the job will shut down and collect the results. Once the job progress reads “Completed”, you can download the results by opening the three-dot menu and clicking **Download**.



Jobs 

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**Completed**

UPLOADED	: Thu Mar 11 2021 15:47:54	SENT TO	: andromeda-21
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 Download  
 Archive

Download results

Let's take a look at the output log:

```

Program PWSCF v.6.5 starts on 11Mar2021 at 20:48:34

This program is part of the open-source Quantum ESPRESSO suite
for quantum simulation of materials; please cite
"P. Giannozzi et al., J. Phys.:Condens. Matter 21 395502 (2009);
"P. Giannozzi et al., J. Phys.:Condens. Matter 29 465901 (2017);
URL http://www.quantum-espresso.org",
in publications or presentations arising from this work. More details at
http://www.quantum-espresso.org/quote

Parallel version (MPI), running on      1 processors

MPI processes distributed on      1 nodes
Waiting for input...
Reading input from standard input

Current dimensions of program PWSCF are:
Max number of different atomic species (ntypx) = 10
Max number of k-points (npk) = 40000
Max angular momentum in pseudopotentials (lmaxx) = 3

gamma-point specific algorithms are used

Subspace diagonalization in iterative solution of the eigenvalue problem:
a serial algorithm will be used

G-vector sticks info
-----
sticks:  dense  smooth   PW   G-vecs:  dense  smooth   PW
Sum      6369   6369   1597      382323  382323  47833

```

Results

## Contact us

We hope this tutorial was helpful. Please let us know if you have any questions or any problems using Galileo. Your feedback is extremely important to us. Contact us anytime at [matthew@hypernetlabs.io](mailto:matthew@hypernetlabs.io) or [alexander@hypernetlabs.io](mailto:alexander@hypernetlabs.io).