pySpawn User Manual

Version 0.0

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# Introduction

## What Is pySpawn?

PySpawn is an implementation of ab initio multiple spawning (AIMS) in python.

## What Are The Primary Goals of pySpawn?

PySpawn is designed with the following goals in mind:

* The core of pySpawn is designed to be small and thus easy to maintain.
* PySpawn is designed to be extensible. At present, the following functionality can, in principle, be implemented without modification of existing code:
  + Interfaces to additional electronic structure codes/methods
  + Quantum mechanical Hamiltonians (e.g. adiabatic representation vs. diabatics vs. diabatized Gaussians on adiabatic surfaces, interpolated vs. analytical derivative couplings, spin orbit Hamiltonian, explicit light field)
  + Quantum mechanical integrators
  + Classical mechanical integrators
* PySpawn is designed with large, shared computer resources in mind. Specifically, pySpawn is designed to facilitate
  + Subdivision of large jobs into small chunks
  + Restart
  + Parallelization

## Features

* Runs FMS/AIMS jobs in the adiabatic representation
* Derivative couplings computed by norm preserving interpolation (NPI)
* PySpawn interfaces to a development version of TeraChem via a protobuf interface

## Citing pySpawn

If you use pyspawn please cite the following:

“pySpawn is a nonadiabatic molecular dynamics software package written by Benjamin G. Levine”

This citation will eventually be replaced by a true publication.

## License

PySpawn is distributed under the MIT License:

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# Obtaining and Installing PySpawn

## Obtaining PySpawn

PySpawn can be obtained from github: <https://github.com/blevine37/pySpawn17>.

## Dependencies

PySpawn requires the following dependencies:

* numpy
* h5py
* google (for TeraChem interface)
* protobuf (for TeraChem interface)

## Installing PySpawn

Having obtained pySpawn and installed the necessary dependencies, cd to the top directory and run

python setup.py install

# Guided Tour of Your First pySpawn Simulation