

Numerical solution of the Vlasov-Poisson problem

April 7th. 2015

Source

- FTP <ftp://ifts.zju.edu.cn> under `incoming/zonca15`.

Part I

An Introduction to Chen Zhao's PIC1D Code

Formulation

- Equation of motion for electron

$$\begin{cases} \frac{dx_i}{dt} = v_i \\ \frac{dv_i}{dt} = -\frac{eE(x_i)}{m_e} \end{cases}$$

- Electric potential

$$E(x) = -\frac{d\phi(x)}{dx}$$

- Poisson equation

$$\frac{d^2\phi(x)}{dx^2} = -\frac{e}{\epsilon_0}[n_0 - n(x)]$$

Formulation

- Weight

$$S_x(x - x_{j\alpha}) = \begin{cases} \frac{h_x - |x - x_{j\alpha}|}{h_x^2} & |x - x_{j\alpha}| < h_x \\ 0 & \text{otherwise} \end{cases}$$

Part II

An Introduction to Wenjun Deng's PIC1D Code

Acquisition Code and Required Lib

- Code
 - You can easily get Wenjun Deng's PIC1D code from:
<https://github.com/wenjundeng/pic1dp>
 - I suggest you setup the code on a Linux based system
- Required Lib
 - Compile & Run
 - Fortran 2003 compiler (e.g. GFortran)
 - MPI-2 (e.g. MPICH2 1.4+, OpenMPI 1.5+)
 - PETSc 3.1+
 - Visualization output results
 - Python 2.7
 - Numpy 1.4+
 - matplotlib 1.1+
 - Scipy (Optional, for calculating dispersion relation analytically)

Installation

- See the [Simple Guide to Run Wenjun Deng's PIC1D Code](#).

Code Structure Overview

- All the **functions / procedures** are **prefixed** by the name of the source file that contained them.
- Most of the parameters are in **src/pic1dp_input.F90** with detailed description.

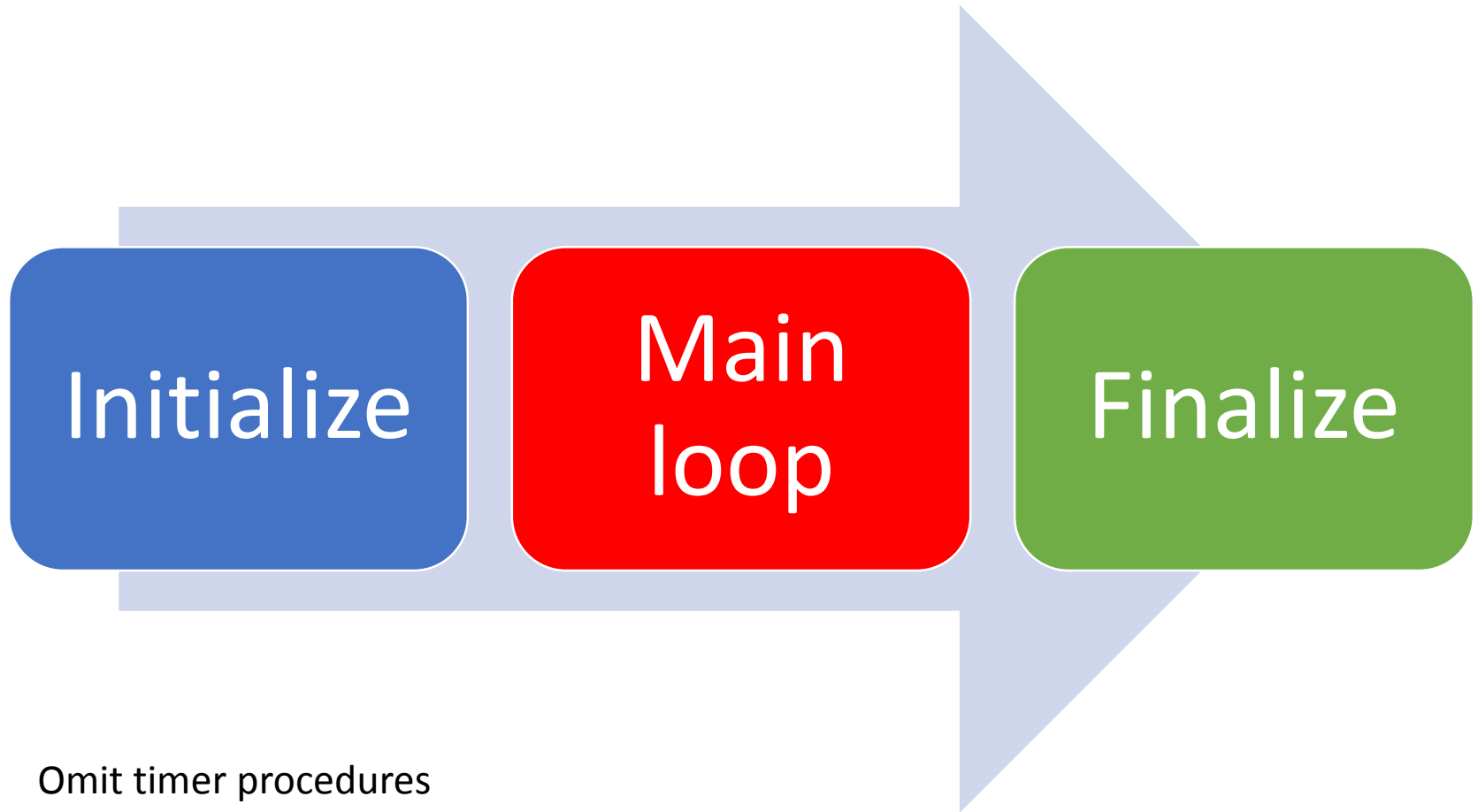
Code Structure Overview

Source files

pic1dp.F90	The main program file.
pic1dp_global.F90	Module for managing global constants, variables and subroutines.
pic1dp_input.F90	Module for managing input parameters and functions. Change input parameters in this file.
pic1dp_particle.F90	Module for managing marker particles.
pic1dp_field.F90	Module for managing field quantities.
pic1dp_interaction.F90	Module for managing field-particle interactions.
pic1dp_output.F90	Module for managing output.
wtimer.F90	A wall clock timer using MPI_Wtime().
multirand.F90	A pseudo-random number generator.

Code Structure Overview

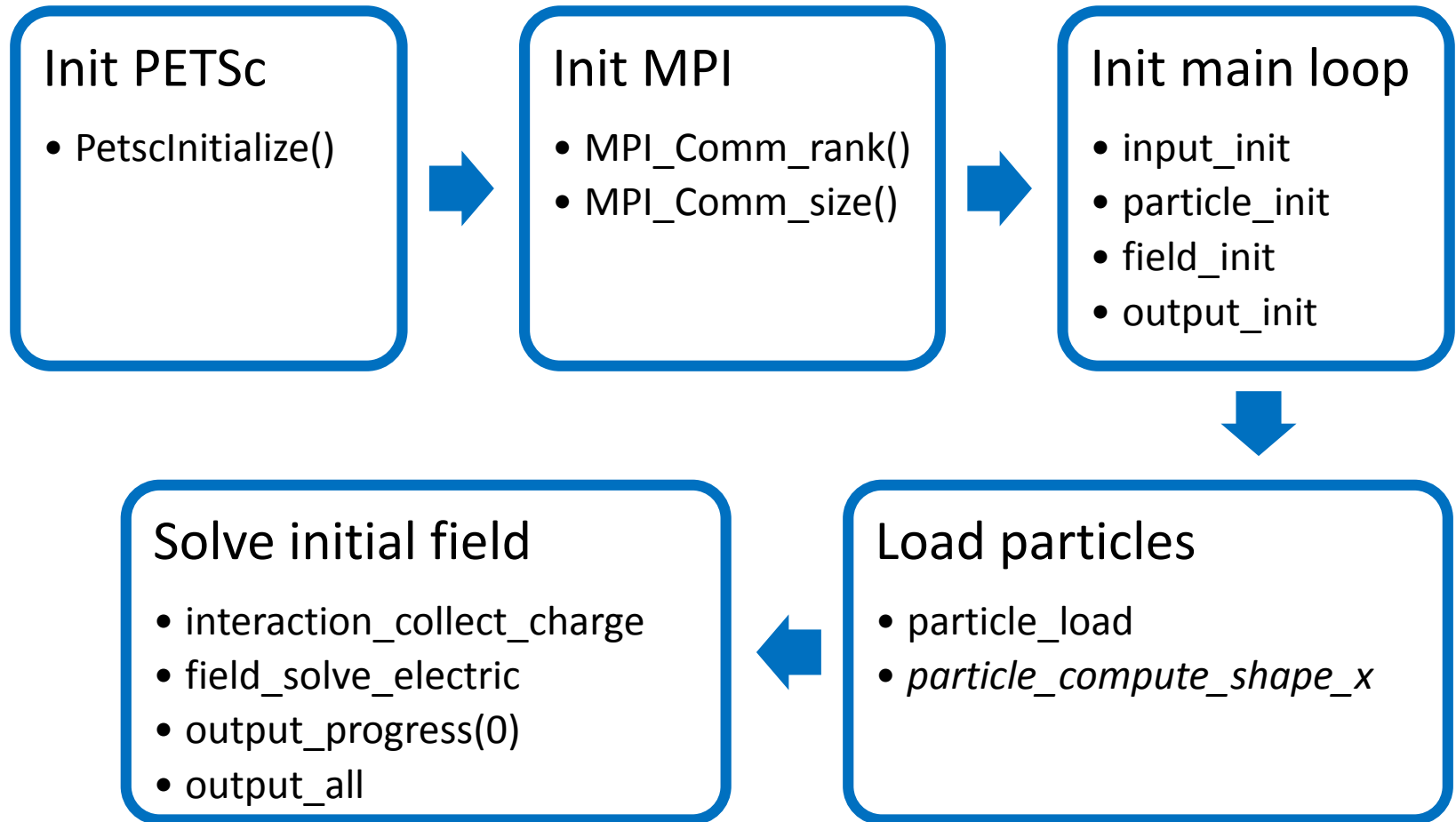
Execution flow



Omit timer procedures

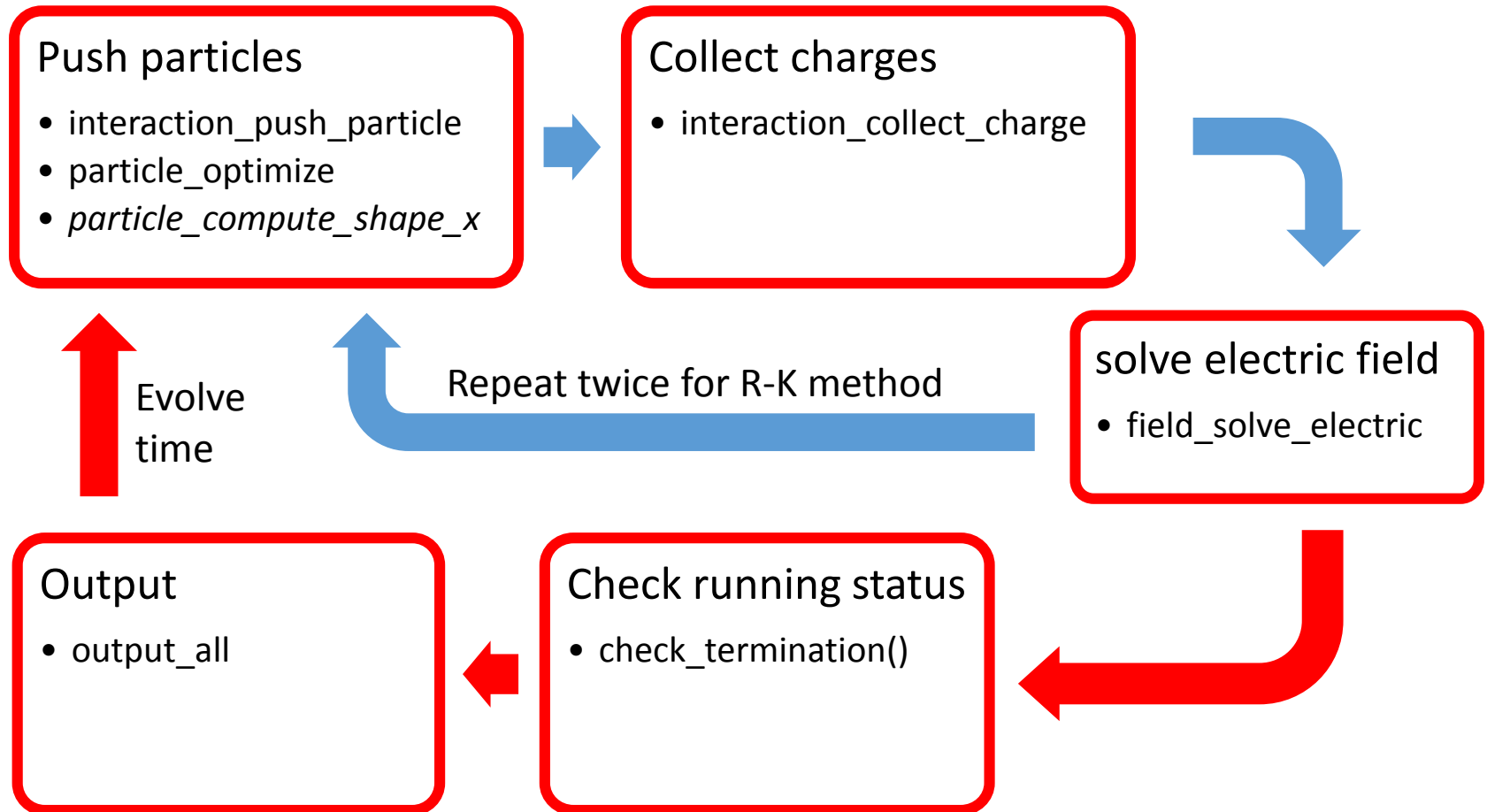
Code Structure Overview

Initialize



Code Structure Overview

Main loop



Code Structure Overview

Finalize

Finalize main loop

- particle_final
- field_final
- output_final



Finalize PETSc

- PetscFinalize()

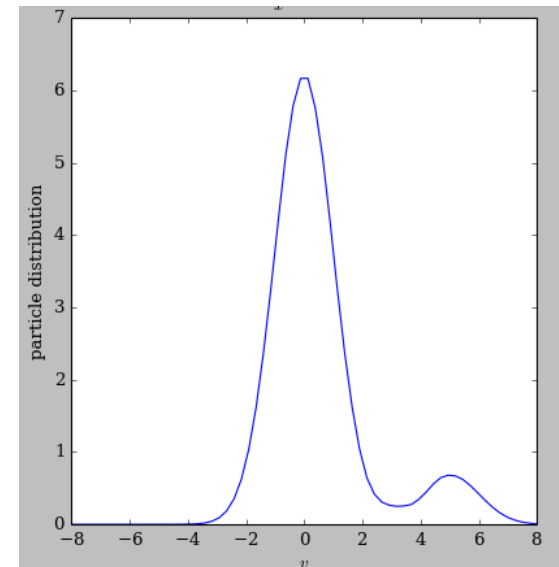
Basic Usage

- Modify input parameters
- Run
- Analyze output results

Basic Usage

Modify input parameters

- Take the default `pic1dp_input.F90` file as an example
 - An electron bump-on-tail instability
 - Section V.A.2 in Siminos et al. (2011) PRE
 - <http://dx.doi.org/10.1103/PhysRevE.83.056402>



Basic Usage

Modify input parameters

termination conditions	physical parameters	initial condition	numerical parameters
• input_ntime_max • input_time_max	• input_linear • input_lx • input_iptcldist • input_nspecies • input_nspecies_xxx • input_nmode • input_modes	• input_init_nmode • input_init_mode • input_init_mode_cos • input_init_mode_sin	• input_deltaf • input_dt • input_nparticle_max • input_species_nparticle_init • input_imarker • input_v_max • input_nx • input_nv • input_iptclshape

Basic Usage

Modify input parameters

marker particle optimizations	random number generator (multirand) parameters	output parameters
• input_nmerge • input_tmerge • input_thshmerge • input_nremove • input_tremove • input_typeremove • input_thshremove • input_remove_frac • input_nsplrit • input_tsplrit • input_thshsplrit • input_split_ngroup • input_split_dv_sig_frac	• input_multirand_al_int • input_multirand_seed_type • input_multirand_warmup • input_multirand_selftest	• input_verbosity • input_output_interval • input_nx_opd • input_nv_opd

Basic Usage

Run

- Compile
 - First time compile
 - Run `make` in the terminal
 - Otherwise
 - Run `make cleanbuild` first in the terminal, then run `make`
- Execute
 - On single machine
 - Run `make run` in the terminal
 - On cluster
 - Change the executive program name in the job script

Basic Usage

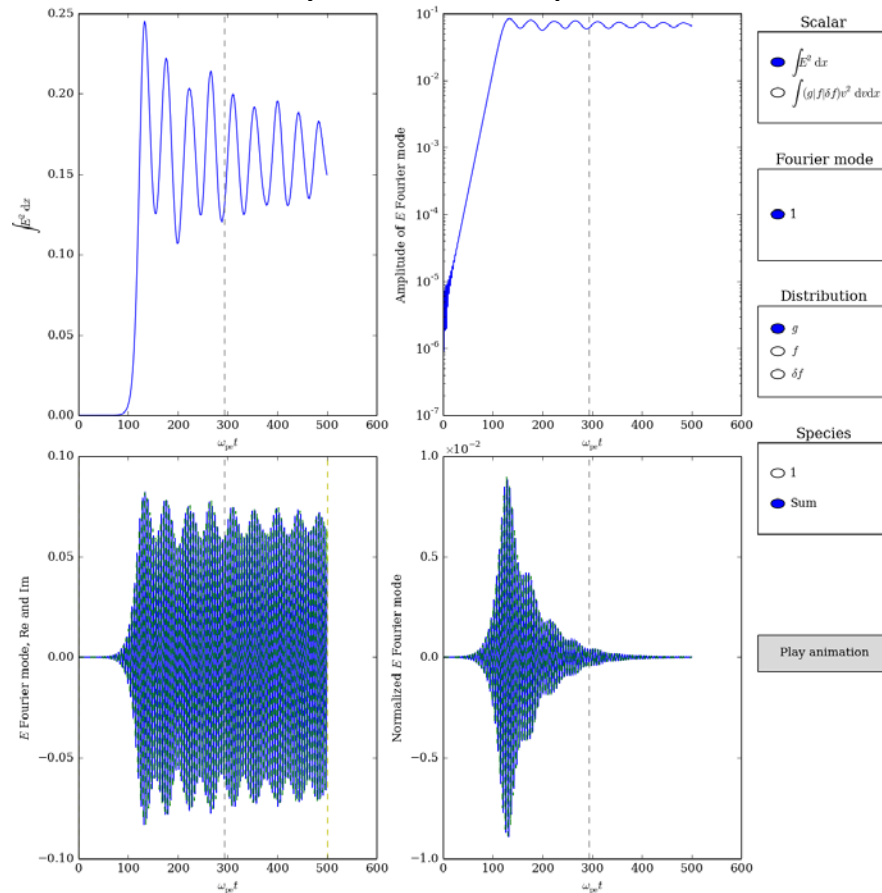
Analyze output results

- The running progress is logged to `run/pic1dp.log`.
- The output data is written to `run/pic1dp.out`.
- To run the visualization app, execute `make visual`.

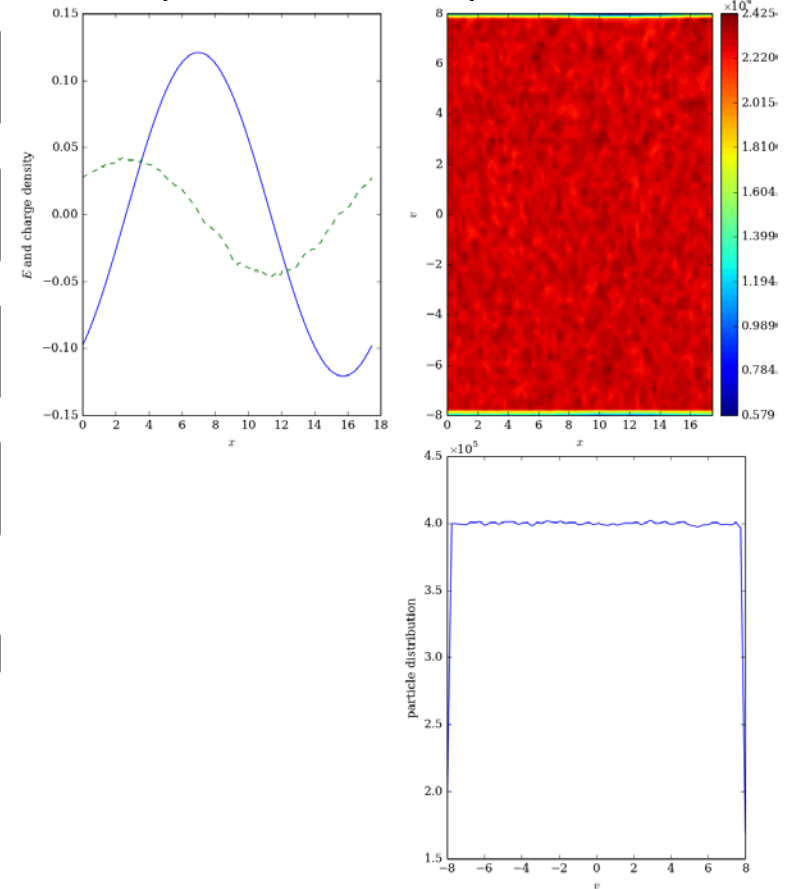
Basic Usage

Analyze output results

time history of various quantities



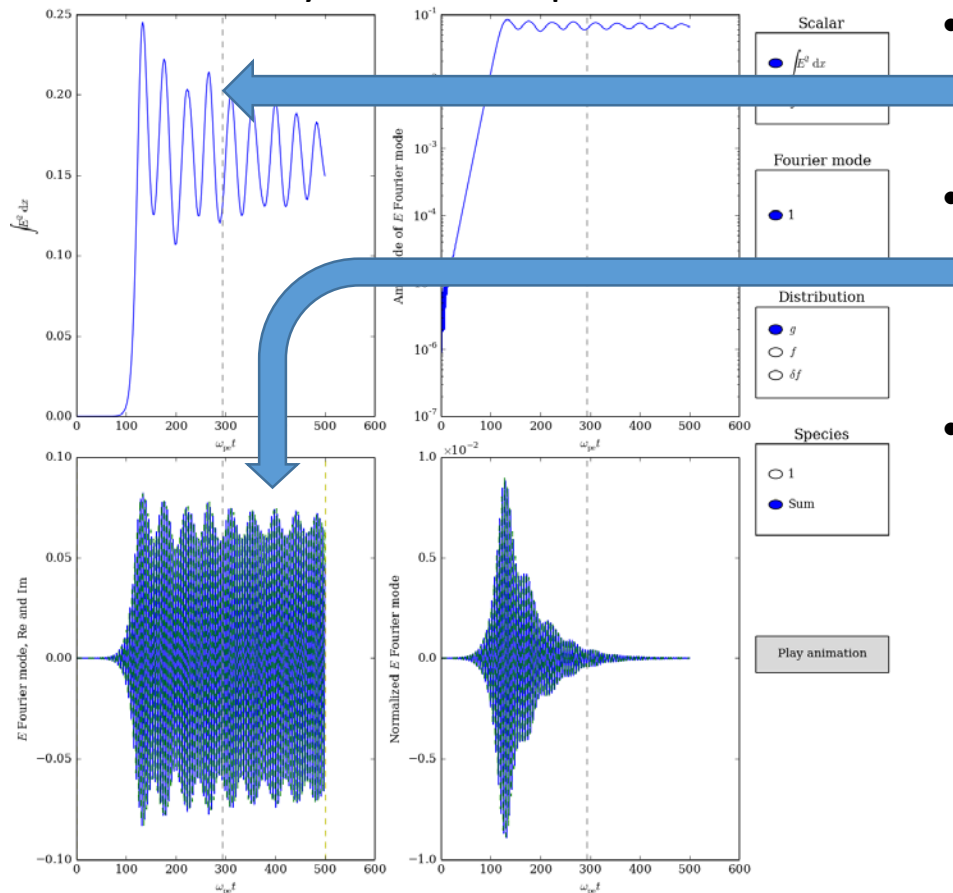
quantities on a specific time



Basic Usage

Analyze output results

time history of various quantities



- Clicking in the upper left panel changes the time for the right 3 panels.
- Clicking and dragging in the lower left panel chooses a time range for the 2 panels in the 2nd column.
- The growth/damping rate calculation is based on the plot end points of the 2nd column upper panel.

Formulation

- See 1D electrostatic δf particle-in-cell (PIC) formulation in vector-matrix form.