

Operation guide for plotting of RANS X transport, flux and variance equations based on ransdat data from PROMPI

Date	Version	Person	Change
25/March/2018	1.0	Miroslav Mocak	Initial instructions for plotting of ransX equations (code still needs to be tested with proper turbulence data)
20/April/2018	1.1	Miroslav Mocak	Introducing parameter α for switching between spherical and Cartesian geometry

Prerequisite:

- Linux/Windows operating system
- Python 2.7 + ipython

1. Go to https://github.com/mmicromegas/PROMPI_DATA
2. To a dedicated directory <DIR> download the following files (or download the whole repo but some classes and still not finished):
 - PROMPI_RANS_xnu.py (class for plotting of the ransX equations. Mapping between calculated fields and equation terms can be found in the ransXtoPROMPI.pdf)
 - PROMPI_data.py (class for reading ransdat data)
 - CALCULUS.py (class with useful calculus methods)
 - plot_rans_xnu.py (control script for plotting)
 - rans_tseries.py (script for calculating of time-averaged fields from ransdat)
3. In <DIR> create two sub-directories DATA and RESULTS
4. Copy your *.ransdat and *.ranshead data to folder DATA
5. Open rans_tseries.py and adjust the following parameters:
 - trange (line 9). Restrict time-range of your ransdat data in DATA folder for time-averaging.
 - tavg (line 10). Set time-averaging window (at least 2 convection turnover timescales)
6. From <DIR>, start ipython and execute > run rans_tseries.py
7. After successful completion of rans_tseries.py, the time-averaged data are stored in a file called tseries_ransout.npy. Check if the file was in the <DIR> created successfully.

8. Open `plot_rans_xnu.py` and adjust the following parameters according to your needs:
 - `intc` . Choose index of central time for which you wish to plot the `ransX` fields.
 - `inuc` . Choose ID of the element for which you want to plot the `ransX` fields. It has to have the format `00xx` , for example `0001` is neutrons, `0002` is protons, `0003` is `he4`, `0004` could be `c12` (all depends on your network)
 - `ig` . Enter geometry of your simulation (1 is Cartesian, 2 is spherical)
 - `LGRID` . Choose whether you want to limit your x-grid. Good if you want to get rid of boundary noise (1-true, 0-false)
 - `xbl`, `xbr` . Set left/right radius for which you want to limit x-grid in your plots. Y-axis will adjust itself automatically.
 - `lc` . Optional. Estimated size of convection zone. This is still work in progress. Set it properly, if you want to get Eulerian diffusivities right.
9. From `<DIR>`, start `ipython` and execute `> run plot_rans_xnu.py`
10. Wait for the plots to be displayed.
11. If you wish to display also radial profiles of element density, flux and variance, uncomment the following lines `#RANSX.plot_Xrho(xbl,xbr,inuc,data_prefix)`,
`#RANSX.plot_Xflux(xbl,xbr,inuc,data_prefix)`, `#RANSX.plot_Xvariance(xbl,xbr,inuc,data_prefix)`
12. If you want to display various diffusivities for the target element, uncomment `#RANSX.plot_X_Ediffusivity(xbl,xbr,inuc,data_prefix)`