FluxEngine v3.0: Changes and User Guide

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1 Overview and major changes

The major changes, new to FluxEngine version 3.0 include:

- Standalone execution Various changes which make it considerably easier to run locally (i.e. without access to a remote server). A big part of this is that Flux-Engine no longer requires input files for data layers which will not be used given a the configuration options specified. For example, you no longer need to specify the location of sea surface skin temperature data, if you're using foundation temperature. This means you do not need access to data which might be irrelevant to your particular use case in order to run FluxEngine.
- Greater flexibility Use of Unix-like glob pattern matching to specify input data means file names no longer need to conform to particular formats. This makes it easier incorporate new and user supplied sources of data. A modular and extendible system for data pre-processing has been added to make it easier to use data from a wider range of sources.
- Substantial restructuring The source code has been substantially restructured. While this is an ongoing process, having a consistent method for reading, processing and outputting data layers reduces the scope for bugs, provides a simpler conceptual model to understanding and extending FluxEngine functionality. For example, additional data pre-processing and gas transfer velocity calculations can be added and tested without modifying the core FluxEngine code.
- Python workflow Python, and some Python modules are now the only dependencies needed by FluxEngine. The Perl driver script has been transcribed into Python and updated. FluxEngine can run from the command-line or imported as a Python module providing a simple API interface for more advanced use-cases.
- User-friendly installation and validation Installation scripts will automatically install missing dependencies for MacOS and Linux based systems. Installation instructions are provided for Windows users. Validation scripts are provided to ensure FluxEngine is operating correctly, and can be used to validate modifications.

1.1 Execution time

Simple benchmarking was conducted using an Intel Core i5 2.7GHz processor with 8GB RAM running MacOS El Capitan. A one year (2010) run using the SOCATv4 validation configuration (Nightingale 2000 k parameterisation with process indicator layers off) took approximately 6 minutes to complete using FluxEngine v3.0. This is compared to over 9 minutes for the same configuration using FluxEngine v2.0. This speed-up can be largely attributed to the removal of non-required input data layers. Run time will therefore differ depending on the options specified by configuration files. For example the Takahashi 2009 validation run took just 4 minutes to complete.

1.2 Feedback and bug reporting

Feedback is greatly appreciated both in terms comments about which aspects of running and using FluxEngine were not intuitive or poorly explained, as well as suggestions about useful potential features to include in the next release. These can be e-mailed to Tom Holding at t.m.holding@exeter.ac.uk. Bugs can be reported via our GitHub page by opening an issue: https://github.com/oceanflux-ghg/FluxEngine/issues or by e-mailing Tom (see above).

2 Downloading FluxEngine v3.0

FluxEngine can now be ran on a standalone computer so you don't need access to a remote host. The latest version can be downloaded from https://github.com/oceanflux-ghg/FluxEngine/archive/master.zip, or if you use Git you can clone the FluxEngine repository here: https://github.com/oceanflux-ghg/FluxEngine.git

After downloading unzip the files into a location of your choice. This should leave you with the directory hierarchy shown below:

```
FluxEngine (or FluxEngine-master) - referred to as the root folder

configs - contains validation and example configuration files

data - input data required to run validations and reference output

fluxengine_src - FluxEngine source code

tools - additional tools which can be run separately to

FluxEngine
```

3 Installing dependencies

One of the changes from v2.0 is that FluxEngine can be run entirely using Python 2.7. That being said there are several modules which are not part of the Python 2.7 standard library. These are netCDF4, numpy and scipy.integrate. If you don't already have these installed, you'll need to install them before running FluxEngine. They can be installed manually (the recommended way to do this is using pip which usually comes bundled with Python 2.7) or they can be installed using the dependency installation scripts which can be found in the root directory for MacOS and Linux.

3.1 MacOS

An installer script is available for MacOS which will automatically install dependencies. While it's provided with no guarantees, the script shouldn't overwrite anything without asking you first. To run it, open Terminal, navigate to the FluxEngine root folder and run

```
./install_dependencies_macos.sh
```

If you see an error about not having permission to execute, chances are you'll need to add this permission before running the install script. To do this run the following:

```
./sudo chmod +x install_dependencies_macos.sh
```

3.2 Linux

An installer script is available for Debian based Linux systems and will automatically install the dependencies needed to run FluxEngine. It has only been tested on Ubuntu 16.4, and is provided with no guarantees. To run it, open Terminal, navigate to the FluxEngine root folder and run

```
./install_dependencies_ubuntu.sh
```

If you see an error about not having permission to execute, chances are you'll need to add this permission before running the install script. To do this run the following:

```
./sudo chmod +x install_dependencies_ubuntu.sh
```

3.3 Windows

An installer script isn't provided for Windows but the following steps should enable you to install the dependencies required to run FluxEngine.

Step 1) Since Python 2.7 doesn't come with Windows you may not have it installed. If you already have Python 2.7 go to step 2, if not then go to https://www.python.org/downloads/ and download the installer. At the time of writing the most recent version of Python 2.7 was 2.7.14. Do not download version 3.x, Flux-Engine will not run on it. Once downloaded run the file and you'll be guided through the installation process. Make a note of your installation directory (it will probably be something like C:\Python27\ or Users\<your username>\AppData\Local\Programs\Python\Python27).

Step 2) Add Python and the Python Scripts folder to your PATH environment variable. This will allow you to run python scripts without having to navigate to specify the whole Python installation directory each time. To do this in Windows 10, follow these instructions: https://www.youtube.com/watch?v=uXqTw5e00Mw using the installation directory from the previous step.

Step 3 Use pip to install three Python modules: netCDF4, numpy and scipy.integrate. Pip is a package manager, which comes with Python and allows you to add/remove additional libraries. First, open a new command prompt window (or close and reopen an existing one - this is important or your PATH environment variable won't have updated). Type the following commands:

```
pip install --upgrade pip
pip install numpy
```

pip install netCDF4

You should now be ready to run FluxEngine. More detailed information on installing and configuring Python 2.7 on Windows can be found here: https://docs.python.org/2.7/using/windows.html

4 Validating FluxEngine

Two validation scripts come bundled with the FluxEngine. These make it easy to verify that everything has installed correctly and that FluxEngine is producing correct output. These scripts compare output generated by your local copy of the FluxEngine with a known reference, and will report any discrepencies.

To validate using Takahashi2009 (T09) and/or SOCATv4 data run the following command/s from FluxEngine's root directory, respectively:

```
python validate_takahashi09.py
python validata_socatv4_sst_salinity_gradients_N00.py
```

Alternatively, you can write python scripts to validate against your own reference data sets using the functions provided in the validation_tools.py file in the tools directory.

5 Running FluxEngine v3.0

In previous versions FluxEngine was ran using a bash script which invoked a Perl script to parse a config file and run the core Python FluxEngine code. This has been considerably simplified in version 3.0 so that everything is handled by Python. This has the advantage of reducing dependencies and moving toward a more API-like interface for executing FluxEngine. An execution script (ofluxghg_run.py, found in the root directory) is still provided, which functions similarly to the previous Perl script (described below). To run the FluxEngine for

5.1 Using the ofluxghg_run command line script

ofluxghg_run.py is found in the FluxEngine root directory can be be used as a standalone tool to run the FluxEngine, for example

```
python ofluxghg_run.py configs/example_config.conf -l
```

will run the flux engine using the options specified in a configuration file called example_config.conf located in the configs subdirectory of the current working directory. This assumes ofluxghg_run.py is being executed from the FluxEngine root directory, but it can be executed from anywhere, provided that 1) the path of the config file is reachable (absolute file paths can be used), 2) ofluxghg_run.py can find the fluxengine_src folder (either it has been added to your Python paths, moved somewhere

which is already in your Python path, or it is a subdirectory of the current working directory) and 3) the input data specified in your config file is reachable from the current working directory (these can also be absolute paths). The -1 option tells FluxEngine to run without process indicator layers, this reduces the time taken to run and results in smaller output files.

The general syntax for running ofluxghg_run is

```
python ofluxghg_run.py config [-options]
```

where [-options] is an option list of options to change how FluxEngine runs. A list of valid options can be listed by using the -h flag, e.g. python ofluxghg_run.py -h. Some options require you to specify a value, for example to set the start and end years you can use

```
python ofluxghg_run.py configs/example_config.conf -year_start 2000 -year_end
2010 -1
```

to run FluxEngine for the years 2000 to 2010. Note: A useful option for testing is -D1 which will only run FluxEngine for a single month. This allows you to check the output and any error messages before committing to a longer run.

5.2 Driving FluxEngine from your own Python scripts

Some tools are provided which allow you to write Python scripts to run FluxEngine in custom ways. To do this you should import the fluxengine_src.fe_setup_tools module and use run_fluxengine function. An example of how to use this function is can be found in ofluxghg_run.py. For more control over how FluxEngine is driven, e.g. if you need to use a single configuration file and overwrite a subset of parameters to run a suite of similar simulations, see the tools additional functions in fe_setup_tools in the fluxengine_src directory. Note that these tools are intended for use by users who are proficient in Python. They haven't undergone intensive testing and likely to contain bugs (especially in Windows environments). Bugs can be reported as described in section 1.2.

6 Creating and modifying configuration files

The format of configuration files has changed a bit from the previous version. Example annotated configuration files can be found in the configs directory. Variables are specified in the config file by name, followed by an equals sign (=) and then the value. In contrast to previous versions the equals sign is now required, but this means that whitespace can be used in variable values. Both variable names and values are case sensitive, and comments can be added using the hash (#) character. Variables can be specified in any order although it is convenient to group related variables together. An example definition of two variables is given below:

6.1 Configuration file specification

src_home

The file path to the fluxengine_src directory. Note that this is no longer used in version 3.0, but is reserved and may be used in the future to distinguish between different versions of FluxEngine which may be installed on the same computer.

flux_calc

The flux calculation to perform. Valid options are:

- bulk i.e. $F = k\alpha_W (pCO_{2_W} pCO_{2_A})$, where F is the air-sea flux, k is the gas transfer velocity, α_W is the solubility of the gas in sea water, pCO_{2_W} is the partial pressure of CO_2 in the surface sea water and pCO_{2_A} is the partial pressure of CO_2 in the atmosphere.
- rapid As described in Woolf et al. (2012) Journal of Geophysical Research: Oceans. $F = k(\alpha_{\scriptscriptstyle W} p C O_{2_{\scriptscriptstyle W}} \alpha_{\scriptscriptstyle A} p C O_{2_{\scriptscriptstyle A}})$.
- equilibrium As described in Woolf et al. (2012) Journal of Geophysical Research: Oceans.

use_sstskin, use_sstfnd, sst_gradients, cool_skin_difference

Determines the source/s of sea surface temperature to use (skin or foundation or both). Valid values for use_sstskin, use_sstfnd and sst_gradients are yes or no. At least one of use_sstskin or use_sstfnd must be set to yes. If only one is specified and sst_gradients is set then the other is estimated using the following equation:

```
sstskin = sstfnd - cool_skin_difference
```

where cool_skin_difference is the difference in temperature between the foundation layer and the skin layer (in Kelvin). The default is 0.17K (see Donlon et al. 2002).

saline_skin_value

Saline skin value is added to salinity. It is an optional entry and will default to 0.0 if not specified.

axes_data_layer, latitude_prod longitude_prod time_prod

Specifies the data layer from which to extract the latitude, longitude and time data from. axes_data_layer must be the name of a data layer, e.g. sstskin, and all other input data layers will have their dimensions checked for consistency with the named data layer.

pco2_reference_year, pco2_annual_extrapolation

These are optional entries which can be used to specify a reference year from which pCO_2 / fCO_2 can be adjusted for. This applies an annual correction according to

datalayername_path, datalayername_prod

These define each input data layer. A full explanation with worked examples is provided below (section 6.3).

These are optional variables which control whether random noise is added for each of their respective data layers. Valid values are yes or no. Note that currently the magnitude of noise must be modified directly in the source code (this is not advised unless absolutely necessary). This system will be replaced with a more flexible system in the near future.

bias_datalayername, bias_datalayername_value

Setting bias_datalayername will add a constant value (defined by bias_datalayername_value) to the named data layer. This is applied after any random noise is added. Currently the following data layer names are supported: windu10, sstskin, sstfnd and pco2. Valid options for bias_datalayername_value are yes or no and default to no if not supplied. bias_datalayername_value must be a valid numeric value and defaults to 0.0 if not supplied. Future releases are likely to expand support to all data layers.

bias_k_bias_k_percent, bias_k_value, bias_k_biology_value, bias_k_wind_value
These variables control the bias applied to k (the gas transfer velocity). bias_k and
bias_k_percent but be set to either yes or no and control whether any bias is applied at all and whether the bias is added as an absolute value or as a percentage
of the original value, respectively. Default values for both of these variables are no.
bias_k_value requires a numeric variable (the default is 0.0) and controls the magnitude of the bias (whether as a percentage or absolute value). Bias is only added to
k if the value of windu10 at the same point in space is above a threshold specified
by bias_k_wind_value (the default value is 0.0). Similarly the corresponding value for
biology is below bias_k_biology_value (the default is 0.0).

k_parameterisation

This controls the way that the gas transfer velocity (k) is calculated. FluxEngine comes bundled with a number of k 'functors' - self-contained Python classes which take as input data layers, and write output to one or more data layers (typically the k datalayer). For a list of available parameterisations you can run the ofluxghg_run.py script with the -list_parameterisations option. Alternatively you can see the Python implementation of each k functor in rate_parameterisation.py file located in the fluxengine_src directory. These can be referred to by name in configuration files. User defined parameterisations can be added to this file and assigned to k_parameterisation by name in the config file. Before writting new k functors you should read the guide for developers (section 8), as this describes best practices and provides information on potential pitfalls.

bias_sstskin_due_rain,

bias_sstskin_due_rain_value,

bias_sstskin_due_rain_intensity, bias_sstskin_due_rain_wind

These control whether and how rain (the rain data layer) influences sea surface temperature. bias_sstskin_due_rain_value turns this feature on or off (valid values are yes and no, with the default being no). If this feature is turned on, a constant bias (bias_sstskin_due_rain_value, default value of 0.0) will be added to sea surface temperature anywhere that rain intensity is greater than bias_sstskin_due_rain_intensity (default value 0.0) and wind intensity (the windu10 data layer) is less than bias_sstskin_due_rain_wind (default is 0.0).

rain_wet_deposition

This option enables wet deposition. Valid values are yes or no, and the default value is no.

k rain linear ho1997

This option enables a linear additive gas transfer velocity term due to rain. A description of the method can be found in Ashton *et al.* (2016) . Valid values are yes or no, and the default value is no.

k_rain_nonlinear_h2012

This option enables a nonlinear additive gas transfer velocity term due to rain. A description of the method can be found in Ashton *et al.* (2016) and Harrison *et al.* (2012). Valid values are yes or no, and the default value is no.

output_dir

This specifies the directory that will be used by FluxEngine to put the output netCDF files in. If the directory doesn't exist it will be created. If other files with the same names and directory structure as the output files already they will be overwritten. A directory will be created for each year that FluxEngine runs, within which a subdirectory will be created for each month run. This directory structure is for convenience, since it is the same as that required by the ofluxghg_net_budget.py tool (see the section 7 for a description of bundled tools).

Full meta data, including a list of data layer names recognised by FluxEngine, default values and expected data types can be accessed in the settings.xml file in the fluxengine_src directory. Note: It is strongly recommended that you do not modify this file.

6.2 k parameterisation specific variables

Several k_parametrisation options require additional variables to be specified which change the way that the gas transfer velocity is calculated. These must be defined in the configuration file. In particular k_generic requires a Schmidt number to be defined k_generic_sc (valid values are 600.0 and 660.0) as well as weightings for each order of the generic gas transfer velocity equation, i.e. k_generic_a0, k_generic_a1, k_generic_a2

and k_generic_a3.

kt_OceanFluxGHG, kt_OceanFluxGHG_kd_wind and k_Wanninkhof2013 require that kb_weighting and kd_weighting be specified. These define the weighting for the bubble and direct components of the gas transfer velocity, as described in Goddijn-Murphy et al., (2015).

Other custom or third-party k parametrisations may require other variables to be specified and you should consult any documentation or guidance specific to the parametrisation being used, or examining the initialiser function (__init__) of the relevant parametrisation functor the rate_parameterisation.py file.

6.3 Data layers

The term 'data layer' is used to describe a 2D dataset, and any accompanying metadata, which is used by FluxEngine either as input, an intermediate product or output. Configuration files must specify all the input data layers which will be needed, but you only need to specify the input data layers which will be used given the specific options youve selected. For example you do not need to you do not need to specify biology input files if the 'process indicator layers off' is set using the -1 flag, and you do not need to specify a sea surface skin temperature data layer if use_sstskin = no is set in the config file. If you try to run the FluxEngine without the required inputs youll get an error message telling you which input data layer was missing. If you specify data layers which are not needed for the calculation options specified they will simply be added to the output data layer.

To specify an input data layer the configuration file must specify a minimum of two attributed: a path to the netCDF / .nc file, and a prod (variable name within the netCDF file). The path can be absolute or relative. Windows users might experience some problems using absolute paths. If you have problems with this please let me know (tom: t.m.holding@exeter.ac.uk) and Ill try to fix it. Path are specified in the config file using the data layer name with the _path suffix, like so:

datalayername_path = path/to/data/file.nc

One important change in this version of FluxEngine is that you should specify the path including a the filename for the netCDF file, rather simply a directory. This provides greater flexibility when working with data from many different sources. To help with this two standard Unix glob patterns can be used to specify patterns of file names: ? and *. These will match any single character/digit, or any number of characters/digits (including no characters), respectively. For example to specify the location of ice coverage data you could use:

ice_path = path/to/data/20100101_???-ice*.nc

This will match any file with a name that starts with 20100101_ followed by any three characters/digits, then the characters -ice, followed by any number of characters/digits and ending in .nc. Note that these cannot be used to match directory names and can

only be used to specify the pattern that FluxEngine will use to match the netCDF file itself.

Several additional tokens can be used to specify file or directory paths which vary according to the current year or month being simulated, and allows different input files or directories to be used for different month/year combinations. The following tokens are supported:

- <YYYY>-four digit year, e.g. 2010
- <YY>-two digit year, e.g. 10 for 2010
- <MM> two digit numerical month, e.g. 01 for January
- <Mmm> three character abbreviation of the month, e.g. Jan for January
- <MMM> three character upper-case abbreviation of the month, e.g. JAN for January
- <mmm> three character lower-case abbreviation of the month, e.g. jan for January

Since these tokens can be used in directory names it can be useful if, for example, you have data for multiple years for one data layer, but only have data covering one year for another. In this case you would specify the file path of the first data layer using the appropriate year token, and the second by hard-coding the year string into the file path. Supposing we only has sea surface foundation temperature for the year 2010, but ice coverage data for each year we're interested in, this might look as follows:

```
ice_path = path/to/data/<YYYYY>/<YYYYY><MM>_???-ice*.nc
sstfnd_path = path/to/data/2010/2010<MM>_OCF-SST-GLO-1M-???-REYNOLDS.nc
```

The second required attribute of a data layer is its product (or 'prod'). This is the name of the variable within the netCDF file. It is specified in the configuration file by using the _prod suffix with the data layer name. The minimal specification (in this case for the 'ice' data layer) could therefore looks like this:

```
ice_path = path/to/data/<YYYY>/<YYYY><MM>_???-ice*.nc
ice_prod = sea_ice_fraction_mean
```

6.4 Optional data layer attributes

There are several optional attributes which can be configured for each data layer. They can be set using the same datalayername_suffix notation used for the path and products above. These are:

- _stddev_prod product name of a variable containing standard deviation data for the data layer
- _count_prod product name containing the number of samples used to calculate standard deviation

- \bullet _netCDFName the variable name used to label this data layer in the output netCDF file/s
- _units a string description of the units
- _minBound minimum allowed value)
- _maxBound maximum allowed value
- _standardName short standardised description of the variable/data layer
- _longName human readable description of the variable/data layer

For example to overwrite the minBound and maxBound attributes for the ice coverage data layer as well as rename it in the output files you can add the following lines to a configuration file:

```
ice_minBound = 0.0
ice_maxBound = 100.0
ice_netCDFName = ice_percent
```

Any value outside of this range will be replaced with missing values.

Default metadata values are stored in settings.xml in the fluxengine_src directory, but shouldn't usually be modified because this would change the behaviour for all FluxEngine runs. These values can always be overwritten in the configuration file as described above.

6.5 Data layer preprocessing

It is sometimes convenient to apply some pre-processing to a data layer before it is used for any computations. There is an additional data layer attribute which allows the user to specify a list of functions to be applied immediately after the data layer read in. A number of simple preprocessing functions are bundled with FluxEngine (these can be listed by running ofluxghg_run.py with the -list_preprocessing flag, or by viewing the functions directly in the data_preprocessing.py file in the fluxengine_src directory). For users comfortable with python, custom pre-processing functions can be added to data_preprocessing.py. These will be automatically detected when running FluxEngine available to use in configuration files. Before modifying this file you should familiarise yourself with the guide for developers notes in section 8.

To specify pre-processing functions the _preprocessing suffix is used with a list of function names separated by commas. Each function will be applied in the order it appears in this list. For example adding the following line to a config file will first transpose the 2D matrix, then convert from Kelvin to Celsius:

```
sstskin_preprocessing = transpose, kelvin_to_celsius
```

Note that no checks are made to ensure the original values are in Kelbin to begin with, and it is up to the user to ensure that any pre-processing functions are applied appropriately.

7 Bundled tools

The tools subdirectory in the flexengine_src directory contains various tools designed to be used in conjunction with FluxEngine. These are discussed below.

resample_netcdf.py - Resamples a 1° by 1° netCDF data to a 5° by 4° grid.

text2ncdf.py - Converts in situ data to a netCDF file.

ofluxghg_flux_budgets.py - Calculates integrated net fluxes.

compare_net_budgets.py - Simple functions which compare net budgets between two runs. Examples of their use can be found in the validation scripts (see section 4). To use these you must first import them into a python project using import fluxengine_src.tools.compare_net_budgets.

validation_tools.py - This file contains several functions to perform common validation tasks, such as comparing the netCDF outputs of two FluxEngine runs, or calculating and comparing the global flux budgets between a new run and a reference data set. To access these functions they must be imported into a python project (e.g. by using import fluxengine_src.tools.validation_tools). The function which is more likely to be useful is validation_run. A description of how to use this function is provided in validation_tools.py.

8 Guide for developers

In FluxEngine v3.0 we have made changes to make it easier for people to modify and extend the functionality of FluxEngine. We have implemented a more consistent structure to the code and added the ability to easily extent certain aspects of FluxEngine's functionality without modifying the core code. The following sections provide some background information on how to do this. This is intended for users who are comfortable programming in Python.

8.1 Adding pre-processing functions

Pre-processing functions are defined in data_preprocessing.py in the fluxengine_src directory. When parsing _preprocessing entries in configuration files FluxEngine searches the function names defined in this file, and so any function which is added will be immediately available for use as a pre-processing function. However, there are certain requirements for the function to operate harmoniously with FluxEngine. These are listed below:

- Pre-processing functions must have a single argument, and this must be the DataLayer instance which corresponds to the input data layer which is being transformed. Detailed of the DataLayer class can be found in DataLayer.py.
- DataLayers should be modified in place. Returned values are ignored.
- Pre-processing functions should only modify the (1D) fdata attribute of the DataLayer instance. The exception to this is if it is convenient to modify 2D view of this (the data attribute), in which case you must call datalayer.calculate_fdata() afterwards. This is because, while in most cased fdata is a view of data, in some cases fdata may be a copy of data and hence any changes to data will not be reflected in fdata. fdata is used for all calculations, so it is important that any changes are copied to this attribute.
- If a pre-processing function modifies the dimensions or any of the meta data associated with a *DataLayer* it must also manually update the relevant attributes as these will not be automatically reflected.

8.2 Adding k-parametrisation functors

The gas transfer velocity calculation is fully customisable by adding a 'functor' class to the rate_parameterisation.py file. Functors are classes which are callable. This class must be derived from the KCalculationBase class and implement four functions:

- __init__ Initialises the class. This must, at a minimum, set self.name. You can add any arguments which the class needs to initialise to the function signature and provided they are added as variables with the same name/s in the configuration file they will be automatically passed to the functor during initialisation (see the notes on adding configuration variables below).
- input_names This should return a list of DataLayer names (strings) which are required as inputs to the gas transfer velocity calculation.
- output_names This should return a list of DataLayer names which are modified or written to. These can be existing or new DataLayers. Any non-existing DataLayers will be created in the correct dimensions (but filled with missing values) by FluxEngine prior to performing the gas transfer velocity calculation.
- __call__ This performs the gas transfer velocity calculation, and will contain all the implementation details for your specific case. In addition to self) an argument called data is passed to this function which contains a dictionary of each DataLayer available to FluxEngine. These can be assessed by using the DataLayer name as a key. Note that you should not create new DataLayer instances, add entries to this dictionary or change DataLayers which are not listed by name in the list returned by output_names.

It is best practice to modify the 1D DataLayer.fdata attribute of output DataLayers. If the 2D DataLayer.data attribute is modified you should update

the fdata attribute by calling DataLayer.calculate_fdata for the DataLayers which have changed. This is because, while in most cases fdata is a view of data to avoid unnecessary duplication, this is not guaranteed on all systems.

The final gas transfer velocity output should usually be written to the k data layer, as this is what will be used by FluxEngine to calculate air-sea gas flux. Additionally, it can be a good idea to set the DataLayer.long_name and DataLayer.short_name attributes of k to provide a description of the parameter-isation used because this will be copied to the output netCDF files.)

An example implementation, as well as all the pre-bundled gas transfer velocity functors, can be found in rate_parameterisation.py in the fluxengine_src directory.

8.2.1 Adding configuration variables

You can add variables to the configuration file and these will be immediately available in the flux engine code (encapsulated in the runParams 'namespace'). For example, if you add

```
my_new_var = 100.0
```

to the configuration file. This can be referenced in the FluxEngine code using runParams.my_new_var. This is utilised by different k-parameterisation functors which require additional variables to initialise correctly (see the definition of k_generic in rate_parameterisation.py for an example).

Additional configuration variables are interpreted as a float if they're formatted as a valid float, otherwise they're interpreted as a string. You can specify the type of input which is required, or add constraints and/or metadata associated with a configuration variables by adding an entry to settings.xml (found in in the fluxengine_src directory). Modifying entries for pre-existing variables is not recommended.

To avoid naming conflicts (since all config variables are imported to the same namespace) it is good practice to add a prefix to the name of any custom configuration variables. Typically this should be the name of the k-parameterisation functor it is associated with. For example, the k_generic) functor requires 5 additional variables each of which begin with k_generic (e.g. k_generic_sc).

8.3 Contributing

If you're a git user you can fork the FluxEngine repository at https://github.com/oceanflux-ghg/FluxEngine and if you develop extensions or functionality which might be useful to the wider community, or spot and fix any bugs, you can share them by sending us pull requests. Alternately, if you don't know what any of that means but you've developed an extension or fixed a bug which you think will be useful to others, you can e-mail me at t.m.holding@exeter.ac.uk.