pyeemd Documentation

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This is the documentation of *pyeemd*, which is a Python wrapper for *libeemd*, a C library for performing the ensemble empirical mode decomposition (EEMD), its complete variant (CEEMDAN) or the regular empirical mode decomposition (EMD).

If you have any questions or problems with *pyeemd*, or you wish to propose enhancements, please check out the main page of *libeemd*. Also, if you want more detail on the inner workings of *libeemd* and/or wish to cite it in scientific work, please see the associated article published in Computational Statistics.
1.1 Installing pyeemd

The *pyeemd* module comes with a regular Python distutils installation script, so installing it should be quite straightforward. The only catch is that you need to first compile *libeemd.so*, since *pyeemd* is only a wrapper for that library. Please see the README file distributed with *libeemd* on more details on how to compile *libeemd*, but if you are unpatient and already have the necessary dependencies installed (GCC, GSL), you can just run `make` in the top-level directory of *libeemd* and you are done.

To install *pyeemd* please run:

```
python2 setup.py install
```

In the top-level directory of *pyeemd* (the one with *setup.py*).

If you want to specify an alternative installation prefix, you can do it as follows:

```
python2 setup.py install --prefix=$HOME/usr
```

1.2 Tutorial

After installing *pyeemd* as described in *Installing pyeemd* you are all set to using it with:

```
import pyeemd
```

The three main decomposition routines implemented in *pyeemd* – EMD, EEMD and CEEMDAN – are available as `emd()`, `eemd()` and `ceemdan()`, respectively. All these methods use similar conventions so interchanging one for another is easy.

Input data to these routines can be any kind of Python sequence that *numpy* can convert to an 1D array of floating point values. The output data will be a 2D *numpy* array, where each row of the array represents a single *intrinsic mode function* (IMF).

As an example, the *examples* subfolder of *pyeemd* contains a file `ecg.csv`, which contains ECG (electrocardiogram) data from the MIT-BIH Normal Sinus Rhythm Database. The data is in CSV (comma separated value) format, which can be read into Python in many ways, one of which is using *numpy.loadtxt()* using the appropriate delimiter:

```
from numpy import loadtxt

ecg = loadtxt("ecg.csv", delimiter=',')
```
Now we have the data in a `numpy` array `ecg`. We can quickly plot what the data looks like using `matplotlib.pyplot`.

![Original signal](image)

Fig. 1.1: Original ECG signal as plotted by `matplotlib.pyplot`.

```python
from matplotlib.pyplot import plot, show, title
title("Original signal")
plot(ecg)
show()
```

The data stored in `ecg` can be decomposed with CEEMDAN using the routine `ceemdan()`. The only thing we need to decide is what to use as the stopping criterion for the sifting iterations. In this example we use a S-number of 4 and a maximum number of siftings of 50:

```python
from pyeemd import ceemdan
imfs = ceemdan(ecg, S_number=4, num_siftings=50)
```

Now the rows of the 2D array `imfs` are the IMFs the original signal decomposes to when applying CEEMDAN. We can plot these IMFs using `matplotlib.pyplot` as before, but `pyeemd` also comes with an utility function `plot_imfs()` for easily plotting the IMFs (using `matplotlib.pyplot`) in separate figures.

```python
from pyeemd.utils import plot_imfs
plot_imfs(imfs, plot_splines=False)
show()
```

The `plot_splines=False` argument prevents the plotting of the envelope curves of the IMFs, which would otherwise be shown.

This concludes our simple tutorial. For more in-depth information about the methods available in `pyeemd` please head to the API documentation. You can also look at example code at the examples subdirectory of `pyeemd`.
How you choose to use or process the IMFs obtained by the decomposition routines is beyond the scope of this document – and beyond the scope of pyeemd – but you might be interested in the Hilbert transform routine offered by scipy.fftpack.hilbert().

## 1.3 API documentation

### 1.3.1 Main decomposition routines

**pyeemd.eemd** *(inp, num_imfs=None, ensemble_size=250, noise_strength=0.2, S_number=None, num_siftings=None, rng_seed=0)*

Decompose input data array *inp* to Intrinsic Mode Functions (IMFs) with the Ensemble Empirical Mode Decomposition algorithm [R1].

The size of the ensemble and the relative magnitude of the added noise are given by parameters *ensemble_size* and *noise_strength*, respectively. The stopping criterion for the decomposition is given by either a S-number or an absolute number of siftings. In the case that both are positive numbers, the sifting ends when either of the conditions is fulfilled. By default, *num_siftings=50* and *S_number=4*. If only *S_number* is set to a positive value, *num_siftings* defaults to 50. If only *num_siftings* is set to a positive value, *S_number* defaults to 0.

**Parameters** *inp*: array_like, shape (N,)

The input signal to decompose. Has to be a one-dimensional array-like object.

**num_imfs**: int, optional

Number of IMFs to extract. If set to *None*, a default value given by *emd_num_imfs(N)* is used. Note that the residual is also counted in this number, so *num_imfs=1* will simply give you the input signal plus noise.
ensemble_size : int, optional
    Number of copies of the input signal to use as the ensemble.

noise_strength : float, optional
    Standard deviation of the Gaussian random numbers used as additional noise. This value is relative to the standard deviation of the input signal.

S_number : int, optional
    Use the S-number stopping criterion \([R2]\) for the EMD procedure with the given values of \(S\). That is, iterate until the number of extrema and zero crossings in the signal differ at most by one, and stay the same for \(S\) consecutive iterations. Typical values are in the range 3 .. 8. If \(S_number\) is zero, this stopping criterion is ignored.

num_siftings : int, optional
    Use a maximum number of siftings as a stopping criterion. If \(num_siftings\) is zero, this stopping criterion is ignored.

rng_seed : int, optional
    A seed for the random number generator. A value of zero denotes an implementation-defined default value.

Returns imfs : ndarray, shape (M, N)
    A \(MxN\) array with \(M = num_imfs\). The rows of the array are the IMFs of the input signal, with the last row being the final residual.

See also:
emd The regular Empirical Mode Decomposition routine.
emd_num_imfs The number of IMFs returned for a given input length \(N\) unless a specific number is set by \(num_imfs\).

Notes

At least one of \(S_number\) and \(num_siftings\) must be positive. If both are positive, the iteration stops when either of the criteria is fulfilled.

References

[R1], [R2]

pyeemd.emd (inp, num_imfs=None, S_number=None, num_siftings=None)
    A convenience function for performing EMD (not EEMD). This simply calls function \(eemd()\) with \(ensemble_size=1\) and \(noise_strength=0\).

pyeemd.ceemdan (inp, num_imfs=None, ensemble_size=250, noise_strength=0.2, S_number=None, num_siftings=None, rng_seed=0)
    Decompose input data array \(inp\) to Intrinsic Mode Functions (IMFs) with the Complete Ensemble Empirical Mode Decomposition with Adaptive Noise (CEEMDAN) algorithm \([R3]\), a variant of EEMD. For description of the input parameters and output, please see documentation of \(eemd()\).

See also:
emd The regular Ensemble Empirical Mode Decomposition routine.
**emd_num_imfs** The number of IMFs returned for a given input length.

**References**

[R3]

### 1.3.2 Auxiliary routines

**pyeemd.emd_num_imfs**($N$)
Return number of IMFs that will be extracted from input data of length $N$, including the final residual.

**pyeemd.emd_find_extrema**($x$)
Find the local minima and maxima from input data $x$. This includes the artificial extrema added to the ends of the data as specified in the original EEMD article [R4].

**Parameters**
- $x$: array_like, shape (N,)
The input data. Has to be a one-dimensional array_like object.

**Returns**
- **maxx**: ndarray
  The x-coordinates of the local maxima.
- **maxy**: ndarray
  The y-coordinates of the local maxima.
- **minx**: ndarray
  The x-coordinates of the local minima.
- **miny**: ndarray
  The y-coordinates of the local minima.

**References**

[R4]

**pyeemd.emd_evaluate_spline**($x$, $y$)
Evaluates a cubic spline with the given ($x$, $y$) points as knots.

**Parameters**
- $x$: array_like, shape (N,)
The x coordinates of the knots. The array must be sorted, start from 0 and end at an integer.
- $y$: array_like, shape (N,)
The y coordinates of the knots.

**Returns**
- **spline_y**: ndarray
  The cubic spline curve defined by the knots and the “not-a-knot” end conditions, evaluated at integer points from 0 to $\max(x)$.

**See also:**

**emd_find_extrema** A method of finding the extrema for spline fitting.
Notes

As you can see from the definition, this method is tuned to work only in the case needed by EMD. This method is made available mainly for visualization and unit testing purposes. Better general purpose spline methods exist already in scipy.interpolate.

1.3.3 Utility code: pyeemd.utils

Some utility functions for visualizing IMFs produced by the (E)EMD methods.

```
utils.plot_imfs(imfs, new_figs=True, plot_splines=True)
```

Plot utility method for plotting IMFs and their envelope splines with pylab.

**Parameters**
- `imfs`: ndarray
  - The IMFs as returned by `pyeemd.emd()`, `pyeemd.eemd()`, or `pyeemd.ceemdan()`.
- `new_figs`: bool, optional
  - Whether to plot the IMFs in separate figures.
- `plot_splines`: bool, optional
  - Whether to plot the envelope spline curves as well.
CHAPTER 2

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