Proper Orthogonal Decomposition Documentation

Release 0.0.5

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ONE

ANALYSIS MODULE

```
class analysis.Timer
    Bases: object
    Allows some basic profiling
analysis.plotResults(Y, T, label=None, legend_loc='upper left', show_legend=False)
analysis.randomRuns(sys, rsys, T, sigma=10.0, integrator='dopri5')
analysis.runAnalysis(n, k, N=1, example='butter', T=20, sigma=1.0, integrator='dopri5')
analysis.x0(n)
```

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TWO

DEBUG MODULE

```
class debug.InputOfCalls (f)
    Bases: object

Save the different calls to functions in order to see what might go wrong somewhere
    classmethod inputs()
        return a dict of {function: [inputs],...}

instances = {}
```

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THREE

EXAMPLE2SYS MODULE

```
example2sys.generateRandomExample (n, m, p=None, distribution=<bound method Random.gauss of <random.Random object at 0x7fd3ec11d620>>, distributionArguments=[0.0, 1.0])

Generate a random example of arbitraty order

How to use: sys = generateRandomExample(n, m, [p, distribution])

Inputs: n system order m number of Inputs p number of outputs [p=m] distribution distribution of the matrix values [distribution=gauss(0., 1.)]

Output: sys random StateSpaceSystem with the parameters set in the input example2sys.heatSystem(N, L=1.0, g0=0.0, gN=0.0)

example2sys.optionPricing(N=None, option='put', r=0.05, T=1.0, K=100.0, L=None)

example2sys.stableRandomSystem(*args, **kwargs)
```

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FOUR

FUTURESCIPY MODULE

Some functions that will be integrated into scipy in a future version. The use of this file should be avoided as soon as the functions are fixed in scipy.

futurescipy.abcd_normalize(A=None, B=None, C=None, D=None)

Check state-space matrices and ensure they are rank-2.

If enough information on the system is provided, that is, enough properly-shaped arrays are passed to the function, the missing ones are built from this information, ensuring the correct number of rows and columns. Otherwise a ValueError is raised.

Parameters B, C, D (*A*,) – State-space matrices. All of them are None (missing) by default.

Returns A, B, C, D – Properly shaped state-space matrices.

Return type array

Raises ValueError – If not enough information on the system was provided.

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POD MODULE

Model order reduction for linear state space systems can be done with Proper Orthogonal Decomposition (POD) methods. Some of them can be simply applied when creating a system.

pod.isStable(A)

class pod.lss (*create_from, **reduction_options)

Bases: object

linear time independent state space system.

Default contstructor is called like lss(A, B, C, D) and a system can easily be copied by calling lss(sys) where sys is a lss object itself.

Parameters

- A (array_like) -
- **B** (array_like) -
- **C** (array_like) –
- **D** (*array_like*) State-Space matrices. If one of the matrices is None, it is replaced by a zero matrix with appropriate dimensions.
- **reduction** (*f'truncation_square_root'*), *optional*) Choose method of reduction. If it isn't provided, matrices are used without reduction.
- **reduction_options (dict, optional) The arguments with which the reduction method is called.

x0 array like, optional

Initial state. Defaults to the zero state.

t0 float

Initial value for t

integrator str

Name of the integrator used by scipy.integrate.ode

integrator_options dict

Options for the specified integrator that can be set.

reduction_functions dict

The functions that can be choosen as an input parameter with the *reduction* keyword.

__call__ (times, control=None, force_ode_reset=False)

Get the output at specified times with a provided control

It is possible to only request the output at one particular time or provide a list of times. If *times* is a sequence, the output will be a list of nparrays at these times, otherwise it's just a single nparray. However the control can either be specified as a function or is a constant array over all times.

Parameters

- times (list or scalar) The output for these timese will be calculated
- **control** (callable control (t, y) or array_like, optional) If it is specified, it will be overwritten in the attributes.
- **force_ode_reset** (*Boolean*, *optional*) If it's called, the ode solver is reset and the current attributes are used.

```
\mathbf{f}(t, y, u)
           Rhs of the differential equation
     integrator = 'dopri5'
     integrator_options = {}
     reduction functions = {'truncation square root': <function truncation square root at 0x107713cf8>}
     setupODE()
          Set the ode solver. All integrator, options and initial value can be set through class attributes.
     solve(t)
     t
           Current time of the system
     t0 = 0.0
           State of the system at current time t
     x0 = None
     У
                                                                          balance=True,
pod.truncation_square_root(A,
                                          В,
                                                C,
                                                     k=0,
                                                              tol=0.0,
                                                                                            scale=True,
```

check_stability=True, length_cache_array=None)

Perform truncation of a system. Scaling and balancing are optional

This allows to reduce a linear state space system by either specifying it to a certain number of states k or by specifying a error tolerance tol relative to the input. In theory the most accurate results are achieved by using balance and scale but the size of the error strongly depends on the particular problem and scaling and balancing may in some cases cost too much.

Parameters

- A (array like) –
- **B** (array_like) –
- C (array_like) State-Space matrices of the system that should be reduced
- k (int, optional) Order of the output system
- tol (float, optional) Error of the output system, based on the Hankel Singular Values
- balance (Boolean, optional) Balance the system before reducing it to make sure, that the
 error is kept small
- scale (Boolean, optional) Scale the system

• **check_stability** (*Boolean*, *optional*) – Checks if all the real parts of the eigenvalues of A are in the left half of the complex plane.

Returns

- Nr (int) Actual size of the system, based on the error: If the Machine error would have been bigger than the error of the reduced system, it may happen that Nr < k and if the error would be inconsiderably bad, It might be the case that Nr > k. In case k was never specified, this is purely based on tol
- Ar, Br, Cr (ndarray) Reduced arrays
- **hsv** (*ndarray*) Hankel singular values of the original system. The size of the error may be calculated based on this.

Raises

- ValueError If the system that's provided is not stable (i.e. *A* has eigenvalues which have non-negative real parts)
- ImportError If the slycot subroutine *ab09ad* can't be found. Occurs if the slycot package is not installed.

SIX

TESTS MODULE

Unit tests and functional tests for the pod.py package

```
class tests.testLss (methodName='runTest')
    Bases: unittest.case.TestCase
    test lss functionalities
    testIdentity()
    test_abcd_normalize()
    test_f()
    test_zero_control()
```

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