serpentTools Documentation

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ONE

SERPENT-TOOLS

A suite of parsers designed to make interacting with SERPENT¹ output files simple and flawless.

The SERPENT Monte Carlo code is developed by VTT Technical Research Centre of Finland, Ltd. More information, including distribution and licensing of SERPENT can be found at montecarlo.vtt.fi

1.1 Issues

If you have issues installing the project, find a bug, or want to add a feature, the GitHub issue page is the best place to do that.

1.2 Contributors

Here are all the wonderful people that helped make this project happen

- · Andrew Johnson
- · Dr. Dan Kotlyar
- Stefano Terlizzi

1.3 References

The Annals of Nuclear Energy article should be cited for all work using SERPENT. If you wish to cite this project, please cite as:

¹ Leppanen, J. et al. (2015) "The Serpent Monte Carlo code: Status, development and applications in 2013." Ann. Nucl. Energy, 82 (2015) 142-150

1.4 License

MIT License

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TWO

CONTRIBUTING

2.1 Contributing

First, thanks for your interest in contributing to this project! This document should help us expedite the process of reviewing issues and pull requests. For a quick look at all the issues that are up for grabs, take a look at the current unclaimed issues. If you claim an issue, use the Assignees setting to let us know that you've got it!

2.1.1 Scope

The scope of this project is to simplify and expedite analysis stemming from SERPENT outputs. In the future we may expand this project to expand to interacting heavily with input files, but that is currently beyond the scope of this project. Any and all issues, features and pull requests will be examined through this scope.

2.1.2 Issues

The goal for this project is to become the de facto method for processing SERPENT outputs and, if you're looking at this, there is some way we can improve. The GitHub issue tracker is the preferred way to post bug reports and feature requests.

Bug Reports

The more information given, the quicker we can reproduce and hopefully resolve the issue. Please see *Issue Template* for a template that should be used for reporting bugs. One of the developers will add a bug label and start moving to resolve the issue posthaste.

Feature Requests

We are very interested in adding functionality from the SERPENT community! Requests can be done through the issue tracker as well. You can create an issue on the issue tracker with [Feature] or [Request] in the title. Describe what you would like to add, some expected results, and the purpose behind the feature. The development team will apply an enhancement label and proceed accordingly.

2.1.3 Pull Requests

Pull requests are how we review, approve, and incorporate changes into the develop and master branches. If you have code you want to contribute, please look at the content in the *Developer's Guide* for things like *Pull Request Checklist*, *Coding Style*, and more.

When your content is ready for the pull request, follow the *Pull Request Template* and make a request! Someone of the core development team will review the changes according to the criteria above and make changes and/or approve for merging!

2.2 Issue Template

Summary of issue

Code for reproducing the issue

Actual outcome including console output and error traceback if applicable

Expected outcome - why this is an issue

- Version from serpentTools.__version__
- Python version python --version
- IPython or Jupyter version if applicable

2.3 Pull Request Template

Title: Short descriptive title summarizing the request

Body: Directly address the issue that this pull request resolves or addresses:

Fixes #99 - some fatal bug that caused things to fail poorly

Include a full description of what is included in this PR, what has changed.

THREE

EXAMPLES

3.1 User Control

The serpentTools package is designed to, without intervention, be able to store all the data contained in each of the various output files. However, the serpentTools.settings module grants great flexibility to the user over what data is obtained through the rc class. This notebook will provide as an intro into using this class.

3.1.1 Basic Usage

```
>>> import serpentTools
>>> from serpentTools.settings import rc, defaultSettings
INFO : serpentTools: Using version 1.0b0+24.g23e6eac.dirty
```

Below are the default values for each setting available

```
>>> for setting in sorted(defaultSettings.keys()):
       print(setting)
>>>
        for key in defaultSettings[setting]:
            print('\t', key, '-', defaultSettings[setting][key])
>>>
depletion.materialVariables
     default - []
     description - Names of variables to store. Empty list -> all variables.
     type - <class 'list'>
depletion.materials
     default - []
     description - Names of materials to store. Empty list -> all materials.
     type - <class 'list'>
depletion.metadataKeys
    default - ['ZAI', 'NAMES', 'DAYS', 'BU']
     description - Non-material data to store, i.e. zai, isotope names, burnup
\rightarrowschedule, etc.
     type - <class 'list'>
     options - default
depletion.processTotal
     default - True
     description - Option to store the depletion data from the TOT block
     type - <class 'bool'>
serpentVersion
     default - 2.1.29
     description - Version of SERPENT
     type - <class 'str'>
     options - ['2.1.29']
```

```
verbosity
   default - warning
   type - <class 'str'>
   description - Set the level of errors to be shown.
   updater - <function updateLevel at 0x000001B7F3DD6598>
   options - ['critical', 'error', 'warning', 'info', 'debug']

xs.variableExtras
   default - []
   description - Full SERPENT name of variables to be read
   type - <class 'list'>

xs.variableGroups
   default - []
   description - Name of variable groups from variables.yaml to be expanded into

SERPENT variable to be stored
   type - <class 'list'>
```

Settings such as depletion.materialVariables are specific for the DepletionReader, while settings that are led with xs are sent to the ResultsReader and BranchingReader, as well as their specific settings. The rc class acts as a dictionary, and updating a value is as simple as

```
>> rc['verbosity'] = 'debug'
DEBUG : serpentTools: Updated setting verbosity to debug
```

The rc object automatically checks to make sure the value is of the correct type, and is an allowable option, if given.

```
>>> try:
>>> rc['depletion.metadataKeys'] = False
>>> except TypeError as te:
>>> print(te)
Setting depletion.metadataKeys should be of type <class 'list'>, not <class 'bool'>
>>> try:
>>> rc['serpentVersion'] = '1.2.3'
>>> except KeyError as ke:
>>> print(ke)
"Setting serpentVersion is
1.2.3
and not one of the allowed options:
['2.1.29']"
```

The rc object can also be used inside a context manager to revert changes.

```
>>> with rc:
>>> rc['depletion.metadataKeys'] = ['ZAI', 'BU']
>>>
>>> rc['depletion.metadataKeys']
DEBUG : serpentTools: Updated setting depletion.metadataKeys to ['ZAI', 'BU']
DEBUG : serpentTools: Updated setting depletion.metadataKeys to ['ZAI', 'NAMES', 'DAYS', 'BU']
['ZAI', 'NAMES', 'DAYS', 'BU']
```

3.2 DepletionReader

3.2.1 Basic Operation

SERPENT produces a burned material file, containing the evolution of material properties through burnup for all burned materials present in the problem. The DepletionReader is capable of reading this file, and storing the data inside <code>DepletedMaterial</code> objects. Each such object has methods and attributes that should ease analysis.

```
>>> import six
>>> import serpentTools
>>> from serpentTools.settings import rc
INFO : serpentTools: Using version 1.0b0+24.g23e6eac.dirty
>>> depFile = 'demo_dep.m'
>>> dep = serpentTools.read(depFile)
INFO : serpentTools: Inferred reader for demo_dep.m: DepletionReader
INFO : serpentTools: Preparing to read demo_dep.m
INFO : serpentTools: Done reading depletion file
```

The materials read in from the file are stored in the materials dictionary, where the keys represent the name of specific materials, and the corresponding values are the depleted material.

```
>>> dep.materials
{'bglass0': <serpentTools.objects.materials.DepletedMaterial at 0x23905154668>,
  'fuel0': <serpentTools.objects.materials.DepletedMaterial at 0x2390578eeb8>,
  'total': <serpentTools.objects.materials.DepletedMaterial at 0x2390579e978>}
```

Metadata, such as the isotopic vector and depletion schedule are also present inside the reader

3.2.2 DepletedMaterial

As mentioned before, all the material data is stored inside these <code>DepletedMaterial</code> objects. These objects share access to the metadata of the reader as well.

```
>>> fuel = dep.materials['fuel0']
>>> fuel.burnup
array([ 0. , 0.02, 0.04, ..., 1.36, 1.38, 1.4 , 1.42])
>>> fuel.days is dep.metadata['days']
True
```

All of the variables present in the depletion file for this material are present, stored in the data dictionary. A few properties commonly used are accessible as attributes as well.

```
[ 0.00000000e+00, 6.06880000e-09, 8.11783000e-09, ..., 8.05991000e-09, 8.96359000e-09, 9.28554000e-09], [ 4.48538000e-06, 4.48486000e-06, 4.48432000e-06, ..., 4.44726000e-06, 4.44668000e-06, 4.44611000e-06], ..., [ 0.00000000e+00, 3.03589000e-11, 7.38022000e-11, ..., 1.62829000e-09, 1.63566000e-09, 1.64477000e-09], [ 0.00000000e+00, 1.15541000e-14, 2.38378000e-14, ..., 8.60736000e-13, 8.73669000e-13, 8.86782000e-13], [ 6.88332000e-02, 6.88334000e-02, 6.88336000e-02, ..., 6.88455000e-02, 6.88457000e-02, 6.88459000e-02]])
```

Similar to the original file, the rows of the matrix correspond to positions in the isotopic vector, and the columns correspond to positions in burnup/day vectors.

```
>>> fuel.mdens.shape # rows, columns
(34, 72)
>>> fuel.burnup.shape
(72,)
>>> len(fuel.names)
34
```

3.2.3 Data Retrieval

At the heart of the <code>DepletedMaterial</code> is the <code>getValues()</code> method. This method acts as an slicing mechanism that returns data for a select number of isotopes at select points in time.

```
>>> dayPoints = [0, 5, 10, 30]

>>> iso = ['Xe135', 'U235']

>>> vals = fuel.getValues('days', 'a', dayPoints, iso)

>>> print(vals.shape)

(2, 4)

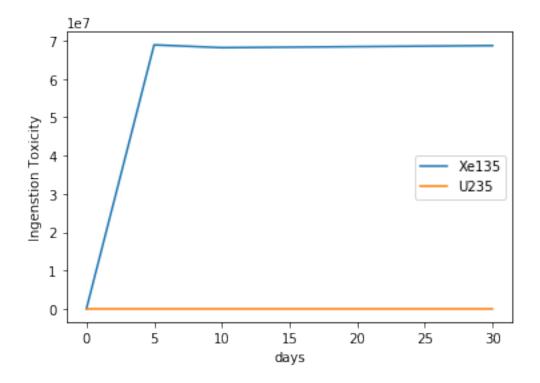
>>> print(vals)

[[ 0.00000000e+00     3.28067000e+14     3.24606000e+14     3.27144000e+14]

[ 5.36447000e+07     5.34519000e+07     5.32499000e+07     5.24766000e+07]]
```

The DepletedMaterial uses this slicing for the built-in plot () method

```
>>> fuel.plot('days', 'ingTox', dayPoints, iso, ylabel='Ingenstion Toxicity');
```



3.2.4 Limitations

Currently, the <code>DepletionReader</code> cannot catch materials with underscore in the name, due to variables like <code>ING_TOX</code> also containing an underscore. Issue #58

3.2.5 Settings

The DepletionReader also has a collection of settings to control what data is stored. If none of these settings are modified, the default is to store all the data from the output file.

```
>>> from serpentTools.settings import rc, defaultSettings
>>> for setting in defaultSettings:
        if 'depletion' in setting:
>>>
>>>
            print(setting)
>>>
            for k, v in six.iteritems(defaultSettings[setting]):
                print('\t', k, v)
depletion.materials
     type <class 'list'>
     description Names of materials to store. Empty list -> all materials.
     default []
depletion.processTotal
    type <class 'bool'>
    description Option to store the depletion data from the TOT block
    default True
depletion.materialVariables
   type <class 'list'>
    description Names of variables to store. Empty list -> all variables.
   default []
depletion.metadataKeys
    type <class 'list'>
```

```
description Non-material data to store, i.e. zai, isotope names, burnup schedule, →etc.
options default
default ['ZAI', 'NAMES', 'DAYS', 'BU']
```

Below is an example of configuring a DepletionReader that only stores the burnup days, and atomic density for all materials that begin with bqlass followed by at least one integer.

Note: Creating the DepletionReader in this manner is functionally equivalent to serpentTools. read(depFile)

```
>>> rc['depletion.processTotal'] = False
>>> rc['depletion.metadataKeys'] = ['BU']
>>> rc['depletion.materialVariables'] = ['ADENS']
>>> rc['depletion.materials'] = [r'bglass\d+']
>>>
>>> bgReader = serpentTools.parsers.DepletionReader(depFile)
>>> bgReader.read()
       : serpentTools: Preparing to read demo_dep.m
       : serpentTools: Done reading depletion file
>>> bgReader.materials
{'bglass0': <serpentTools.objects.materials.DepletedMaterial at 0x239057dcb00>}
>>> bglass = bgReader.materials['bglass0']
>>> bglass.data
                      , 0. , 0. , ..., 0.
{'adens': array([[ 0.
                                                                    , 0.
         0.
             ],
        [ 0.
                  , 0.
                              , 0.
                                          , . . . ,
                                                  0.
                                                              0.
         0.
                  ],
                  , 0.
        [ 0.
                              , 0.
                                          , . . . ,
         0.
                  ],
        . . . ,
        0.
                              , 0.
                                          , ..., 0.
                     0.
                                                              0.
         0.
                  ],
        [ 0.
                              , 0.
                     0.
                                          , ..., 0.
                                                              0.
         0.
        [0.0715841, 0.0715843, 0.0715845, ..., 0.0715968, 0.0715969,
         0.0715971]])}
>>> bglass.data.keys()
dict_keys(['adens'])
```

3.2.6 Conclusion

The DepletionReader is capable of reading and storing all the data from the SERPENT burned materials file. Upon reading, the reader creates custom DepletedMaterial objects that are responsible for storing and retrieving the data. These objects also have a handy plot method for quick analysis. Use of the rc settings control object allows increased control over the data selected from the output file.

3.2.7 References

1. J. Leppänen, M. Pusa, T. Viitanen, V. Valtavirta, and T. Kaltiaisenaho. "The Serpent Monte Carlo code: Status, development and applications in 2013." Ann. Nucl. Energy, 82 (2015) 142-150

FOUR

DEVELOPER'S GUIDE

4.1 Coding Style

For the most part, this project follows the PEP 8 standard with a few differences. Some points are included here

- 79 characters per line
- Four spaces per indentation level
- Avoiding extraneous whitespace:

```
Yes: spam(ham[1], {eggs: 2})
No: spam( ham[ 1 ], { eggs: 2 } )
```

Some of the specific style points for this project are included below

- mixedCase for variables, methods, and functions
- CamelCase for classes:

```
class DemoClass(object):
    def doSomething(self, arg0, longerArgumentName):
        pass
```

• Directly call the __init__ method from a parent class, e.g.:

```
class MyQueue(list):

    def __init__(self, items):
        list.__init__(self)
        self.extend(items)
```

- Arrange imports in the following order:
 - 1. imports from the standard library: os, sys, etc.
 - 2. imports from third party code: numpy, matplotlib, etc.
 - 3. imports from the serpentTools package
- Longer import paths are preferred to shorter:

```
# yes
from really.long.path.to.a import function
function()
# not preferred
```

```
import really
really.long.path.to.a.function()
```

4.2 Documentation

All public functions, methods, and classes should have adequate documentation through docstrings, examples, or inclusion in the appropriate file in the docs directory. Good forethought into documenting the code helps resolve issues and, in the case of docstrings, helps produce a full manual.

4.2.1 Docstrings

Docstrings are defined in PEP 257 and are characterized by """triple double quotes"". These can be used to reduce the effort in creating a full manual, and can be viewed through python consoles to give the user insight into what is done, what is required, and what is returned from a particular object. This project uses numpy style docstrings, examples of which are given below

Functions and Methods

Below is the depmtx() function annotated using short and longer docstrings:

```
def depmtx(filePath):
    """Return t, no, zai, a, and n1 values from the depmtx file."""
```

or:

```
def depmtx(filePath):
   Read the contents of the ``depmtx`` file and return contents
    .. note::
       If ``scipy`` is not installed, matrix ``A`` will be full.
       This can cause some warnings or errors if sparse or
       non-sparse solvers are used.
   Parameters
   fileP: str
       Path to depletion matrix file
   Returns
   t: float
       Length of time
   n0: numpy.ndarray
       Initial isotopic vector
   zai: numpy.array
       String identifiers for each isotope in ``n0`` and ``n1``
   a: numpy.array or scipy.sparse.csc_matrix
       Decay matrix. Will be sparse if scipy is installed
   n1: numpy.array
       Final isotopic vector
```

Both docstrings indicate what the function does, and what is returned. The latter, while more verbose, is preferred in most cases, for the following reasons. First, far more information is yielded to the reader. Second, the types of the inputs and outputs are given and clear. Some IDEs can obtain the expected types from the docstrings and inform the user if they are using an incorrect type.

More content can be added to the docstring, including

- · Raises Errors/warnings raised by this object
- See Also follow up information that may be useful
- Yields If this object is a generator. Similar to Returns

Classes

Classes can have a more lengthy docstring, as they often include attributes to which the class has access. Below is an example of the docstring for the <code>DepletionReader</code>:

```
Parser responsible for reading and working with depletion files.

Parameters
------
filePath: str
    path to the depletion file

Attributes
------
materials: dict
    Dictionary with material names as keys and the corresponding
    :py:class:`~serpentTools.objects.DepletedMaterial` class
    for that material as values

metadata: dict
    Dictionary with file-wide data names as keys and the
    corresponding dataas values, e.g. 'zai': [list of zai numbers]

settings: dict
    names and values of the settings used to control operations
    of this reader
```

Class docstrings can be added to the class signature, or to the __init__ method, as:

```
class Demo(object):
    """
    Demonstration class

Parameters
------
x: str
    Just a string

Attributes
------
capX: str
    Capitalized x
"""
```

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or:

```
def __init__(self, x):
    """

    Demonstration class

    Parameters
    ------
    x: str
        Just a string

Attributes
    ------
    capX: str
        Capitalized x
    """
```

Deprecation

If an object is deprecated or will be modified in future versions, then the deprecated() and willChange() decorators should be applied to the object, and a note should be added to the docstring indicating as much.

4.2.2 Examples

When possible, features should be demonstrated, either through Jupyter notebooks in the examples/ directory, or with an Examples section in the docstring. Specifically, all readers should be demonstrated as Jupyter notebooks that detail the typical usage, user control settings, and examples of how the data is stored and accessed.

4.3 Data Model

This project has two key offerings to the SERPENT community.

- 1. Scripting tools that are capable of interpreting the output files
- 2. Purpose-built containers for storing, accessing, and utilizing the data in the output files

4.3.1 Readers

The readers are the core of this package, and should have a logical and consistent method for structuring data. Each reader contained in this project should:

- 1. Be capable of reading the intended output file
- 2. Yield to the user the data in a logical manner
- 3. Default to storing all the data without user input
- 4. Receive input from the user regarding what data to restrict/include
- 5. Report activity at varying levels, including failure, to the user

The third and fourth points refer to use of the serpentTools.settings module to receive user input. The final point refers to use of the logging and messaging functions included in <code>serpentTools.messages</code>.

For example, the <code>DepletionReader</code> stores data in the following manner:

```
Reader
| - filePath
| - fileMetadata
| - isotope names
| - isotope ZAIs
| - depletion schedule in days and burnup
| - materials
# for each material:
| - DepletedMaterial-n
| - name
| - atomic density
| - mass density
| - volume
| etc
```

Here, the reader primarily creates and stores material objects, and the useful data is stored on these objects. Some files, like the fission matrix and depletion matrix files may not have a structure that naturally favors this object-oriented approach.

4.3.2 Containers

The readers are primarily responsible for creating and populating containers, that are responsible for storing and retrieving data. Just like the <code>DepletedMaterial</code> has shortcuts for analysis like <code>getValues()</code> and <code>plot()</code>, each such container should easily and naturally provide access to the data, and also perform some common analysis on the data.

4.4 Pull Request Checklist

Below is the criteria that will be used in the process of reviewing pull requests (PR):

- 1. The content of the PR fits within the scope of the project Scope
- 2. The code included in the PR is written in good pythonic fashion, and follows the style of the project *Coding Style*
- 3. The code directly resolves a previously raised issue Issues
- 4. PR does not cause unit tests and builds to fail
- 5. Changes are reflected in documentation Documentation

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FIVE

API

5.1 Settings

```
class serpentTools.settings.UserSettingsLoader
      Class that stores the active user settings.
      clear() \rightarrow None. Remove all items from D.
      copy () \rightarrow a shallow copy of D
      expandVariables()
           Extend the keyword groups into lists of serpent variables.
                Returns Names of all variables to be scraped
               Return type set
      fromkeys()
           Returns a new dict with keys from iterable and values equal to value.
      get (k[,d]) \rightarrow D[k] if k in D, else d. d defaults to None.
      getReaderSettings (settingsPreffix)
           Get all module-wide and reader-specific settings.
               Parameters settingsPreffix (str or list) - Name of the specific reader. Will look
                    for settings that lead with readerName, e.g. depletion.metadataKeys or xs.
                    variables
               Returns Single level dictionary with settingName: settingValue pairs
               Return type dict
               Raises KeyError – If the reader name is not located in the readers settings dictionary
      items () \rightarrow a set-like object providing a view on D's items
      keys () \rightarrow a set-like object providing a view on D's keys
      pop(k|, d|) \rightarrow v, remove specified key and return the corresponding value.
           If key is not found, d is returned if given, otherwise KeyError is raised
      popitem () \rightarrow (k, v), remove and return some (key, value) pair as a
           2-tuple; but raise KeyError if D is empty.
      setValue (name, value)
           Set the value of a specific setting.
               Parameters
```

• name (str) - Full name of the setting

```
• value (value to be set) -
```

Raises

- KeyError If the value is not one of the allowable options or if the setting does not match an existing setting
- TypeError If the value is not of the correct type

```
setdefault (k[,d]) \to D.get(k,d), also set D[k]=d if k not in D 

update ([E], **F) \to None. Update D from dict/iterable E and F.

If E is present and has a .keys() method, then does: for k in E: D[k] = E[k] If E is present and lacks a .keys() method, then does: for k, v in E: D[k] = v In either case, this is followed by: for k in F: D[k] = F[k]

values () \to an object providing a view on D's values
```

5.2 Messages

System-wide methods for producing status update and errors.

See also:

- https://docs.python.org/2/library/logging.html
- https://www.python.org/dev/peps/pep-0391/

```
exception serpentTools.messages.SerpentToolsException
     Base-class for all exceptions in this project
     with_traceback()
         Exception.with_traceback(tb) - set self.__traceback__ to tb and return self.
serpentTools.messages.critical(message)
     Log that something has gone horribly wrong.
serpentTools.messages.debug(message)
     Log a debug message.
serpentTools.messages.deprecated(useInstead)
     Decorator that warns that different function should be used instead.
serpentTools.messages.error(message)
     Log that something caused an exception but was supressed.
serpentTools.messages.info(message)
     Log an info message, e.g. status update.
serpentTools.messages.updateLevel(level)
     Set the level of the logger.
serpentTools.messages.warning(message)
     Log a warning that something that could go wrong or should be avoided.
serpentTools.messages.willChange(changeMsg)
```

Decorator that warns that some functionality may change.

18 Chapter 5. API

5.3 Parser Module

The main module associated with reading SERPENT files. For information on each individual reader, please see the dedicated file listed below

Reader	Hyperlink
Branching	Branching Reader
Bumat	Bumat Reader
Depletion	Depletion Reader
Detector	Detector Reader
FissionMatrix	Fission Matrix Reader
Results	Results Reader

serpentTools.parsers.inferReader(filePath)

Attempt to infer the correct reader type.

Parameters filePath (str) – File to be read.

Raises SerpentToolsException - If a reader cannot be inferred

serpentTools.parsers.read(filePath, reader='infer')

Simple entry point to read a file and obtain the processed reader.

Note: If you know the type of reader you will be using, it is best to either pass in the string argument, or directly use the appropriate reader class

Parameters

- **filePath** (str) Path to the file to be reader
- **reader** (*str* or *callable*) Type of reader to use. If a string is given, then the actions described below will happen. If callable, then that function will be used with the file path as the first argument.

String argument	Action	
infer	Infer the correct reader based on the file	
dep	DepletionReader	
branch	BranchingReader	
det	DetectorReader	
results	ResultsReader	
bumat	BumatReader	
fission	FissionMatrixReader	

Returns Correct subclass corresponding to the file type

Return type serpentTools.objects.readers.BaseReader

Raises

- AttributeError If the object created by the reader through reader (filePath) does not have a read method.
- SerpentToolsException If the reader could not be inferred or if the requested reader string is not supported
- NotImplementedError This has the ability to load in readers that may not be complete, and thus the read method may raise this error.

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```
serpentTools.parsers.depmtx(fileP)
```

Read the contents of the depmtx file and return contents

Note: If scipy is not installed, matrix A will be full. This can cause some warnings or errors if sparse or non-sparse solvers are used.

Parameters fileP (str) – Path to depletion matrix file

Returns

- **t** (*float*) Length of time
- **n0** (*numpy.ndarray*) Initial isotopic vector
- zai (numpy.array) String identifiers for each isotope in n0 and n1
- a (numpy.array or scipy.sparse.csc_matrix) Decay matrix. Will be sparse if scipy is installed
- n1 (numpy.array) Final isotopic vector

5.4 Containers

Many of the readers utilize custom built objects for storing data in a logical and accessible format. These containers are detailed here.

Note: Variables taken from SERPENT will be converted to mixedCase and stored under those names. for example, INF_TOT would be stored as infTot

5.4.1 Homogenized Universe

class serpentTools.objects.containers.**HomogUniv** (*container*, *name*, *bu*, *step*, *day*)

Class for storing homogenized universe specifications and retrieving them

Parameters

```
• container (str) – name of the parser
```

- name (str) name of the universe
- **bu** (float) burnup value
- **step** (float) temporal step
- day (float) depletion day

name

str - name of the universe

bu

float - burnup value

step

float - temporal step

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day

float - depletion day

addData (variableName, variableValue, uncertainty=False)

Sets the value of the variable and, optionally, the associate s.d.

Warning: This method will overwrite data for variables that already exist

Parameters

- variableName (str) Variable Name
- variableValue Variable Value
- uncertainty (bool) Set to True in order to retrieve the uncertainty associated to the expected values

Raises TypeError - If the uncertainty flag is not boolean

get (variableName, uncertainty=False)

Gets the value of the variable VariableName from the dictionaries

Parameters

- variableName (str) Variable Name
- uncertainty (bool) Boolean Variable- set to True in order to retrieve the uncertainty associated to the expected values

Returns

- x Variable Value
- dx Associated uncertainty

Raises

- TypeError If the uncertainty flag is not boolean
- KeyError If the variable requested is not stored on the object

5.4.2 Material Objects

```
class serpentTools.objects.materials.DepletedMaterial (parser, name)
    Class for storing material data from _dep.m files.
```

Parameters

- parser (DepletionReader) Parser that found this material. Used to obtain file metadata like isotope names and burnup
- name (str) Name of this material

zai

numpy.array or None - Isotope id's

names

numpy.array or None - Names of isotopes

days

numpy.array or None - Days overwhich the material was depleted

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adens

numpy.array or None – Atomic density over time for each nuclide

mdens

numpy.array or None – Mass density over time for each nuclide

burnup

numpy.array or None – Burnup of the material over time

addData(variable, rawData)

Add data straight from the file onto a variable.

Parameters

- **variable** (str) Name of the variable directly from SERPENT
- rawData (list) List of strings corresponding to the raw data from the file

getValues (xUnits, yUnits, timePoints=None, names=None)

Return x values for given time, and corresponding isotope values.

Parameters

- xUnits (str) name of x value to obtain, e.g. 'days', 'burnup'
- yUnits (str) name of y value to return, e.g. 'adens', 'burnup'
- timePoints (list or None) If given, select the time points according to those specified here. Otherwise, select all points
- names (list or None) If given, return y values corresponding to these isotope names. Otherwise, return values for all isotopes.

Returns Array of values.

Return type numpy.array

Raises

- AttributeError If the names of the isotopes have not been obtained and specific isotopes have been requested
- KeyError If at least one of the days requested is not present

getXY (xUnits, yUnits, timePoints=None, names=None)

Return x values for given time, and corresponding isotope values.

Deprecated since version 0.1.0: Use getValues() instead.

plot (xUnits, yUnits, timePoints=None, names=None, ax=None, legend=True, label=True, xlabel=None, ylabel=None)

Plot some data as a function of time for some or all isotopes.

Parameters

- xUnits (str) name of x value to obtain, e.g. 'days', 'burnup'
- yUnits (str) name of y value to return, e.g. 'adens', 'burnup'
- timePoints (list or None) If given, select the time points according to those specified here. Otherwise, select all points
- names (list or None) If given, return y values corresponding to these isotope names. Otherwise, return values for all isotopes.
- **ax** (None or matplotlib axes) If given, add the data to this plot. Otherwise, create a new plot

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- legend (bool) Automatically add the legend
- label (bool) Automatically label the axis
- xlabel (None or str) If given, use this as the label for the x-axis. Otherwise, use xUnits
- ylabel (None or str) If given, use this as the label for the y-axis. Otherwise, use yUnits

Returns Axes corresponding to the figure that was plotted

Return type matplotlib axes

See also:

getValues()

5.5 Branching Reader

Warning: This reader is not implemented yet. This merely serves as a placeholder

class serpentTools.parsers.BranchingReader(filePath)

Parser responsible for reading and working with automated branching files.

Parameters filePath (str) – path to the depletion file

write (template=None)

Write the data from the branching sequence using a specified template.

Parameters template (str) – Will attempt to make a template in the following order:

- 1. By reading the template from the file pointed to by template, or
- 2. Making a template directly from the string

If neither of these succeeds, the method will fail and raise an exception.

5.6 Bumat Reader

Warning: This reader is not implemented yet. This merely serves as a placeholder

class serpentTools.parsers.bumat.BumatReader(filePath, readerSettingsLevel)

Parser responsible for reading and working with burned material files.

Parameters filePath (str) – path to the depletion file

5.7 Depletion Reader

Warning: Does not support depleted materials with underscores, i.e. fuel_1 will not be matched with the current methods

class serpentTools.parsers.DepletionReader(filePath)

Parser responsible for reading and working with depletion files.

Parameters filePath (str) – path to the depletion file

materials

dict – Dictionary with material names as keys and the corresponding DepletedMaterial class for that material as values

metadata

dict – Dictionary with file-wide data names as keys and the corresponding dataas values, e.g. 'zai': [list of zai numbers]

settings

dict - names and values of the settings used to control operations of this reader

read()

Read through the depletion file and store requested data.

5.8 Detector Reader

Warning: This reader is not implemented yet. This merely serves as a placeholder

class serpentTools.parsers.detector.DetectorReader (filePath, readerSettingsLevel)
 Parser responsible for reading and working with detector files.

Parameters filePath (str) – path to the depletion file

read()

Read the file and store the data.

Warning This read function has not been implemented yet

5.9 Fission Matrix Reader

Warning: This reader is not implemented yet. This merely serves as a placeholder

Parser responsible for reading and working with fission matrix files.

Parameters filePath (str) – path to the depletion file

read()

Read the file and store the data.

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Warning This read function has not been implemented yet

5.10 Results Reader

Warning: This reader is not implemented yet. This merely serves as a placeholder

class serpentTools.parsers.results.ResultsReader(filePath, readerSettingsLevel)
 Parser responsible for reading and working with result files.

Parameters filePath (str) – path to the depletion file

read()

Read the file and store the data.

Warning This read function has not been implemented yet

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