SIMPA

version 0.2.0

CAMI (Computer Assisted Medical Interventions), DKFZ, Heidelberg

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Welcome to the SIMPA documentation!



Simulation and Image Processing for Photoacoustic Imaging

README

The Simulation and Image Processing for Photoacoustic Imaging (SIMPA) toolkit.

SIMPA Install Instructions

These install instructions are made under the assumption that you have access to the phabricator simpa project. When you are reading these instructions there is a 99% chance that is the case (or someone send these instructions to you).

So, for the 1% of you: Please also follow steps 1 - 3:

- 1. git clone https://phabricator.mitk.org/source/simpa.git
- 2. git checkout master
- 3. git pull

Now open a python instance in the 'simpa' folder that you have just downloaded. Make sure that you have your preferred virtual environment activated

- 1. cd simpa
- 2. python -m setup.py build install
- 3. Test if the installation worked by using python followed by import simpa then exit()

If no error messages arise, you are now setup to use simpa in your project.

Building the documentation

When the installation went fine and you want to make sure that you have the latest documentation you should do the following steps in a command line:

1. Navigate to the simpa source directory (same level where the setup.py is in)

2. Execute the command sphinx-build -b pdf -a simpa_documentation/src simpa_documentation

3. Find the PDF file in simpa_documentation/simpa_documantation.pdf

External Tools installation instructions

mcx (Optical Forward Model)

Either download suitable executables or build yourself from the following sources: http://mcx.space/

k-Wave (Acoustic Forward Model)

Please follow the following steps and use the k-Wave install instructions for further (and much better) guidance under http://www.k-wave.org/!

- 1. Install MATLAB with the core and parallel computing toolboxes activated at the minimum.
- 2. Download the kWave toolbox
- 3. Add the kWave toolbox base bath to the toolbox paths in MATLAB
- 4. If wanted: Download the CPP and CUDA binary files and place them inthe k-Wave/binaries folder
- 5. Note down the system path to the matlab executable file.

On MATLAB r2020a or newer there is a bug when using the GPU binaries with kWave. Please follow these instructions http://www.k-wave.org/forum/topic/error-reading-h5-files-when-using-binaries to fix this bug.

MITK

Overview

The main use case for the simpa framework is the simulation of photoacoustic images. However, it can also be used for image processing.

Simulating photoacoustic images

A basic example on how to use simpa in you project to run an optical forward simulation is given in the samples/minimal_optical_simulation.py file.

Performance profiling

Do you wish to know which parts of the simulation pipeline cost the most amount of time? If that is the case then you can use the following commands to profile the execution of your simulation script. You simply need to replace the myscript name with your script name.

```
python -m cProfile -o myscript.cprof myscript.py
pyprof2calltree -k -i myscript.cprof
```

Developer Guide

Dear SIMPA developers, Dear person who wants to contribute to the SIMPA toolkit,

First of all: Thank you for your participation and help! It is much appreciated! This Guide is meant to be used as a collection of How-To's to contribute to the framework. In case you have any questions, do not hesitate to get in touch with the members of the core development team:

Kris K. Dreher (k.dreher@dkfz-heidelberg.de)

Janek M. Groehl (janek.grohl@cruk.cam.ac.uk)

How to contribute

The SIMPA code is written and maintained on a closed repository that is hosted on a server of the German Cancer Research Center. The current master branch of the repository is open source and mirrored on github.

To contribute to SIMPA, please fork the SIMPA github repository and create a pull request with a branch containing your suggested changes. The core team developers will then review the suggested changes and integrate these into the code base.

Please see the github guidelines for creating pull requests: https://docs.github.com/en/github/collaborating-with-issues-and-pull-requests/about-pull-requests

Coding style

When writing code for SIMPA, please use the PEP 8 python coding conventions (https://www.python.org/dev/peps/pep-0008/) and consider to use the following structures in your code in order to make a new developer or someone external always know exactly what to expect.

- Classnames are written in camel-case notation ClassName
- Function names are written in small letter with _ as the delimiter function_name
- Function parameters are always annotated with their type arg1: type = default
- Only use primitive types as defaults. If a non-primitive type is used, then the default should be None and the parameter should be initialized in the beginning of a function.
- A single line of code should not be longer than 120 characters.
- Functions should follow the following simple structure:
 - Input validation (arguments all not None, correct type, and acceptable value ranges?)
 - 2. Processing (clean handling of errors that might occur)
 - 3. Output generation (sanity checking of the output before handing it off to the caller)

Documenting your code

Only documented code will appear in the sphinx generated documentation.

A class should be documented using the following syntax:

```
class ClassName(Superclass):
    """
    Explain how the class is used and what it does.
    """
```

For functions, a lot of extra attributes can be added to the documentation:

```
def function_name(self, arg1:type = default, arg2:type = default) -> return_type:
    """
    Explain how the function is used and what it does.

:param arg1: type, value range, Null acceptable?
:param arg2: type, value range, Null acceptable?
:returns: type, value range, does it return Null?
:raises ExceptionType: explain when and why this exception is raised
    """
```

Adding literature absorption spectra

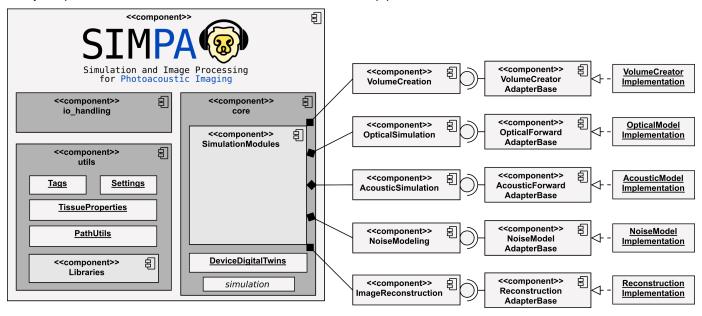
handled central point, where absorption spectra are collected and in The simpa.utils.libraries.spectra_library.py. file comprises the class AbsorptionSpectrumLibrary, in which the new absorption spectra can be added using the following two steps:

- 1. In the beginning of the class, there is a bunch of constants that define spectra using the AbsorptionSpectrum class. Add a new constant here: NEW_SPECTRUM = AbsorptionSpectrum(absorber_name, wavelengths, absorptions). By convention, the naming of the constant should be the same as the absorber_name field. The wavelengths and absorptions arrays must be of the same length and contain corresponding values.
- 2. In the <u>__init__</u> method of the AbsorptionSpectrumLibrary class, the class constants are added to an internal list. This has the benefit of enabling the Library class to be iterable. Add your newly added constant field to the list here.

3. Your absorption spectrum is now usable throughout all of simpa and is accessible using the SPECTRAL_LIBRARY sngleton that can be imported using from simpa.utils import SPECTRAL_LIBRARY.

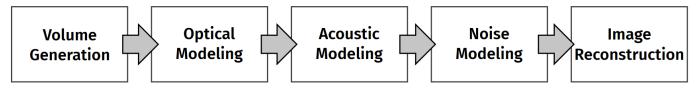
Class references

This component diagram shows the three principle modules of the SIMPA toolkit and gives an insight into their constituents. The core is concerned with providing interfaces for the simulation tools, while the utils module contains many scripts and classes to facilitate the use of the simulation pipeline.



Module: core

The purpose of the core module is to provide interfaces that facilitate the integration of toolboxes and code for photoacoustic modeling into a single continuous pipeline:



Volume creation

The core contribution of the SIMPA toolkit is the creation of in silico tissue-mimicking phantoms. This feature is represented by the volume_creation module, that two main volume creation modules: | Model-based creation of volumes using a set of rules | Segmentation-based creation of volumes

class simpa.core.volume_creation.VolumeCreatorBase
 Use this class to define your own volume creation adapter.

abstract create_simulation_volume (settings: simpa.utils.settings_generator.Settings) \rightarrow dict

This method will be called to create a simulation volume.

Parameters: settings – the settings dictionary containing the simulation instructions.

simpa.core.volume_creation.volume_creation.run_volume_creation(global_settings:
simpa.utils.settings_generator.Settings)

This method is the main entry point of volume creation for the SIMPA framework. It uses the Tags.VOLUME_CREATOR tag to determine which of the volume creators should be used to create the simulation phantom.

Parameters: global settings – the settings dictionary that contains the simulation instructions

Model-based volume creation

class simpa.core.volume_creation.versatile_volume_creator.ModelBasedVolumeCreator
The model-based volume creator uses a set of rules how to generate structures to create a simulation volume.
These structures are added to the dictionary and later combined by the algorithm:

```
# Initialise settings dictionaries
simulation_settings = Settings()
all_structures = Settings()
structure = Settings()
# Definition of en example structure.
# The concrete structure parameters will change depending on the
# structure type
structure[Tags.PRIORITY] = 1
structure[Tags.STRUCTURE_START_MM] = [0, 0, 0]
structure[Tags.STRUCTURE_END_MM] = [0, 0, 100]
structure[Tags.MOLECULE_COMPOSITION] = TISSUE_LIBRARY.muscle()
structure[Tags.CONSIDER_PARTIAL_VOLUME] = True
structure[Tags.ADHERE_TO_DEFORMATION] = True
structure[Tags.STRUCTURE_TYPE] = Tags.HORIZONTAL_LAYER_STRUCTURE
all_structures["arbitrary_identifier"] = structure
simulation_settings[Tags.STRUCTURES] = all_structures
# Define further simulation settings
 . . .
simulate(simulation_settings)
```

create simulation volume (settings) → dict

This method creates a in silico respresentation of a tissue as described in the settings file that is given.

Parameters: settings - a dictionary containing all relevant Tags for the simulation to be able to

instantiate a tissue.

Returns: a path to a npz file containing characteristics of the simulated volume: absorption,

scattering, anisotropy, oxygenation, and a segmentation mask. All of these are given as

3d numpy arrays.

Segmentation-based volume creation

class simpa.core.volume_creation.segmentation_based_volume_creator.SegmentationBasedVol
umeCreator

This volume creator expects a np.ndarray to be in the settigs under the Tags.INPUT_SEGMENTATION_VOLUME tag and uses this array together with a SegmentationClass mapping which is a dict defined in the settings under Tags.SEGMENTATION_CLASS_MAPPING.

With this, an even greater utility is warranted.

 $\label{create_simulation_volume} \textbf{(} \texttt{settings: simpa.utils.settings_generator.Settings)} \rightarrow \textbf{dict}$ This method will be called to create a simulation volume.

Parameters: settings – the settings dictionary containing the simulation instructions.

Optical forward modeling

simpa.core.optical_simulation.optical_modelling.run_optical_forward_model (settings)

This method is the main entry point for running optical forward simulations with the SIMPA toolkit. It is important, that the Tags.OPTICAL_MODEL tag is set in the settings dictionary, as well as any tags that have to be present for the specific model.

class simpa.core.optical_simulation.OpticalForwardAdapterBase
 Use this class as a base for implementations of optical forward models.

abstract forward_model (absorption_cm, scattering_cm, anisotropy, settings)
A deriving class needs to implement this method according to its model.

Parameters:

- absorption_cm Absorption in units of per centimeter
- scattering_cm Scattering in units of per centimeter
- anisotropy Dimensionless scattering anisotropy
- settings Setting dictionary

Returns: Fluence in units of J/cm^2

simulate (optical_properties_path, settings)

Call this method to invoke the simulation process.

A adapter that implements the forward_model method, will take optical properties of absorption, scattering, and scattering anisotropy as input and return the light fluence as output.

Parameters:

- optical_properties_path path to a .npz file that contains the following tags: Tags.PROPERTY_ABSORPTION_PER_CM -> contains the optical absorptions in units of one per centimeter Tags.PROPERTY_SCATTERING_PER_CM -> contains the optical scattering in units of one per centimeter Tags.PROPERTY_ANISOTROPY -> contains the dimensionless optical scattering anisotropy
- settings –

Returns:

simpa.core.optical_simulation.illumination_definition.define_illumination (settings, nx,
ny, nz)

This method creates a dictionary that represents the illumination geometry in a way that it can be used with the respective illumination framework.

Parameters:

- settings The settings file containing the simulation instructions
- nx number of voxels along the x dimension of the volume
- ny number of voxels along the y dimension of the volume
- ullet nz number of voxels along the z dimension of the volume

 $\label{eq:core.optical_simulation} simpa.core.optical_simulation.illumination_definition.define_illumination_mcx (settings, nx, ny, nz) \rightarrow dict$

This method creates a dictionary that contains tags as they are expected for the mcx simulation tool to represent the illumination geometry.

Parameters:

- settings The settings file containing the simulation instructions
- **nx** number of voxels along the x dimension of the volume
- ny number of voxels along the y dimension of the volume
- nz number of voxels along the z dimension of the volume

mcx integration

class simpa.core.optical_simulation.mcx_adapter.McxAdapter

This class implements a bridge to the mcx framework to integrate mcx into SIMPA. MCX is a GPU-enabled Monte-Carlo model simulation of photon transport in tissue:

Fang, Qianqian, and David A. Boas. "Monte Carlo simulation of photon migration in 3D turbid media accelerated by graphics processing units."

Optics express 17.22 (2009): 20178-20190.

forward_model (absorption_cm, scattering_cm, anisotropy, settings)

A deriving class needs to implement this method according to its model.

Parameters:

- absorption_cm Absorption in units of per centimeter
- scattering_cm Scattering in units of per centimeter
- anisotropy Dimensionless scattering anisotropy
- settings Setting dictionary

Returns: Fluence in units of J/cm^2

Acoustic forward modeling

simpa.core.acoustic_simulation.acoustic_modelling.run_acoustic_forward_model (settings)
This method is the entry method for running an acoustic forward model. It is invoked in the simpa.core.simulation.simulate method, but can also be called individually for the purposes of performing acoustic forward modeling only or in a different context.

The concrete will be chosen based on the:

Tags.ACOUSTIC MODEL

tag in the settings dictionary.

Parameters: settings – The settings dictionary containing key-value pairs that determine the simulation.

Here, it must contain the Tags.ACOUSTIC_MODEL tag and any tags that might be required

by the specific acoustic model.

Raises: AssertionError - an assertion error is raised if the Tags.ACOUSTIC MODEL tag is not

given or points to an unknown acoustic forward model.

Returns: returns the path to the simulated data within the saved HDF5 container.

class simpa.core.acoustic_simulation.AcousticForwardAdapterBase

This class should be used as a base for implementations of acoustic forward models.

abstract forward_model (settings) → numpy.ndarray

This method performs the acoustic forward modeling given the initial pressure distribution and the acoustic tissue properties contained in the settings file. A deriving class needs to implement this method according to its model.

Parameters: settings – Setting dictionary Returns: time series pressure data

 $\textbf{simulate (} \texttt{settings)} \rightarrow \textbf{numpy.ndarray}$

Call this method to invoke the simulation process.

Parameters: settings – the settings dictionary containing all simulation parameters.

Returns: a numpy array containing the time series pressure data per detection element

k-Wave integration

class simpa.core.acoustic_simulation.k_wave_adapter.KwaveAcousticForwardModel

The KwaveAcousticForwardModel adapter enables acoustic simulations to be run with the k-wave MATLAB toolbox. k-Wave is a free toolbox (http://www.k-wave.org/) developed by Bradley Treeby and Ben Cox (University College London) and Jiri Jaros (Brno University of Technology).

In order to use this toolbox, MATLAB needs to be installed on your system and the path to the MATLAB binary needs to be specified in the settings dictionary.

In order to use the toolbox from with SIMPA, a number of parameters have to be specified in the settings dictionary:

```
The initial pressure distribution:
    Tags.OPTICAL_MODEL_INITIAL_PRESSURE
Acoustic tissue properties:
    Tags.PROPERTY SPEED OF SOUND
    Tags.PROPERTY_DENSITY
    Tags.PROPERTY_ALPHA_COEFF
The digital twin of the imaging device:
    Tags.DIGITAL DEVICE
Other parameters:
    Tags.PERFORM UPSAMPLING
    Tags.SPACING MM
    Tags.UPSCALE_FACTOR
    Tags.MEDIUM_ALPHA_POWER
    Tags.GPU
    Tags.PMLInside
    Tags.PMLAlpha
    Tags.PlotPML
    Tags.RECORDMOVIE
    Tags.MOVIENAME
    Tags.ACOUSTIC_LOG_SCALE
    Tags.SENSOR DIRECTIVITY PATTERN
```

Many of these will be set automatically by SIMPA, but you may use the simpa.utils.settings_generator convenience methods to generate settings files that contain sensible defaults for these parameters.

Please also refer to the simpa_examples scripts to see how the settings file can be parametrized successfully.

forward_model (settings) → numpy.ndarray

This method performs the acoustic forward modeling given the initial pressure distribution and the acoustic tissue properties contained in the settings file. A deriving class needs to implement this method according to its model.

Parameters: settings – Setting dictionary Returns: time series pressure data

Noise modeling

```
simpa.core.noise_simulation.noise_modelling.apply_noise_model_to_time_series_data
(settings, acoustic_model_result_path)
```

This is the primary method for performing noise pertubation of data.

The noise model can be activated using the Tags.NOISE_MODEL and Tags.APPLY_NOISE_MODEL tags.

Parameters:

- settings the settings dictionary containing the simulation instructions
- acoustic_model_result_path path where the data is within the HDF5 file.

Returns:

```
class simpa.core.noise_simulation.GaussianNoiseModel
```

The purpose of the GaussianNoiseModel class is to apply an additive gaussian noise to the input data.

The mean and standard deviation of the model can be defined either by using the Tags.NOISE_MEAN and Tags.NOISE_STD tags, but they can also be set using a pandas dataframe that contains mean and standard deviation of noise for different wavelengths and temperatures. This can be done using the Tags.NOISE_MODEL_PATH tag.

```
apply_noise_model (time_series_data, settings)
```

Applies the defined noise model to the input time series data.

Parameters:

- time_series_data the data the noise should be applied to.
- settings the settings dictionary that contains the simulation instructions.

Returns: a numpy array of the same shape as the input data.

class simpa.core.noise_simulation.NoiseModelAdapterBase

This class functions as a base class that can be used to easily define different noise models by extending the apply_noise_model function.

abstract apply_noise_model (time_series_data: numpy.ndarray, settings: dict) \rightarrow numpy.ndarray

Applies the defined noise model to the input time series data.

Parameters:

- time_series_data the data the noise should be applied to.
- settings the settings dictionary that contains the simulation instructions.

Returns: a numpy array of the same shape as the input data.

Image reconstruction

 ${\it class} \ {\it simpa.core.image_reconstruction.} \textbf{ReconstructionAdapterBase} \\ \ TODO$

abstract reconstruction_algorithm (time_series_sensor_data, settings)

A deriving class needs to implement this method according to its model.

Parameters:

- time_series_sensor_data the time series sensor data
- settings Setting dictionary

Returns: a reconstructed photoacoustic image

simulate (settings)

Parameters: settings -

Returns:

Backprojection

Time Revearsal

class simpa.core.image_reconstruction.TimeReversalAdapter.TimeReversalAdapter

The time reversal adapter includes the time reversal reconstruction algorithm implemented by the k-Wave toolkit into SIMPA.

Time reversal reconstruction uses the time series data and computes the forward simulation model backwards in time:

```
Treeby, Bradley E., Edward Z. Zhang, and Benjamin T. Cox.
"Photoacoustic tomography in absorbing acoustic media using time reversal." Inverse Problems 26.11 (2010): 115003.
```

```
Static get_acoustic_properties (global_settings: dict, input_data: dict)
```

This method extracts the acoustic tissue properties from the settings dictionary and amends the information to the input_data.

Parameters:

- **global_settings** the settings dictionary containing key value pairs with the simulation instructions.
- input_data a dictionary containing the information needed for time reversal.

reconstruction_algorithm (time_series_sensor_data, settings)

A deriving class needs to implement this method according to its model.

Parameters:

• time_series_sensor_data - the time series sensor data

settings – Setting dictionary

Returns: a reconstructed photoacoustic image

Digital device twins

At every step along the forward simulation, knowledge of the photoacoustic device that is used for the measurements is needed. This is important to reflect characteristic artefacts and challenges for the respective device.

To this end, we have included digital twins of commonly used devices into the SIMPA core.

MSOT Acuity Echo

class simpa.core.device_digital_twins.msot_devices.MSOTAcuityEcho

This class represents a digital twin of the MSOT Acuity Echo, manufactured by iThera Medical, Munich, Germany (https://www.ithera-medical.com/products/msot-acuity/). It is based on the real specifications of the device, but due to the limitations of the possibilities how to represent a device in the software frameworks, constitutes only an approximation.

Some important publications that showcase the use cases of the MSOT Acuity and Acuity Echo device are:

Regensburger, Adrian P., et al. "Detection of collagens by multispectral optoacoustic tomography as an imaging biomarker for Duchenne muscular dystrophy."

Nature Medicine 25.12 (2019): 1905-1915.

Knieling, Ferdinand, et al. "Multispectral Optoacoustic Tomography for Assessment of Crohn's Disease Activity."

The New England journal of medicine 376.13 (2017): 1292.

adjust_simulation_volume_and_settings

(global settings:

simpa.utils.settings_generator.Settings)

In case that the PAI device needs space for the arrangement of detectors or illuminators in the volume, this method will update the volume accordingly.

check_settings_prerequisites

(global_settings:

 ${\tt simpa.utils.settings_generator.Settings)} \rightarrow {\tt bool}$

It might be that certain device geometries need a certain dimensionality of the simulated PAI volume, or that it required the existence of certain Tags in the global global_settings. To this end, a PAI device should use this method to inform the user about a mismatch of the desired device and throw a ValueError if that is the case.

Raises: ValueError – raises a value error if the prerequisites are not matched.

Returns: True if the prerequisites are met, False if they are not met, but no exception has been

raised.

get_detector_element_orientations

(global_settings:

simpa.utils.settings_generator.Settings $) \rightarrow numpy.ndarray$

This method yields a normalised orientation vector for each detection element. The length of this vector is the same as the one obtained via the position methods:

```
get_detector_element_positions_base_mm
get_detector_element_positions_accounting_for_device_position_mm
```

Returns: a numpy array that contains normalised orientation vectors for each detection element

Similar to:

```
get_detector_element_positions_base_mm
```

This method returns the absolute positions of the detection elements relative to the device position in the imaged volume, where the device position is defined by the following tag:

```
Tags.DIGITAL DEVICE POSITION
```

Returns: A numpy array containing the coordinates of the detection elements

```
get_detector_element_positions_base_mm () → numpy.ndarray
```

Defines the abstract positions of the detection elements in an arbitraty coordinate system. Typically, the center of the field of view is defined as the origin.

To obtain the positions in an interpretable coordinate system, please use the other method:

```
get_detector_element_positions_accounting_for_device_position_mm
```

Returns: A numpy array containing the position vestors of the detection elements.

get_illuminator_definition (global_settings: simpa.utils.settings_generator.Settings)
Defines the illumination geometry of the device in the settings dictionary.

RSOM Explorer P50

This class represents an approximation of the Raster-scanning Optoacoustic Mesoscopy (RSOM) device built by iThera Medical (Munich, Germany). Please refer to the companie's website for more information (https://www.ithera-medical.com/products/rsom-explorer-p50/).

Since simulating thousands of individual forward modeling steps to obtain a single raster-scanned image is computationally not feasible, we approximate the process with a device design that has detection elements across the entire field of view. Because of this limitation we also need to approximate the light source with a homogeneous illumination across the field of view.

The digital device is modeled based on the reported specifications of the RSOM Explorer P50 system. Technical details of the system can be found in the dissertation of Mathias Schwarz (https://mediatum.ub.tum.de/doc/1324031/1324031.pdf) and you can find more details on use cases of the device in the following literature sources:

Yew, Yik Weng, et al. "Raster-scanning optoacoustic mesoscopy (RSOM) imaging as an objective disease severity tool in atopic dermatitis patients." Journal of the American Academy of Dermatology (2020).

Hindelang, B., et al. "Non-invasive imaging in dermatology and the unique potential of raster-scan optoacoustic mesoscopy."

Journal of the European Academy of Dermatology and Venereology 33.6 (2019): 1051-1061.

adjust_simulation_volume_and_settings

(global_settings:

simpa.utils.settings_generator.Settings)

In case that the PAI device needs space for the arrangement of detectors or illuminators in the volume, this method will update the volume accordingly.

check_settings_prerequisites

(global_settings:

simpa.utils.settings_generator.Settings $) \rightarrow bool$

It might be that certain device geometries need a certain dimensionality of the simulated PAI volume, or that it required the existence of certain Tags in the global global_settings. To this end, a PAI device should use this method to inform the user about a mismatch of the desired device and throw a ValueError if that is the case.

Raises: ValueError – raises a value error if the prerequisites are not matched.

Returns: True if the prerequisites are met, False if they are not met, but no exception has been

raised.

get_detector_element_orientations

(global_settings:

simpa.utils.settings_generator.Settings)

This method yields a normalised orientation vector for each detection element. The length of this vector is the same as the one obtained via the position methods:

```
get_detector_element_positions_base_mm
get_detector_element_positions_accounting_for_device_position_mm
```

Returns: a numpy array that contains normalised orientation vectors for each detection element

This method returns the absolute positions of the detection elements relative to the device position in the imaged volume, where the device position is defined by the following tag:

```
Tags.DIGITAL_DEVICE_POSITION
```

Returns: A numpy array containing the coordinates of the detection elements

```
get_detector_element_positions_base_mm()
```

Defines the abstract positions of the detection elements in an arbitraty coordinate system. Typically, the center of the field of view is defined as the origin.

To obtain the positions in an interpretable coordinate system, please use the other method:

```
get_detector_element_positions_accounting_for_device_position_mm
```

Returns: A numpy array containing the position vestors of the detection elements.

get_illuminator_definition (global_settings: simpa.utils.settings_generator.Settings)
Defines the illumination geometry of the device in the settings dictionary.

Module: utils

The utils module contains several general-purpose utility functions whose purpose it is to facilitate the use of SIMPA. The most important of these is the Tags class, which defines the strings and data types that have to be used for the keys and values of the settings dictionary.

```
class simpa.utils.tags.Tags
```

This class contains all 'Tags' for the use in the settings dictionary as well as strings that are used in SIMPA as naming conventions. Every Tag that is intended to be used as a key in the settings dictionary is represented by a tuple. The first element of the tuple is a string that corresponds to the name of the Tag. The second element of the tuple is a data type or a tuple of data types. The values that are assigned to the keys in the settings should match these data types. Their usage within the SIMPA package is divided in "SIMPA package", "module X", "adapter Y", "class Z" and "naming convention".

```
ACOUSTIC_MODEL = ('acoustic_model', <class 'str'>)
Choice of the used acoustic model.
Usage: module acoustic simulation
```

ACOUSTIC MODEL K WAVE = 'kwave'

Corresponds to the kwave simulaiton.

Usage: module acoustic_simulation, naming convention

ACOUSTIC MODEL SCRIPT LOCATION = ('acoustic model script location', <class 'str'>)

Absolute path of the location of the acoustic_simulation folder in the SIMPA core module.

Usage: module acoustic_simulation

ACOUSTIC_MODEL_TEST = 'simpa_tests'

Corresponds to an adapter for testing purposes only.

Usage: module acoustic_simulation, naming convention

ACOUSTIC_SIMULATION_3D = ('acoustic_simulation_3d', <class 'bool'>)

If True, simulates the acoustic forward model in 3D.

Usage: SIMPA package

ADHERE_TO_DEFORMATION = ('adhere_to_deformation', <class 'bool'>)

If True, a structure will be shifted according to the deformation.

Usage: adapter versatile_volume_creation

BACKGROUND = 'Background'

Corresponds to the name of a structure.

Usage: adapter versatile_volume_creation, naming convention

CONSIDER_PARTIAL_VOLUME = ('consider_partial_volume', <class 'bool'>)

If True, the structure will be generated with its edges only occupying a partial volume of the voxel.

Usage: adapter versatile_volume_creation

DEFORMATION_X_COORDINATES_MM = 'deformation_x_coordinates'

Mesh that defines the x coordinates of the deformation.

Usage: adapter versatile_volume_creation, naming convention

DEFORMATION_Y_COORDINATES_MM = 'deformation_y_coordinates'

Mesh that defines the y coordinates of the deformation.

Usage: adapter versatile_volume_creation, naming convention

DEFORMATION Z ELEVATIONS MM = 'deformation z elevation'

Mesh that defines the z coordinates of the deformation.

Usage: adapter versatile_volume_creation, naming convention

DEFORMED LAYERS SETTINGS = ('deformed layers settings', <class 'dict'>)

Settings that contain the functional which defines the deformation of the layers.

Usage: adapter versatile_volume_creation

DIGITAL DEVICE = ('digital device', <class 'str'>)

Digital device that is chosen as illumination source and detector for the simulation.

Usage: SIMPA package

DIGITAL_DEVICE_MSOT = 'digital_device_msot'

Corresponds to the MSOTAcuityEcho device.

Usage: SIMPA package, naming convention

DIGITAL_DEVICE_POSITION = ('digital_device_position', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Position in [x, y, z] coordinates of the device in the generated volume.

Usage: SIMPA package

DIGITAL_DEVICE_RSOM = 'digital_device_rsom'

Corresponds to the RSOMExplorerP50 device.

Usage: SIMPA package, naming convention

GPU = ('gpu', (<class 'bool'>, <class 'bool'>, <class 'numpy.bool '>))

If True, uses all available gpu options of the used modules.

Usage: SIMPA package

ILLUMINATION_DIRECTION = ('illumination_direction', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Direction of the photon source as [x, y, z] vector used in mcx.

Usage: module optical_modelling, adapter mcx_adapter

ILLUMINATION_PARAM1 = ('illumination_param1', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

First parameter group of the specified illumination type as [x, y, z, w] vector used in mcx.

Usage: module optical_modelling, adapter mcx_adapter

ILLUMINATION_PARAM2 = ('illumination_param2', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Second parameter group of the specified illumination type as [x, y, z, w] vector used in mcx.

Usage: module optical_modelling, adapter mcx_adapter

ILLUMINATION_POSITION = ('illumination_position', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Position of the photon source in [x, y, z] coordinates used in mcx.

Usage: module optical_modelling, adapter mcx_adapter

ILLUMINATION_TYPE = ('optical_model_illumination_type', <class 'str'>)

Type of the illumination geometry used in mcx.

Usage: module optical_modelling, adapter mcx_adapter

ILLUMINATION_TYPE_DISK = 'disk'

Corresponds to disk source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION_TYPE_DKFZ_PAUS = 'pasetup'

Corresponds to pasetup source in mcx. The geometrical definition is described in:

Usage: adapter mcx_adapter, naming convention

ILLUMINATION TYPE FOURIER = 'fourier'

Corresponds to fourier source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION TYPE FOURIER X = 'fourierx'

Corresponds to fourierx source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION TYPE FOURIER X 2D = 'fourierx2d'

Corresponds to fourierx2d source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION TYPE GAUSSIAN = 'gaussian'

Corresponds to gaussian source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION_TYPE_MSOT_ACUITY_ECHO = 'msot_acuity_echo'

Corresponds to msot_acuity_echo source in mcx. The device is manufactured by iThera Medical, Munich, Germany (https://www.ithera-medical.com/products/msot-acuity/).

Usage: adapter mcx_adapter, naming convention

ILLUMINATION TYPE PATTERN = 'pattern'

Corresponds to pattern source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION_TYPE_PATTERN_3D = 'pattern3d'

Corresponds to pattern3d source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION TYPE PENCIL = 'pencil'

Corresponds to pencil source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION TYPE PENCILARRAY = 'pencilarray'

Corresponds to pencilarray source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION_TYPE_PLANAR = 'planar'

Corresponds to planar source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION_TYPE_RING = 'ring'

Corresponds to ring source in mcx.

Usage: adapter mcx_adapter, naming convention

ILLUMINATION_TYPE_SLIT = 'slit'

Corresponds to slit source in mcx.

Usage: adapter mcx_adapter, naming convention

INPUT_SEGMENTATION_VOLUME = ('input_segmentation_volume', <class 'numpy.ndarray'>)

Array that defines a segmented volume.

Usage: adapter segmentation_based_volume_creator

K_WAVE_SPECIFIC_DT = ('dt_acoustic_sim', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>)) Temporal resolution of kwave.

Usage: adapter k_wave_adapter, adapter TimeReversalAdapter

K_WAVE_SPECIFIC_NT = ('Nt_acoustic_sim', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>))

Total time steps simulated by kwave.

Usage: adapter k_wave_adapter, adapter TimeReversalAdapter

LASER_PULSE_ENERGY_IN_MILLIJOULE = ('laser_pulse_energy_in_millijoule', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>, <class 'list'>, <class 'range'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Laser pulse energy used in the optical simulation.

Usage: module optical_simulation

MAX DEFORMATION MM = 'max deformation'

Maximum deformation in z-direction.

Usage: adapter versatile_volume_creation, naming convention

MEDIUM_TEMPERATURE_CELCIUS = ('medium_temperature', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>))

Temperature of the simulated volume.

Usage: module noise_simulation

MOLECULE_COMPOSITION = ('molecule_composition', <class 'list'>)

List that contains all the molecules within a structure.

Usage: module volume_creation

OPTICAL_MODEL = ('optical_model', <class 'str'>)

Choice of the used optical model.

Usage: module optical_simulation

OPTICAL_MODEL_BINARY_PATH = ('optical_model_binary_path', <class 'str'>)

Absolute path of the location of the optical forward model binary.

Usage: module optical_simulation

OPTICAL_MODEL_FLUENCE = 'fluence'

Location of the optical forward model output fluence in the SIMPA output file.

Usage: naming convention

OPTICAL_MODEL_ILLUMINATION_GEOMETRY_XML_FILE = ('optical_model_illumination_geometry_xml_file', <class 'str'>)

Absolute path of the location of the optical forward model illumination geometry.

Usage: module optical_simulation

OPTICAL_MODEL_INITIAL_PRESSURE = 'initial_pressure'

Location of the optical forward model output initial pressure in the SIMPA output file.

Usage: naming convention

OPTICAL_MODEL_MCX = 'mcx'

Corresponds to the mcx simulation.

Usage: module optical_simulation, naming convention

OPTICAL_MODEL_MCXYZ = 'mcxyz'

Corresponds to the mcxyz simulation.

Usage: module optical_simulation, naming convention

OPTICAL_MODEL_NUMBER_PHOTONS = ('optical_model_number_of_photons', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>))

Number of photons used in the optical simulation.

Usage: module optical simulation

OPTICAL_MODEL_OUTPUT_NAME = 'optical_forward_model_output'

Location of the optical forward model output in the SIMPA output file.

Usage: naming convention

OPTICAL_MODEL_TEST = 'simpa_tests'

Corresponds to an adapter for testing purposes only.

Usage: module optical_simulation, naming convention

OPTICAL MODEL UNITS = 'units'

Location of the optical forward model output units in the SIMPA output file.

Usage: naming convention

PRIORITY = ('priority', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>))

Number that corresponds to a priority of the assigned structure. If another structure occupies the same voxel in a volume, the structure with a higher priority will be preferred.

Usage: adapter versatile_volume_creator

PROPERTY_SEGMENTATION = 'seg'

We define PROPERTY_GRUNEISEN_PARAMETER to contain all wavelength-independent constituents of the PA signal. This means that it contains the percentage of absorbed light converted into heat. Naturally, one could make an argument that this should not be the case, however, it simplifies the usage of this tool.

RANDOM_SEED = ('random_seed', (<class 'int'>, <class 'numpy.integer'>))

Random seed for numpy and torch.

Usage: SIMPA package

RECONSTRUCTION_MODE_FULL = 'full'

Upsampling settings

RUN_OPTICAL_MODEL = ('run_optical_forward_model', <class 'bool'>)

If True, the simulation will run the optical forward model.

Usage: module core (simulate.py)

SEGMENTATION_CLASS_MAPPING = ('segmentation_class_mapping', <class 'dict'>)

Mapping that assigns every class in the INPUT_SEGMENTATION_VOLUME a MOLECULE_COMPOSITION.

Usage: adapter segmentation_based_volume_creator

SIMULATE DEFORMED LAYERS = ('simulate deformed layers', <class 'bool'>)

If True, the horizontal layers are deformed according to the DEFORMED LAYERS SETTINGS.

Usage: adapter versatile_volume_creation

SIMULATION EXTRACT FIELD OF VIEW = ('extract field of view', <class 'bool'>)

If True, converts a 3D volume to a 2D volume by extracting the middle slice along the y-axis.

Usage: SIMPA package

SIMULATION_PATH = ('simulation_path', <class 'str'>)

Absolute path to the folder where the SIMPA output is saved.

Usage: SIMPA package

STRUCTURE_BIFURCATION_LENGTH_MM = ('structure_bifurcation_length_mm', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>))

Length after which a VesselStructure will bifurcate.

Usage: adapter versatile_volume_creation, class VesselStructure

STRUCTURE_CURVATURE_FACTOR = ('structure_curvature_factor', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>))

Factor that determines how strongly a vessel tree is curved.

Usage: adapter versatile_volume_creation, class VesselStructure

STRUCTURE_DIRECTION = ('structure_direction', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Direction as [x, y, z] vector starting from STRUCTURE_START_MM in which the vessel will grow.

Usage: adapter versatile_volume_creation, class VesselStructure

STRUCTURE_ECCENTRICITY = ('structure_excentricity', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'numpy.ndarray'>))

Eccentricity of the structure.

Usage: adapter versatile_volume_creation, class EllipticalTubularStructure

STRUCTURE_END_MM = ('structure_end', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Ending of the structure as [x, y, z] coordinates in the generated volume.

Usage: adapter versatile_volume_creation, class GeometricalStructure

STRUCTURE_FIRST_EDGE_MM = ('structure_first_edge_mm', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Edge of the structure as [x, y, z] vector starting from STRUCTURE_START_MM in the generated volume.

Usage: adapter versatile_volume_creation, class ParallelepipedStructure

STRUCTURE_RADIUS_MM = ('structure_radius', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>, <class 'numpy.ndarrav'>))

Radius of the structure.

Usage: adapter versatile volume creation, class GeometricalStructure

STRUCTURE_RADIUS_VARIATION_FACTOR = ('structure_radius_variation_factor', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>))

Factor that determines how strongly a the radius of vessel tree varies.

Usage: adapter versatile_volume_creation, class VesselStructure

STRUCTURE_SECOND_EDGE_MM = ('structure_second_edge_mm', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Edge of the structure as [x, y, z] vector starting from STRUCTURE_START_MM in the generated volume.

Usage: adapter versatile volume creation, class ParallelepipedStructure

STRUCTURE_START_MM = ('structure_start', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Beginning of the structure as [x, y, z] coordinates in the generated volume.

Usage: adapter versatile_volume_creation, class GeometricalStructure

STRUCTURE_THIRD_EDGE_MM = ('structure_third_edge_mm', (<class 'list'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Edge of the structure as [x, y, z] vector starting from STRUCTURE_START_MM in the generated volume. Usage: adapter versatile volume creation, class ParallelepipedStructure

STRUCTURE_X_EXTENT_MM = ('structure_x_extent_mm', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>))

X-extent of the structure in the generated volume.

Usage: adapter versatile_volume_creation, class RectangularCuboidStructure

STRUCTURE_Y_EXTENT_MM = ('structure_y_extent_mm', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>))

Y-extent of the structure in the generated volume.

Usage: adapter versatile volume creation, class RectangularCuboidStructure

STRUCTURE_Z_EXTENT_MM = ('structure_z_extent_mm', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>, <class 'float'>))

Z-extent of the structure in the generated volume.

Usage: adapter versatile_volume_creation, class RectangularCuboidStructure

TIME_REVEARSAL_SCRIPT_LOCATION = ('time_revearsal_script_location', <class 'str'>)

Absolute path of the location of the image_reconstruction folder in the SIMPA core module.

Usage: adapter TimeReversalAdapter

TIME_STEP = ('time_step', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>))

Temporal resolution of mcx. Usage: adapter mcx_adapter

TISSUE_PROPERTIES_OUPUT_NAME = 'properties'

Location of the simulation properties in the SIMPA output file

Usage: naming convention

TOTAL_TIME = ('total_time', (<class 'int'>, <class 'numpy.integer'>, <class 'float'>))

Total simulated time in mcx. Usage: adapter mcx_adapter

UNITS_PRESSURE = 'newton_per_meters_squared'

IO settings

VOLUME_CREATOR = ('volume_creator', <class 'str'>)

Choice of the volume creator adapter.

Usage: module volume_creation, module device_digital_twins

VOLUME_CREATOR_SEGMENTATION_BASED = 'volume_creator_segmentation_based'

Corresponds to the SegmentationBasedVolumeCreator.

Usage: module volume_creation, naming convention

VOLUME_CREATOR_VERSATILE = 'volume_creator_versatile'

Corresponds to the ModelBasedVolumeCreator.

Usage: module volume creation, naming convention

VOLUME_NAME = ('volume_name', <class 'str'>)

Name of the SIMPA output file.

Usage: SIMPA package

WAVELENGTH = ('wavelength', (<class 'int'>, <class 'numpy.integer'>))

Single wavelength used for the current simulation.

Usage: SIMPA package

WAVELENGTHS = ('wavelengths', (<class 'list'>, <class 'range'>, <class 'tuple'>, <class 'numpy.ndarray'>))

Iterable of all the wavelengths used for the simulation.

Usage: SIMPA package

```
class simpa.utils.constants.SaveFilePaths
```

The save file paths specify the path of a specific data structure in the dictionary of the simpa output hdf5. All of these paths have to be used like: SaveFilePaths.PATH.format(Tags.UPSAMPLED_DATA or Tags.ORIGINAL_DATA, wavelength)

class simpa.utils.constants.SegmentationClasses

The segmentation classes define which "tissue types" are modelled in the simulation volumes.

```
simpa.utils.deformation_manager.create_deformation_settings (bounds_mm, maximum_z_elevation_mm=1, filter_sigma=1, cosine_scaling_factor=4)
```

FIXME

simpa.utils.deformation_manager.get_functional_from_deformation_settings
(deformation_settings: dict)

FIXME

class simpa.utils.settings_generator.Settings (dictionary: dict = None)

simpa.utils.calculate.calculate_gruneisen_parameter_from_temperature
(temperature_in_celcius)

This function returns the dimensionless gruneisen parameter based on a heuristic formula that was determined experimentally:

```
@book{wang2012biomedical,
    title={Biomedical optics: principles and imaging},
    author={Wang, Lihong V and Wu, Hsin-i},
    year={2012},
    publisher={John Wiley \& Sons}
}
```

Parameters: temperature_in_celcius - the temperature in degrees celcius

Returns: a floating point number, if temperature_in_celcius is a number or a float array, if

temperature_in_celcius is an array

simpa.utils.calculate.calculate_oxygenation (molecule_list)

Returns: an oxygenation value between 0 and 1 if possible, or None, if not computable.

```
simpa.utils.calculate.create_spline_for_range (xmin_mm=0, xmax_mm=10,
maximum_y_elevation_mm=1, spacing=0.1)
```

Creates a functional that simulates distortion along the y position between the minimum and maximum x positions. The elevation can never be smaller than 0 or bigger than maximum_y_elevation_mm.

Parameters:

- xmin mm the minimum x axis value the return functional is defined in
- xmax_mm the maximum x axis value the return functional is defined in
- maximum_y_elevation_mm the maximum y axis value the return functional will yield

Returns: a functional that describes a distortion field along the y axis

simpa.utils.calculate.randomize_uniform (min_value: float, max_value: float)
returns a uniformly drawn random number in [min_value, max_value]

Parameters:

- min_value minimum value
- max_value maximum value

Returns: random number in [min_value, max_value]

class simpa.utils.tissue_properties.TissueProperties

Libraries

Another important aspect of the utils class is the libraries that are being provided. These contain compilations of literature values for the acoustic and optical properties of commonly used tissue.

update_internal_properties()
FIXME

class simpa.utils.libraries.literature_values.MorphologicalTissueProperties

This class contains a listing of morphological tissue parameters as reported in literature. The listing is not the result of a meta analysis, but rather uses the best fitting paper at the time pf implementation. Each of the fields is annotated with a literature reference or a descriptions of how the particular values were derived for tissue modelling.

class simpa.utils.libraries.literature_values.OpticalTissueProperties

This class contains a listing of optical tissue parameters as reported in literature. The listing is not the result of a meta analysis, but rather uses the best fitting paper at the time pf implementation. Each of the fields is annotated with a literature reference or a descriptions of how the particular values were derived for tissue modelling.

class simpa.utils.libraries.literature_values.StandardProperties

This class contains a listing of default parameters that can be used. These values are sensible default values but are generally not backed up by proper scientific references, or are rather specific for internal use cases.

class simpa.utils.libraries.spectra_library.AbsorptionSpectrum (spectrum_name: str,
wavelengths: numpy.ndarray, absorption_per_centimeter: numpy.ndarray)

An instance of this class represents the absorption spectrum over wavelength for a particular

get_absorption_for_wavelength (wavelength: int) → float

Parameters: wavelength – the wavelength to retrieve a optical absorption value for [cm^{-1}]. Must be

an integer value between the minimum and maximum wavelength.

Returns: the best matching linearly interpolated absorption value for the given wavelength.

get_absorption_over_wavelength()

Returns: numpy array with the available wavelengths and the corresponding absorption properties

simpa.utils.libraries.spectra_library.view_absorption_spectra (save_path=None)
Opens a matplotlib plot and visualizes the available absorption spectra.

Parameters: save_path – If not None, then the figure will be saved as a png file to the destination.

class simpa.utils.libraries.tissue_library.MolecularCompositionGenerator

The MolecularCompositionGenerator is a helper class to facilitate the creation of a MolecularComposition instance.

class simpa.utils.libraries.tissue_library.TissueLibrary
TODO

blood_arterial()

Returns: a settings dictionary containing all min and max parameters fitting for full blood.

blood_generic (oxygenation=None)

Returns: a settings dictionary containing all min and max parameters fitting for full blood.

blood_venous()

```
a settings dictionary containing all min and max parameters fitting for full blood.
 bone ()
                     a settings dictionary containing all min and max parameters fitting for full blood.
 constant (mua, mus, g)
   TODO
 dermis (background_oxy=0.5)
                     a settings dictionary containing all min and max parameters fitting for dermis tissue.
 epidermis ()
           Returns: a settings dictionary containing all min and max parameters fitting for epidermis tissue.
 get_blood_volume_fractions (total_blood_volume_fraction, oxygenation)
   TODO
 muscle (background_oxy=0.5)
                     a settings dictionary containing all min and max parameters fitting for generic background
           Returns:
 subcutaneous_fat (background_oxy=0.5)
                     a settings dictionary containing all min and max parameters fitting for subcutaneous fat
                     tissue.
class
             simpa.utils.libraries.structure_library.Background
                                                                                (global_settings:
                                                                            background_settings:
simpa.utils.settings_generator.Settings,
simpa.utils.settings_generator.Settings = None)
 to settings () \rightarrow dict
   TODO :return : A tuple containing the settings key and the needed entries
class simpa.utils.libraries.structure_library.CircularTubularStructure (global_settings:
simpa.utils.settings_generator.Settings,
                                                                     single_structure_settings:
simpa.utils.settings_generator.Settings = None)
 to_settings()
   TODO :return : A tuple containing the settings key and the needed entries
                       simpa.utils.libraries.structure_library.EllipticalTubularStructure
                     simpa.utils.settings_generator.Settings, single_structure_settings:
(global settings:
simpa.utils.settings_generator.Settings = None)
 to settings ()
   TODO :return : A tuple containing the settings key and the needed entries
       simpa.utils.libraries.structure_library.GeometricalStructure
                                                                                (global_settings:
simpa.utils.settings generator.Settings,
                                                                     single structure settings:
simpa.utils.settings_generator.Settings = None)
 TODO
 abstract to_settings () → simpa.utils.settings_generator.Settings
   TODO :return : A tuple containing the settings key and the needed entries
class simpa.utils.libraries.structure_library.HorizontalLayerStructure (global_settings:
simpa.utils.settings_generator.Settings,
                                                                     single_structure_settings:
simpa.utils.settings_generator.Settings = None)
```

Module: io_handling

to_settings()

TODO :return : A tuple containing the settings key and the needed entries

This class currently has no partial volume effects implemented. TODO

to_settings()

TODO :return : A tuple containing the settings key and the needed entries

to_settings()

TODO :return : A tuple containing the settings key and the needed entries

class simpa.utils.libraries.structure_library.SphericalStructure (global_settings:
simpa.utils.settings_generator.Settings, simpa.utils.settings_generator.Settings = None)

to_settings()

TODO :return : A tuple containing the settings key and the needed entries

class simpa.utils.libraries.structure_library.VesselStructure (global_settings:
simpa.utils.settings_generator.Settings, simpa.utils.settings_generator.Settings = None)

to_settings()

TODO :return : A tuple containing the settings key and the needed entries

Module: io_handling

simpa.io_handling.io_hdf5.load_hdf5 (file_path, file_dictionary_path='/')
Loads a dictionary from an hdf5 file.

Parameters:

- file_path Path of the file to load the dictionary from.
- file_dictionary_path Path in dictionary structure of hdf5 file to lo the dictionary in.

Returns: Dictionary

simpa.io_handling.io_hdf5.save_hdf5 (dictionary: dict, file_path: str,
file_dictionary_path: str = '/', file_compression: str = None)
Saves a dictionary with arbitrary content to an hdf5-file with given filepath.

Parameters:

- dictionary Dictionary to save.
- file path Path of the file to save the dictionary in.
- **file_dictionary_path** Path in dictionary structure of existing hdf5 file to store the dictionary in.
- **file_compression** possible file compression for the hdf5 output file. Values are: gzip, lzf and szip.

Returns: Null

Examples

Performing a complete forward simulation with acoustic modeling, optical modeling, as well as image reconstruction

The file can be found in simpa_examples/minimal_optical_simulation.py:

```
from simpa.utils import Tags, TISSUE_LIBRARY
from simpa.core.simulation import simulate
from simpa.utils.settings_generator import Settings
from simpa.utils.libraries.structure_library import HorizontalLayerStructure
import numpy as np
# TODO change these paths to the desired executable and save folder
SAVE_PATH = "D:/save/"
MCX_BINARY_PATH = "D:/bin/Release/mcx.exe"
VOLUME_TRANSDUCER_DIM_IN_MM = 75
VOLUME_PLANAR_DIM_IN_MM = 20
VOLUME_HEIGHT_IN_MM = 25
SPACING = 0.15
RANDOM_SEED = 4711
def create_example_tissue():
    This is a very simple example script of how to create a tissue definition.
    It contains a muscular background, an epidermis layer on top of the muscles
    and a blood vessel.
    background_dictionary = Settings()
    background_dictionary[Tags.MOLECULE_COMPOSITION] = TISSUE_LIBRARY.muscle()
    background_dictionary[Tags.STRUCTURE_TYPE] = Tags.BACKGROUND
    muscle_dictionary = Settings()
    muscle_dictionary[Tags.PRIORITY] = 1
    muscle dictionary[Tags.STRUCTURE START MM] = [0, 0, 0]
    muscle_dictionary[Tags.STRUCTURE_END_MM] = [0, 0, 100]
    muscle_dictionary[Tags.MOLECULE_COMPOSITION] = TISSUE_LIBRARY.muscle()
    muscle_dictionary[Tags.CONSIDER_PARTIAL_VOLUME] = True
    muscle_dictionary[Tags.ADHERE_TO_DEFORMATION] = True
    muscle_dictionary[Tags.STRUCTURE_TYPE] = Tags.HORIZONTAL_LAYER_STRUCTURE
    vessel 1 dictionary = Settings()
    vessel_1_dictionary[Tags.PRIORITY] = 3
    vessel_1_dictionary[Tags.STRUCTURE_START_MM] = [VOLUME_TRANSDUCER_DIM_IN_MM/2,
                                                    0, 10]
    vessel_1_dictionary[Tags.STRUCTURE_END_MM] = [VOLUME_TRANSDUCER_DIM_IN_MM/2, VOLUME_PLAN
    vessel_1_dictionary[Tags.STRUCTURE_RADIUS_MM] = 3
    vessel_1_dictionary[Tags.MOLECULE_COMPOSITION] = TISSUE_LIBRARY.blood_generic()
    vessel_1_dictionary[Tags.CONSIDER_PARTIAL_VOLUME] = True
    vessel_1_dictionary[Tags.STRUCTURE_TYPE] = Tags.CIRCULAR_TUBULAR_STRUCTURE
    epidermis_dictionary = Settings()
    epidermis dictionary[Tags.PRIORITY] = 8
    epidermis_dictionary[Tags.STRUCTURE_START_MM] = [0, 0, 0]
    epidermis_dictionary[Tags.STRUCTURE_END_MM] = [0, 0, 1]
    epidermis_dictionary[Tags.MOLECULE_COMPOSITION] = TISSUE_LIBRARY.epidermis()
```

```
epidermis_dictionary[Tags.CONSIDER_PARTIAL_VOLUME] = True
    epidermis_dictionary[Tags.ADHERE_TO_DEFORMATION] = True
    epidermis_dictionary[Tags.STRUCTURE_TYPE] = Tags.HORIZONTAL_LAYER_STRUCTURE
    tissue_dict = Settings()
    tissue_dict[Tags.BACKGROUND] = background_dictionary
    tissue_dict["muscle"] = muscle_dictionary
    tissue_dict["epidermis"] = epidermis_dictionary
    tissue_dict["vessel_1"] = vessel_1_dictionary
    return tissue_dict
# Seed the numpy random configuration prior to creating the global_settings file in
# order to ensure that the same volume
# is generated with the same random seed every time.
np.random.seed(RANDOM_SEED)
settings = {
    # These parameters set the general propeties of the simulated volume
    Tags.RANDOM_SEED: RANDOM_SEED,
    Tags.VOLUME_NAME: "CompletePipelineTestMSOT_"+str(RANDOM_SEED),
    Tags.SIMULATION_PATH: SAVE_PATH,
    Tags.SPACING_MM: SPACING,
    Tags.DIM_VOLUME_Z_MM: VOLUME_HEIGHT_IN_MM,
    Tags.DIM_VOLUME_X_MM: VOLUME_TRANSDUCER_DIM_IN_MM,
    Tags.DIM_VOLUME_Y_MM: VOLUME_PLANAR_DIM_IN_MM,
    Tags.VOLUME_CREATOR: Tags.VOLUME_CREATOR_VERSATILE,
    Tags.SIMULATE_DEFORMED_LAYERS: True,
    # Tags.DEFORMED_LAYERS_SETTINGS: create_deformation_settings([[0, VOLUME_TRANSDUCER_DIM_
                                                                  [0, VOLUME_PLANAR_DIM_IN_MM
    #
                                                                 maximum_z_elevation_mm=10,
    #
                                                                 filter_sigma=0,
                                                                  cosine_scaling_factor=1),
    # Simulation Device
    Tags.DIGITAL_DEVICE: Tags.DIGITAL_DEVICE_MSOT,
    # The following parameters set the optical forward model
    Tags.RUN_OPTICAL_MODEL: True,
    Tags.WAVELENGTHS: [700],
    Tags.OPTICAL_MODEL_NUMBER_PHOTONS: 1e7,
    Tags.OPTICAL_MODEL_BINARY_PATH: MCX_BINARY_PATH,
    Tags.OPTICAL_MODEL: Tags.OPTICAL_MODEL_MCX,
    Tags.ILLUMINATION_TYPE: Tags.ILLUMINATION_TYPE_MSOT_ACUITY_ECHO,
    Tags.LASER_PULSE_ENERGY_IN_MILLIJOULE: 50,
    # The following parameters tell the script that we do not want any extra
    # modelling steps
    Tags.RUN_ACOUSTIC_MODEL: True,
    Tags.ACOUSTIC_SIMULATION_3D: False,
    Tags.ACOUSTIC_MODEL: Tags.ACOUSTIC_MODEL_K_WAVE,
    Tags.ACOUSTIC_MODEL_BINARY_PATH: "C:/Program Files/MATLAB/R2020b/bin/matlab.exe",
    Tags.ACOUSTIC_MODEL_SCRIPT_LOCATION: "C:/simpa/simpa/core/acoustic_simulation",
    Tags.GPU: True,
    Tags.MEDIUM_ALPHA_POWER: 1.05,
    Tags.SENSOR_RECORD: "p",
    # Tags.SENSOR_DIRECTIVITY_PATTERN: "pressure",
```

```
Tags.PMLInside: False,
    Tags.PMLSize: [31, 32],
    Tags.PMLAlpha: 1.5,
    Tags.PlotPML: False,
    Tags.RECORDMOVIE: False,
    Tags.MOVIENAME: "visualization log",
    Tags.ACOUSTIC_LOG_SCALE: True,
    Tags.APPLY_NOISE_MODEL: False,
    Tags.SIMULATION_EXTRACT_FIELD_OF_VIEW: True,
    Tags.PERFORM_IMAGE_RECONSTRUCTION: True,
    Tags.RECONSTRUCTION_ALGORITHM: Tags.RECONSTRUCTION_ALGORITHM_BACKPROJECTION
settings = Settings(settings)
# global_settings[Tags.SIMULATE_DEFORMED_LAYERS] = True
np.random.seed(RANDOM_SEED)
settings[Tags.STRUCTURES] = create_example_tissue()
print("Simulating ", RANDOM_SEED)
import time
timer = time.time()
simulate(settings)
print("Needed", time.time()-timer, "seconds")
print("Simulating ", RANDOM_SEED, "[Done]")
```

Reading the HDF5 simulation output

The file can be found in simpa_examples/access_saved_PAI_data.py:

```
from simpa.io_handling import load_hdf5, save_hdf5
import matplotlib.pyplot as plt
import matplotlib as mpl
import numpy as np
from simpa.utils import SegmentationClasses, Tags
from simpa.utils.settings_generator import Settings
values = []
names = []
for string in SegmentationClasses.__dict__:
    if string[0:2] != "___":
        values.append(SegmentationClasses.__dict__[string])
        names.append(string)
values = np.asarray(values)
names = np.asarray(names)
sort_indexes = np.argsort(values)
values = values[sort_indexes]
names = names[sort_indexes]
colors = [list(np.random.random(3)) for _ in range(len(names))]
cmap = mpl.colors.LinearSegmentedColormap.from_list(
    'Custom cmap', colors, len(names))
PATH = "D:/save/LNetOpticalForward_planar_0.hdf5"
WAVELENGTH = 532
file = load_hdf5(PATH)
settings = Settings(file["settings"])
```

```
fluence = (file['simulations']['original_data']['optical_forward_model_output']
           [str(WAVELENGTH)]['fluence'])
initial_pressure = (file['simulations']['original_data']
                    ['optical_forward_model_output']
                    [str(WAVELENGTH)]['initial_pressure'])
absorption = (file['simulations']['original_data']['simulation_properties']
              [str(WAVELENGTH)]['mua'])
segmentation = (file['simulations']['original_data']['simulation_properties']
              [str(WAVELENGTH)]['seg'])
reconstruction = None
speed_of_sound = None
if Tags.PERFORM IMAGE RECONSTRUCTION in settings and settings[Tags.PERFORM IMAGE RECONSTRUCT
    time_series = np.squeeze(
        file["simulations"]["original_data"]["time_series_data"][str(WAVELENGTH)]["time_seri
    reconstruction = np.squeeze(
            file["simulations"]["original_data"]["reconstructed_data"][str(WAVELENGTH)]["rec
    speed_of_sound = file['simulations']['original_data']['simulation_properties'][str(WAVEI
reconstruction = reconstruction.T
shape = np.shape(reconstruction)
x_pos = int(shape[0]/2)
y_pos = int(shape[1]/2)
z_pos = int(shape[2]/2)
plt.figure()
plt.subplot(161)
{\tt plt.imshow(np.fliplr(np.rot90(reconstruction[x\_pos, :, :], -1)))}
plt.subplot(162)
plt.imshow(np.rot90(np.log10(initial_pressure[x_pos, :, :]), -1))
plt.subplot(163)
plt.imshow(np.fliplr(np.rot90(reconstruction[:, y_pos, :], -1)))
plt.subplot(164)
plt.imshow(np.rot90(np.log10(initial_pressure[:, y_pos, :]), -1))
plt.subplot(165)
plt.imshow(np.fliplr(np.rot90(reconstruction[:, :, z_pos], -1)))
plt.subplot(166)
plt.imshow(np.rot90(np.log10(initial_pressure[:, :, z_pos]), -3))
plt.show()
exit()
if Tags.PERFORM_IMAGE_RECONSTRUCTION in settings and settings[Tags.PERFORM_IMAGE_RECONSTRUCT
    if len(shape) > 2:
        plt.figure()
        plt.subplot(141)
        plt.imshow(np.rot90(np.log10(np.log10(time_series[:, :]-np.min(time_series))), -1),
        plt.subplot(142)
        plt.imshow(np.rot90((reconstruction[:, y_pos, :]), -2))
        plt.subplot(143)
        plt.imshow(np.rot90(np.log10(initial_pressure[:, y_pos, :]), -1))
        plt.subplot(144)
        plt.imshow(np.rot90(segmentation[:, y_pos, :], -1), vmin=values[0], vmax=values[-1],
        plt.show()
    else:
        plt.figure()
```

```
plt.subplot(141)
        plt.imshow(np.rot90((reconstruction[:, :]), -1))
        plt.subplot(142)
        plt.imshow(np.rot90((speed_of_sound), -1))
        plt.subplot(143)
        plt.imshow(np.rot90(np.log10(initial_pressure), -1))
        plt.subplot(144)
        plt.imshow(np.rot90(segmentation, -1), vmin=values[0], vmax=values[-1], cmap=cmap)
        plt.show()
else:
    if len(shape) > 2:
        plt.figure()
        plt.subplot(241)
        plt.title("Fluence")
        plt.imshow(np.rot90((fluence[x_pos, :, :]), -1))
        plt.subplot(242)
        plt.title("Absorption")
        plt.imshow(np.rot90(np.log10(absorption[x_pos, :, :]), -1))
        plt.subplot(243)
        plt.title("Initial Pressure")
        plt.imshow(np.rot90(np.log10(initial_pressure[x_pos, :, :]), -1))
        plt.subplot(244)
        plt.title("Segmentation")
        plt.imshow(np.rot90(segmentation[x_pos, :, :], -1), vmin=values[0], vmax=values[-1],
        cbar = plt.colorbar(ticks=values)
        cbar.ax.set_yticklabels(names)
        plt.subplot(245)
        plt.imshow(np.rot90(fluence[:, y_pos, :], -1))
        plt.subplot(246)
        plt.imshow(np.rot90(np.log10(absorption[:, y_pos, :]), -1))
        plt.subplot(247)
        plt.imshow(np.rot90(np.log10(initial_pressure[:, y_pos, :]), -1))
        plt.subplot(248)
        plt.imshow(np.rot90(segmentation[:, y_pos, :], -1), vmin=values[0], vmax=values[-1],
        cbar = plt.colorbar(ticks=values)
        cbar.ax.set_yticklabels(names)
        plt.show()
    else:
        plt.figure()
        plt.subplot(141)
        plt.imshow(np.rot90(np.log10(fluence), -1))
        plt.subplot(142)
        plt.imshow(np.rot90(np.log10(absorption), -1))
        plt.subplot(143)
        plt.imshow(np.rot90(np.log10(initial_pressure), -1))
        plt.subplot(144)
        plt.imshow(np.rot90(segmentation, -1))
        plt.show()
```

Defining custom tissue structures and properties

The file can be found in simpa_examples/create_custom_tissues.py:

```
from simpa.utils import MolecularCompositionGenerator
from simpa.utils import MOLECULE_LIBRARY
from simpa.utils import Molecule
from simpa.utils import AbsorptionSpectrum
import numpy as np
```

```
def create_custom_absorber():
   wavelengths = np.linspace(200, 1500, 100)
    absorber = AbsorptionSpectrum(spectrum_name="random absorber",
                                  wavelengths=wavelengths,
                                  absorption_per_centimeter=np.random.random(
                                      np.shape(wavelengths)))
    return absorber
def create_custom_chromophore(volume_fraction: float = 1.0):
    chromophore = Molecule(
            spectrum=create_custom_absorber(),
            volume_fraction=volume_fraction,
            mus500=40.0,
            b mie=1.1,
            f_ray=0.9,
            anisotropy=0.9
    return chromophore
def create_custom_tissue_type():
    # First create an instance of a TissueSettingsGenerator
    tissue_settings_generator = MolecularCompositionGenerator()
    water_volume_fraction = 0.4
    bvf = 0.5
    oxy = 0.4
    # Then append chromophores that you want
    tissue_settings_generator.append(key="oxyhemoglobin", value=
                            MOLECULE_LIBRARY.oxyhemoglobin(oxy * bvf))
    tissue_settings_generator.append(key="deoxyhemoglobin", value=
                            MOLECULE_LIBRARY.deoxyhemoglobin(oxy * bvf))
    tissue_settings_generator.append(key="water", value=
                            MOLECULE_LIBRARY.water(water_volume_fraction))
    tissue_settings_generator.append(key="custom", value=
                            create_custom_chromophore(0.1))
    return_tissue_settings_generator.get_settings()
```

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