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**Spylce**

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## CONTENTS

The SPyIce package is a software tool that enables 1D finite difference simulation for vertical transport equations. It specifically focuses on thermal diffusion with the influence of salinity and physical properties. The package utilizes the Thomas tridiagonal solver as the solver algorithm. With SPyIce, users can model and analyze the behavior of temperature, salinity, and other relevant variables in a vertical system. It provides a comprehensive framework for studying the thermal diffusion process and its interaction with salinity in various scenarios. Hydra is used to automate the simulation runs of the Sea-Ice Model. It is used to manage and run sea ice simulations, making it easier for users to explore different scenarios and optimize their models.

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## 1.1 Introduction

Multiphysics simulation has been growing in the past decades due to computational capabilities. It involves translating the physical partial differential equations to an efficient numerical code. Implementing good Scientific software development practices help in reducing computational power. Testing applicability of multiphysical computational models can prove to be challenging due to thermodynamic and numerical constraints. Therefore, improving the code reproducibility through automation and object oriented concepts is beneficial.

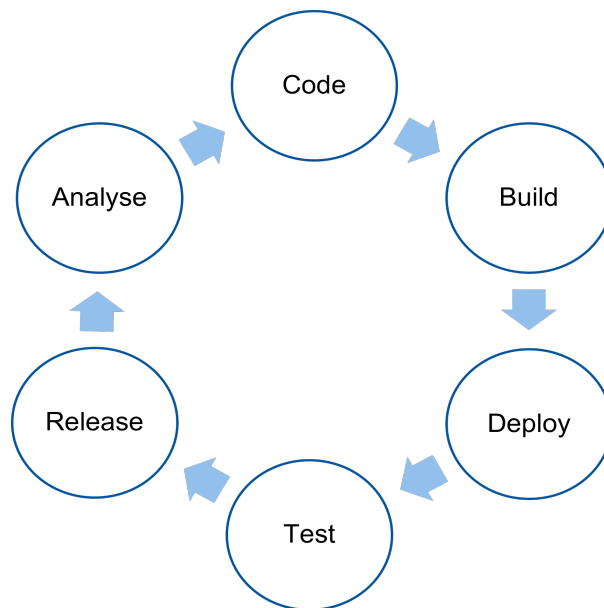


Fig. 1: Caption

This project aims to improve an existing 1D advection-diffusion sea ice simulation model us-

ing python to improve the benchmark testing for various model parameters. The automation of the model is performed using hydra python package. Hatch python project manager is used for testing, static analysis checks, and for creating reproducible build ecosystem. OOP concepts are leveraged and the mediator design pattern is implemented to improve sustainability of code.

## 1.2 Problem Statement

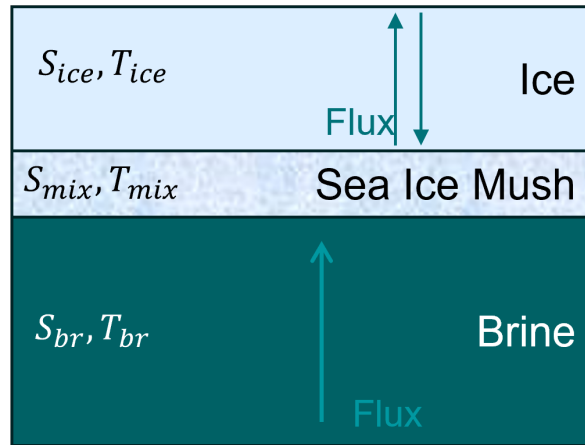


Fig. 2: caption

Modelling of Sea ice freezing involves ice-brine interface tracking which is modelling using the generalised Stefan problem. In this project, a 1D advective-diffusion model predominantly diffusive in nature is implemented using finite differences. The numerical equations models thermal and brine effects on the sea ice system with the following set of equations (??):

$$\begin{aligned} (\rho c)_{eff} \frac{\partial T}{\partial t} &= \frac{\partial}{\partial z} \left( k_{eff} \frac{\partial T}{\partial z} \right) - \rho_i L \frac{\partial \phi}{\partial t} \\ \phi \frac{\partial S_{br}}{\partial t} &= \frac{\partial}{\partial z} \left( D_{eff} \frac{\partial S_{br}}{\partial z} \right) - \frac{\rho_i}{\rho_{br}} P_S S_{br} \frac{\partial \phi}{\partial t} \end{aligned} \quad (1.1)$$

Brine in a sea ice system propagates along a vertical column and the flux across boundaries is considered to be constant (CHECK!!) as shown in Fig(). The Dirichlet boundary conditions (essential boundary condition) is implemented at the top layer of the vertical column and the system is assumed to be at melting temperature of sea ice as (??):.

$$T_m(S) = T_m - \Gamma S \quad (1.2)$$

The linear numerical model is solved using the *Thomas tri-diagonal solver* (REF) which is for a positive definite diagonally dominant matrix (REF). The conditional stability of the implicit numerical difference system is verified using the *Fourier stability criteria* (REF). The model parameters used in this project are number of iterations, sea ice freezing duration, and initial salinity.

**Stefan Problem:** So  $\lambda$  giving the minimum absolute result is determined and used to compute the location of the phase change.

$$-\frac{2sle^{-l^2}}{\sqrt{\pi} \operatorname{erf}(l)} - \frac{2se^{-2l^2}}{\pi \operatorname{erf}^2(l)} - 1 \quad (1.3)$$

## 1.3 Implementation

Hydra configuration tool is used to feed multiple input combinations to the Sea ice model and generate a structured output directory based on the given input combination. For example: all combinations of user defined input values for salinity, maximum iterations and time step size are sequentially run using hydra which improves the traceability of simulation along with logging files in their respective output directory. The inputs are initialised and processed by the Preprocessing class. Here the respective initial and boundary conditions are applied to the user-defined discrete finite difference mesh. The simulation model is executed for the given number of maximum iterations which allows to model for a time period of (time step size) \* (max iterations). At a given time  $t$ , the numerical model is solved until it attains a convergence for field parameters temperature, salinity and volumetric liquid fraction whose pseudo code is given below in code\_pseudo:

Listing 1: Python example

```
def run_model(self) -> None:
    """Runs the model using the provided configuration and output directory.
    ↪ """
    # apply boundary and initial conditions during the pre-processing stage
    ↪ and get the pre-processed dataclass
    preprocess_data, userinput_data = PreProcess.get_variables(
        self.config, self.out_dir_final
    )
    # run the sea ice model and get the results dataclass
    results_data = SeaIceModelClass.get_results(preprocess_data, userinput_
    ↪ data)
    # error analysis of results and get the analysis dataclass
    analysis_data = Analysis.get_error_results(
        t_k_diff=results_data.t_k_diff, t_stefan_diff=results_data.t_stefan_
    ↪ diff
    )
    # plot the sea ice model using the user input, results, and analysis
    ↪ dataclasses
    self.plot_model(userinput_data, results_data, analysis_data)
```

Once the field values are obtained, an error analysis is performed using Analytical class to verify discrepancies between numerical and analytical results. The analytical results are verified with the one-phase Stefan problem which keeps one of the two phases constant (liquid phase in this project) while modelling. The temperature field can be visualised using the Visualise-model class where the temperature fields can be compared at different spatial nodes points and their nodal time evolution in comparison to the analytical results.

## QUICKSTART

### 2.1 Installation Guide

#### 2.1.1 Prerequisites

Before installing Spyice, make sure you have the following prerequisites installed:

- **Python 3.11 or above:** Spyice requires Python 3.11 or above to run. You can download the latest version of Python from the [official website](#).
- **Hatch:** Hatch is a Python package manager that we'll use to create a virtual environment for Spyice. You can install Hatch by following the instructions on the [official website](#).
- **Sphinx:** Sphinx is a documentation generator that we'll use to build the Spyice documentation. You can install Sphinx using the appropriate package manager for your system by following the instructions on the [official website](#).

#### 2.1.2 Installation Steps

Follow these steps to install Spyice and set up the environment using hatch:

1. Clone this project repository to your local machine.
2. (Optional) The required wheels can be built with the help of *hatch* without the worries of cross-compilation and native architecture support.

```
` hatch build `
```

This command will build the wheels for the project and store them under the name *dist/spyice-1.0.0.dev0-py3-none-any.whl*

3. Create a the new default python virtual environment

```
` hatch env create `
```

The project will be automatically installed in editable mode by *hatch* when the environment is created. Confirm the installation by running *pip show spyice* in the shell.

4. You're all set! You can now start using Spyice.

#### 2.1.3 Viewing the documentation

Enter the project directory and run the following command in the terminal to view the pre-built static version of the sphinx documentation.

```
` python -m http.server --directory=docs/build/html/ `
```

This will start a local server at *http://localhost:8000/* where you can view the documentation by opening the link in your browser.



## 3.1 spyice package

### 3.1.1 Subpackages

spyice.models package

Submodules

src.spyice.models.advection\_diffusion

**class** src.spyice.models.advection\_diffusion.**AdvectionDiffusion**

Bases: object

Class representing an advection-diffusion model.

#### Parameters

- **argument** (*str*) – The argument for the advection-diffusion equation. Must be either “temperature” or “salinity”.
- **X** (*float*) – The value of X ( Field values T or S).
- **source** (*float*) – The source value.
- **X\_initial** (*float*) – The initial value of X (Field values T or S).
- **W** (*float*) – Liquid fraction current value.
- **W\_initial** (*float*) – The initial value of Liquid fraction.
- **w** (*float*) – The value of w convective upwind velocity.
- **dt** (*float*) – The time step size.
- **dz** (*float*) – The spatial step size.
- **nz** (*int*) – The number of spatial steps.
- **t\_passed** (*float*) – The time passed.
- **S\_IC** (*float*) – The initial condition for S.
- **Stefan** (*bool*, *optional*) – Whether to use Stefan condition. Defaults to False.
- **Buffo** (*bool*, *optional*) – Whether to use Buffo condition. Defaults to False.

- **bc\_neumann** (*float*, *optional*) – The Neumann boundary condition. Defaults to None.

**Raises**

- **AssertionError** – If the argument is not “temperature” or “salinity”.
- **AssertionError** – If both Stefan and Buffo conditions are set to True.

**Buffosolver**(*a*, *b*, *c*, *f*)

Solves a tridiagonal linear system using the Thomas algorithm. :param a: Lower diagonal of the tridiagonal matrix. :type a: np.ndarray :param b: Upper diagonal of the tridiagonal matrix. :type b: np.ndarray :param c: Main diagonal of the tridiagonal matrix. :type c: np.ndarray :param f: Right-hand side vector. :type f: np.ndarray

**Returns**

Solution vector of the linear system.

**Return type**

np.ndarray

**TDMA solver**(*a*, *b*, *c*, *d*)

Solves a tridiagonal matrix system using the TDMA (Thomas algorithm). :param a: Coefficients of the sub-diagonal elements. :type a: array-like :param b: Coefficients of the diagonal elements. :type b: array-like :param c: Coefficients of the super-diagonal elements. :type c: array-like :param d: Right-hand side vector. :type d: array-like

**Returns**

Solution vector.

**Return type**

array-like

**References**

- Tridiagonal matrix algorithm: [http://en.wikipedia.org/wiki/Tridiagonal\\_matrix\\_algorithm](http://en.wikipedia.org/wiki/Tridiagonal_matrix_algorithm)
- TDMA (Thomas algorithm): [http://www.cfd-online.com/Wiki/Tridiagonal\\_matrix\\_algorithm\\_-\\_TDMA\\_\(Thomas\\_algorithm\)](http://www.cfd-online.com/Wiki/Tridiagonal_matrix_algorithm_-_TDMA_(Thomas_algorithm))

**\_\_init\_\_**(*argument*, *X*, *source*, *X\_initial*, *W*, *W\_initial*, *w*, *dt*, *dz*, *nz*, *t\_passed*, *S\_IC*, *Stefan=False*, *Buffo=False*, *Voller=False*, *bc\_neumann=None*)

**Parameters**

- **argument** (*str*) – The argument for the advection-diffusion equation. Must be either “temperature” or “salinity”.
- **X** (*float*) – The value of X (Field values T or S).
- **source** (*float*) – The source value.
- **X\_initial** (*float*) – The initial value of X (Field values T or S).
- **W** (*float*) – Liquid fraction current value.
- **W\_initial** (*float*) – The initial value of Liquid fraction.

- **w** (*float*) – The value of *w* convective upwind velocity.
- **dt** (*float*) – The time step size.
- **dz** (*float*) – The spatial step size.
- **nz** (*int*) – The number of spatial steps.
- **t\_passed** (*float*) – The time passed.
- **S\_IC** (*float*) – The initial condition for *S*.
- **Stefan** (*bool*, *optional*) – Whether to use Stefan condition. Defaults to False.
- **Buffo** (*bool*, *optional*) – Whether to use Buffo condition. Defaults to False.
- **bc\_neumann** (*float*, *optional*) – The Neumann boundary condition. Defaults to None.

#### Raises

- **AssertionError** – If the argument is not “temperature” or “salinity”.
- **AssertionError** – If both Stefan and Buffo conditions are set to True.

#### **assemble\_tridiagonal()**

#### **factor\_1**(*argument*, *a*, *c*, *dt*, *dz*, *nz*)

Factor 1 and avoid zero division error

#### Parameters

- **argument** (*str*) – The argument value.
- **a** (*numpy.ndarray*) – The array containing values for ‘a’.
- **c** (*numpy.ndarray*) – The array containing values for ‘c’.
- **dt** (*float*) – The value of ‘dt’.
- **dz** (*float*) – The value of ‘dz’.
- **nz** (*int*) – The value of ‘nz’.

#### Returns

**The calculated factor(s) based on the given argument.**

If the argument is not “salinity”, returns a *numpy.ndarray*. If the argument is “salinity”, returns a list containing two *numpy.ndarrays*.

#### Return type

*numpy.ndarray* or list

#### **factor\_2**(*a*, *b*, *dt*, *dz*, *nz*)

Calculates the factor2 array for advection-diffusion model. :param a: Array containing values for ‘a’. :type a: *numpy.ndarray* :param b: Array containing values for ‘b’. :type b: *numpy.ndarray* :param dt: Time step. :type dt: *float* :param dz: Spatial step. :type dz: *float* :param nz: Number of elements in the array. :type nz: *int*

**Returns**

Array containing the calculated factor2 values.

**Return type**

numpy.ndarray

**factor\_3**(*a*, *d*, *nz*)

Factor 3 and avoid zero division error.

**Parameters**

- **a** (*numpy.ndarray*) – Array containing values.
- **d** (*numpy.ndarray*) – Array containing values.
- **nz** (*int*) – Number of elements.

**Returns**

Array containing factor 3 values.

**Return type**

numpy.ndarray

**modify\_tridiagonal\_voller\_scheme**()

**set\_up\_tridiagonal**()

Set up the tridiagonal matrix for Salinity and Temperature. This method sets up the tridiagonal matrix for Salinity and Temperature calculations. It initializes the main diagonal (main\_A), upper diagonal (upper\_A), and lower diagonal (lower\_A) of the matrix based on the given arguments. :param self: The object instance.

**Returns**

None

**unknowns\_matrix**(*temperature\_melt*, *non\_constant\_physical\_properties=False*)

Calculates the unknowns matrix for the advection-diffusion model. :returns:

**A tuple containing the following elements:**

- **X\_new** (ndarray): The solution vector for the unknowns.
- **X\_wind** (ndarray): The solution vector corrected for brine movement.
- **dt** (float): The time step used in the calculation.

**Return type**

tuple

**voller\_X\_array\_set\_zero\_to\_melt**(*X*, *X\_boundary*)

src.spyice.models.advection\_diffusion.top\_temp = 'Stefan'

$a * (dU/dt) + b * (dU/dz) + d/dz(c * dU/dz) + d * (dW/dt) = 0$

**Type**

Solves for X in Advection- Diffusion- Equation of the form

**src.spyice.models.sea\_ice\_model****class** src.spyice.models.sea\_ice\_model.**SeaIceModel**

Bases: object

SeaIceModelClass represents a class that models the behavior of sea ice.

**Parameters**

- **dataclass** (PreprocessData (page ??)) – The preprocessed data for the model.
- **user\_dataclass** (UserInput (page ??)) – The user input data for the model.

**\_\_init\_\_**(*dataclass*, *user\_dataclass*)**Parameters**

- **dataclass** (PreprocessData (page ??)) – The preprocessed data for the model.
- **user\_dataclass** (UserInput (page ??)) – The user input data for the model.

**add\_new\_parameter\_results**()**Return type**

None

**bc\_neumann**(*phi\_k*, *nz*, *bc\_condition=None*)

Apply Neumann boundary condition to the sea ice model.

**Parameters**

- **phi\_k** (*float*) – The value of phi at the k-th layer.
- **nz** (*int*) – The number of layers in the sea ice model.
- **bc\_condition** (*str*, *optional*) – The type of boundary condition to apply. Defaults to None.

**Returns**

None

**check\_and\_reset\_any\_iteration\_data**(*parameter\_list*, *parameter\_all\_list*)

Check the iteration data for temperature and phi values. This method appends the temperature and phi values from the current iteration to the respective arrays. The arrays are used to store the iteration data for further analysis.

**Parameters**

- **parameter\_list** (*list*) – The list of parameter values.
- **parameter\_all\_list** (*list*) – The list of all parameter values.

**Returns**

None

**Return type**

None

**check\_convergence**(*t, counter, t\_prev, s\_prev, phi\_prev, phi\_k, t\_k, s\_k, t\_initial, phi\_initial, s\_initial, vollen=False, \*\*kwargs*)

**choose\_phase\_type\_iteration**(*t*)

Choose the phase type iteration based on the one-phase and two-phase generalised Stefan Problem.

**Parameters**

**t** (*int*) – The time index.

**Returns**

**A tuple containing the following values:**

- **t\_stefan** (float): The Stefan temperature.
- **error\_depth\_t** (float): The error in depth.
- **depth\_stefan\_t** (float): The depth at time t.

**Return type**

tuple

**Raises**

**InvalidPhaseError** (page ??) – If the phase type is invalid (not 1 or 2).

**convergence\_loop\_iteration**(*t, t\_km1, s\_km1, phi\_km1, buffo=False, stefan=False, vollen=False, temp\_grad=None, salinity\_equation=False*)

Performs a single iteration of the convergence loop by first resetting the parameters to their previous time step values and then running the convergence loop until convergence is reached. :param t: Current temperature. :type t: float :param t\_km1: Temperature at the previous time step. :type t\_km1: float :param s\_km1: Salinity at the previous time step. :type s\_km1: float :param phi\_km1: Porosity at the previous time step. :type phi\_km1: float :param buffo: Flag indicating whether to use the buffo method. Defaults to False. :type buffo: bool, optional :param stefan: Flag indicating whether to use the Stefan method. Defaults to False. :type stefan: bool, optional :param temp\_grad: Temperature gradient. Defaults to None. :type temp\_grad: float, optional

**Methods called:**

- **reset\_iteration\_parameters**: Resets the iteration parameters for the sea ice model with values at previous time steps.
- **run\_while\_convergence\_iteration**: Runs the convergence loop until convergence is reached.

**Returns**

**A tuple containing the following values:**

- **t\_k** (float): Current temperature.
- **t\_prev** (float): Temperature at the previous time step.
- **s\_k** (float): Current salinity.
- **s\_prev** (float): Salinity at the previous time step.
- **phi\_k** (float): Current porosity.

- `phi_prev` (float): Porosity at the previous time step.
- `h_k` (float): Current heat flux.
- `h_solid` (float): Heat flux at the solid-liquid interface.
- `thickness` (float): Current thickness.
- `thickness_index` (int): Index of the thickness.
- `t_km1` (float): Temperature at the previous time step.
- `s_km1` (float): Salinity at the previous time step.
- `phi_km1` (float): Porosity at the previous time step.

**Return type**

tuple

**classmethod** `get_results(dataclass, user_dataclass)`

Runs the sea ice model and returns the results.

**Parameters**

- **`cls`** (*class*) – The class object.
- **`dataclass`** (PreprocessData (page ??)) – The dataclass containing preprocessed data.
- **`user_dataclass`** (UserInput (page ??)) – The dataclass containing user input.

**Returns**

The results dataclass object generated by running the sea ice model.

**Return type**

Results

**initialize\_state\_variables**(*t, t\_km1, s\_km1, phi\_km1*)

Initializes the state variables for the sea ice model.

**Parameters**

- **`t`** (*int*) – The current time step.
- **`t_km1`** (*float*) – The temperature at the previous time step.
- **`s_km1`** (*float*) – The salinity at the previous time step.
- **`phi_km1`** (*float*) – The liquid fraction at the previous time step.

**Returns****A tuple containing the initialized state variables:**

- `t_initial` (float): The initial temperature.
- `t_km1` (float): The temperature at the previous time step.
- `s_km1` (float): The salinity at the previous time step.
- `phi_initial` (float): The initial liquid fraction.
- `phi_km1` (float): The liquid fraction at the previous time step.

- **temp\_grad** (*float*): The temperature gradient.

**Return type**

tuple

**phi\_all\_mush\_list**(*phi\_k*, *phi\_all\_mush\_list*)

Calculates the number of elements in **phi\_k** that fall within the *mush\_lowerbound* and *mush\_upperbound* range.

**Parameters**

- **phi\_k** (*numpy.ndarray*) – The input array containing the values to be checked.
- **phi\_all\_mush\_list** (*list*) – The list to which the count of elements within the range will be appended.

**Returns**

The updated *phi\_all\_mush\_list* with the count of elements within the range appended.

**Return type**

list

**record\_iteration\_data**()

Records the iteration data for temperature and phi values. This method appends the temperature and phi values from the current iteration to the respective arrays. The arrays are used to store the iteration data for further analysis.

**Parameters**

**None**

**Returns**

**None**

**record\_mushy\_layer\_data**(*t*, *t\_km1*, *stefan*, *phi\_k*, *residual\_voller*, *s\_prev*)

Records the mushy layer data for temperature and phi values at time iterations *t* at specific time steps corresponding to initial stages, middle and final stages of the process.

**reset\_iteration\_parameters**(*t*, *tkm1*, *s\_km1*, *phi\_km1*)

Reset the iteration parameters for the sea ice model.

**Parameters**

- **t** (*float*) – Current temperature.
- **tkm1** (*float*) – Temperature at the previous time step.
- **s\_km1** (*float*) – Salinity at the previous time step.
- **phi\_km1** (*float*) – Liquid fraction at the previous time step.

**Returns**

**A tuple containing the following iteration parameters:**

- **t\_km1** (*float*): Temperature at the previous time step.
- **s\_km1** (*float*): Salinity at the previous time step.



- `phi_km1` (float): Liquid fraction at the previous time step.
- `temp_grad` (float): Temperature gradient.
- `t_err` (float): Temperature error.
- `s_err` (float): Salinity error.
- `phi_err` (float): Liquid fraction error.
- `t_initial` (float): Initial temperature.
- `s_initial` (float): Initial salinity.
- `phi_initial` (float): Initial liquid fraction.
- `source_term_array` (ndarray): Array of source-term values.
- `counter` (int): Iteration counter.

**Return type**

tuple

**run\_sea\_ice\_model()**

Runs the sea ice model.

This function iterates over a specified number of time steps and performs calculations to simulate the behavior of sea ice. It updates the results and saves a temperature profile plot at the end.

**Parameters**

None

**Returns**

None

**run\_while\_convergence\_iteration**(*t, t\_initial, s\_initial, phi\_initial, buffo, stefan, voller, t\_err, s\_err, phi\_err, source\_term, counter, \_is\_salinity\_equation=False*)

Runs the convergence loop until convergence is reached.

**Parameters**

- **t** – Time step
- **t\_km1** – Previous temperature array
- **s\_km1** – Previous salinity array
- **phi\_km1** – Previous phi array
- **buffo** – Flag for Buffo
- **stefan** – Flag for Stefan
- **t\_err** – Temperature error
- **s\_err** – Salinity errorhatch
- **phi\_err** – Phi error
- **t\_initial** – Initial temperature
- **phi\_initial** – Initial phi

- **source\_term** – RHS Source term of PDE equation
- **counter** – Iteration counter

**Methods called:**

- **initialise\_previous\_statevariables**: Defines the previous state variables for convergence iteration (temperature, salinity, liquid\_fraction).
- **update\_state\_variables**: Update the state variables (Enthalpy, Enthalpy Solid, Liquid Fraction, Temperature, Salinity).
- **locate\_ice\_ocean\_interface**: Locate the ice-ocean interface based on the liquid fraction.
- **overwrite\_statevariables**:
- **track\_mush\_for\_parameter**
- **phi\_all\_mush\_list**
- **check\_convergence**: computes error for convergence

**Returns**

**A tuple containing the following values:**

- **t\_km1** (float): Temperature at the previous time step.
- **s\_km1** (float): Salinity at the previous time step.
- **phi\_km1** (float): Liquid fraction at the previous time step.
- **t\_prev** (float): Temperature at the previous time step.
- **s\_prev** (float): Salinity at the previous time step.
- **phi\_prev** (float): Liquid fraction at the previous time step.
- **h\_k** (float): Current heat flux.
- **h\_solid** (float): Heat flux at the solid-liquid interface.
- **phi\_k** (float): Liquid fraction at the current time step.
- **t\_k** (float): Current temperature.
- **s\_k** (float): Current salinity.
- **thickness** (float): Current thickness.
- **thickness\_index** (int): Index of the thickness

**Return type**

tuple

**set\_boundary\_condition\_type(*critical\_depth*, *bc\_type*='Neumann')**

Sets the boundary condition type for the model. This method sets the boundary condition type for the model. It calculates the temperature gradient based on the given critical depth and boundary condition type. If the boundary condition type is “Neumann”, the temperature gradient is calculated using the formula:  $\text{temp\_grad} = dz * (\text{temperature\_melt} - \text{boundary\_top\_temperature}) / \text{critical\_depth}$ . If the boundary condition type is not “Neumann”, the temperature gradient is set to None.

**Parameters**

- **critical\_depth** (*float*) – The critical depth value.
- **bc\_type** (*str, optional*) – The type of boundary condition. Defaults to “Neumann”.

**Example**

```
model.set_boundary_condition_type(10.0, “Neumann”)
```

```
set_dataclass(_dataclass)
```

Sets the dataclass attributes of the object.

**Parameters**

**\_dataclass** – An instance of the dataclass.

**Returns**

None

```
t_running(fig, ax1, t_stefan, t_k, t_k_buffo=None, count=0)
```

Plot the temperature profile against depth.

**Parameters**

- **fig** (*matplotlib.figure.Figure*) – The figure object to plot on.
- **ax1** (*matplotlib.axes.Axes*) – The axes object to plot on.
- **t\_stefan** (*numpy.ndarray*) – The temperature profile obtained analytically.
- **t\_k** (*numpy.ndarray*) – The temperature profile obtained numerically.
- **t\_k\_buffo** (*numpy.ndarray, optional*) – The temperature profile obtained using Buffo method. Defaults to None.
- **count** (*int, optional*) – The count value. Defaults to 0.

**Returns**

None

```
track_mush_for_parameter(phi_k_, param, param_iterlist)
```

Track the mush for a given parameter.

**Parameters**

- **phi\_k** – numpy array representing the values of phi\_k
- **param** – numpy array representing the parameter values
- **param\_iterlist** – list to store the tracked mush values for the parameter
- **parameter** (*Updated list with the tracked mush values for the*)

```
src.spyice.models.sea_ice_model.locate_ice_ocean_interface(phi, dz, nz, **kwargs)
```

Locate ice ocean interface, based on liquid fraction equivalent ice thickness

**Parameters**

- **phi** (*array-like*) – Liquid fraction [-]
- **dz** (*float*) – Spatial discretization [m]
- **nz** (*int*) – Number of computational nodes
- **\*\*kwargs** – Additional keyword arguments Stefan (bool): Validation with Stefan problem (default: True)

**Returns**

A tuple containing:

- **if\_depth** (*float*): Location of the ice-water interface/sea ice total thickness [m]
- **if\_depth\_index** (*int*): Index of the ‘transition cell’ from ice to ocean (freezing) or water to ice (melting)

**Return type**

tuple

**src.spyice.models.stefan\_problem****class** src.spyice.models.stefan\_problem.**StefanProblem**

Bases: object

A class to solve the Stefan problem.

**static** **calculate\_temperature\_profile**(*depth\_stefan, t, dz, nz, ui*)

Calculates the temperature profile for the Stefan problem.

**Parameters**

- **depth\_stefan** (*float*) – The depth of the Stefan problem.
- **t** (*float*) – The time.
- **dz** (*float*) – The step size in the z-direction.
- **nz** (*int*) – The number of grid points in the z-direction.
- **ui** (UserInput (page ??)) – An instance of the UserInput class containing the required constants and parameters.

**Returns**

An array containing the temperature profile for the Stefan problem.

**Return type**

numpy.ndarray

**static** **calculate\_temperature\_twophase\_profiles**(*depth\_stefan, t, dz, nz, ui*)

Calculate the temperature and salinity profiles for the Stefan problem in a two-phase system.

**Parameters**

- **depth\_stefan** (*float*) – The depth of the Stefan problem.
- **t** (*float*) – The time.
- **dz** (*float*) – The grid spacing.

- **nz** (*int*) – The number of grid points.
- **ui** (UserInput (page ??)) – An instance of the UserInput class containing the input parameters.

**Returns**

A tuple containing the temperature and salinity profiles as numpy arrays.

**Return type**

tuple

**static stefan\_problem**(*t, ui*)

Calculates the Stefan problem solution for a given time and user input.

**Parameters**

- **t** (*float*) – The time value.
- **ui** (UserInput (page ??)) – An instance of the UserInput class containing the necessary input parameters.

**Returns**

The calculated Stefan problem solution.

**Return type**

float

**Raises**

**None** –

**static stefan\_problem\_twophase**(*t, ui*)

Solves the Stefan problem for a two-phase system.

**Parameters**

- **t** (*float*) – Time parameter.
- **ui** (UserInput (page ??)) – User input object containing boundary conditions and constants.

**Returns**

The result of the Stefan problem calculation.

**Return type**

float

**Module contents**

**spyice.parameters package**

**Submodules**

**src.spyice.parameters.constants**

**class** src.spyice.parameters.constants.**Constants**

Bases: Enum

Enumeration class for constants.

```
DEBUG = DebugConstants(phi_ini=1, phi_ini_Stefan=1, beta=1, kappa=1,
mu=1, Ra_c=0, alpha=1, k_i=1, k_br=1, k_w=1, D_s=0, c_br=1, c_i=1,
c_w=1, L=1, rho_i=1, rho_br=1, rho_w=1, m=1, g=0, phi_c=1, P_s=1,
a_phi=1, b_phi=1)
```

```
REAL = RealConstants(param_const=False, phi_ini=1, phi_ini_Stefan=1,
beta=0.0005836, kappa=1.37e-07, mu=0.00188, Ra_c=1.01,
alpha=0.0015600000000000002, L=334774, D_s=2e-09, k_i=2.0, k_br=0.6,
k_w=0.6, c_i=2000, c_br=3985, c_w=3985, rho_br=1028, rho_i=917,
rho_w=1028, m=2, g=9.8, phi_c=0.06823, P_s=0.01, a_phi=5.9e-06,
b_phi=169491.5254237288)
```

#### `src.spyice.parameters.debug_constants`

**class** `src.spyice.parameters.debug_constants.DebugConstants`

Bases: `object`

Class representing debug constants used in the model.

##### Parameters

- **phi\_ini** (*int*) – Freezing constant.
- **phi\_ini\_Stefan** (*int*) – Freezing constant for Stefan condition.
- **beta** (*int*) – Coefficient for density dependence on salinity.
- **kappa** (*int*) – Thermal diffusivity.
- **mu** (*int*) – Viscosity.
- **Ra\_c** (*int*) – Critical Rayleigh number.
- **alpha** (*int*) – Linear coefficient for Rayleigh number driven advection.
- **k\_i** (*int*) – Thermal conductivity of ice [W/m/K].
- **k\_br** (*int*) – Thermal conductivity of brine [W/m/K].
- **k\_w** (*int*) – Thermal conductivity of water.
- **D\_s** (*int*) – Diffusivity for salt.
- **c\_br** (*int*) – Specific heat of seawater (J/kg/K).
- **c\_i** (*int*) – Specific heat of ice.
- **c\_w** (*int*) – Specific heat of water.
- **L** (*int*) – Latent heat of fusion ice<->water (J/Kg).
- **rho\_i** (*int*) – Density of ice (Kg/m<sup>3</sup>).
- **rho\_br** (*int*) – Density of ocean used in volume averaging.
- **rho\_w** (*int*) – Density of water.
- **m** (*int*) – Cementation exponent for Archie’s equation.
- **g** (*int*) – Gravity constant.
- **phi\_c** (*int*) – Constant for phi.

- **P\_s** (*int*) – Constant for P\_s.
- **a\_phi** (*int*) – Constant for a\_phi.
- **b\_phi** (*int*) – Constant for b\_phi.

**\_\_init\_\_**(*phi\_ini=1, phi\_ini\_Stefan=1, beta=1, kappa=1, mu=1, Ra\_c=0, alpha=1, k\_i=1, k\_br=1, k\_w=1, D\_s=0, c\_br=1, c\_i=1, c\_w=1, L=1, rho\_i=1, rho\_br=1, rho\_w=1, m=1, g=0, phi\_c=1, P\_s=1, a\_phi=1, b\_phi=1*)

#### Parameters

- **phi\_ini** (*int*)
- **phi\_ini\_Stefan** (*int*)
- **beta** (*int*)
- **kappa** (*int*)
- **mu** (*int*)
- **Ra\_c** (*int*)
- **alpha** (*int*)
- **k\_i** (*int*)
- **k\_br** (*int*)
- **k\_w** (*int*)
- **D\_s** (*int*)
- **c\_br** (*int*)
- **c\_i** (*int*)
- **c\_w** (*int*)
- **L** (*int*)
- **rho\_i** (*int*)
- **rho\_br** (*int*)
- **rho\_w** (*int*)
- **m** (*int*)
- **g** (*int*)
- **phi\_c** (*int*)
- **P\_s** (*int*)
- **a\_phi** (*int*)
- **b\_phi** (*int*)

#### Return type

None

**D\_s:** *int*

```
L: int
P_s: int
Ra_c: int
a_phi: int
alpha: int
b_phi: int
beta: int
c_br: int
c_i: int
c_w: int
g: int
k_br: int
k_i: int
k_w: int
kappa: int
m: int
mu: int
phi_c: int
phi_ini: int
phi_ini_Stefan: int
rho_br: int
rho_i: int
rho_w: int
```

`src.spyice.parameters.real_constants`

**class** `src.spyice.parameters.real_constants.RealConstants`

Bases: `object`

Class representing real-valued constants used in the model.

Parameters

- **phi\_ini** (*int*) – Initial freezing value.
- **phi\_ini\_Stefan** (*int*) – Initial freezing value according to Stefan condition.



- **beta** (*float*) – Coefficient for density dependence on salinity.
- **kappa** (*float*) – Thermal diffusivity.
- **mu** (*float*) – Viscosity.
- **Ra\_c** (*float*) – Critical Rayleigh number.
- **alpha** (*float*) – Linear coefficient for Rayleigh number driven advection.
- **k\_i** (*float*) – Thermal conductivity of ice.
- **k\_br** (*float*) – Thermal conductivity of brine.
- **k\_w** (*float*) – Thermal conductivity of water.
- **D\_s** (*float*) – Diffusivity for salt.
- **c\_i** (*int*) – Specific heat of ice.
- **c\_br** (*int*) – Specific heat of seawater.
- **c\_w** (*int*) – Specific heat of water.
- **L** (*int*) – Latent heat of fusion between ice and water.
- **rho\_br** (*int*) – Density of ocean.
- **rho\_i** (*int*) – Density of ice.
- **rho\_w** (*int*) – Density of water.
- **m** (*int*) – Cementation exponent for Archie’s equation.
- **g** (*float*) – Earth gravity.
- **phi\_c** (*float*) – Critical porosity.
- **P\_s** (*float*) – Partition coefficient.
- **a\_phi** (*float*) – Coefficient a for porosity calculation.
- **b\_phi** (*float*) – Coefficient b for porosity calculation.
- **critical\_depth** (*float*) – Critical depth value.

```
__init__(param_const=False, phi_ini=1, phi_ini_Stefan=1, beta=0.0005836,
          kappa=1.37e-07, mu=0.00188, Ra_c=1.01, alpha=0.00156000000000000002,
          L=334774, D_s=2e-09, k_i=2.0, k_br=0.6, k_w=0.6, c_i=2000, c_br=3985,
          c_w=3985, rho_br=1028, rho_i=917, rho_w=1028, m=2, g=9.8,
          phi_c=0.06823, P_s=0.01, a_phi=5.9e-06, b_phi=169491.5254237288)
```

#### Parameters

- **param\_const** (*bool*)
- **phi\_ini** (*int*)
- **phi\_ini\_Stefan** (*int*)
- **beta** (*float*)
- **kappa** (*float*)
- **mu** (*float*)

- **Ra\_c** (*float*)
- **alpha** (*float*)
- **L** (*int*)
- **D\_s** (*float*)
- **k\_i** (*float*)
- **k\_br** (*float*)
- **k\_w** (*float*)
- **c\_i** (*int*)
- **c\_br** (*int*)
- **c\_w** (*int*)
- **rho\_br** (*int*)
- **rho\_i** (*int*)
- **rho\_w** (*int*)
- **m** (*int*)
- **g** (*float*)
- **phi\_c** (*float*)
- **P\_s** (*float*)
- **a\_phi** (*float*)
- **b\_phi** (*float*)

Return type

None

**D\_s:** float

**L:** int

**P\_s:** float

**Ra\_c:** float

**a\_phi:** float

**alpha:** float

**b\_phi:** float

**beta:** float

**c\_br:** int

**c\_i:** int

**c\_w:** int

**critical\_depth = 0.01**

```

g: float
k_br: float
k_i: float
k_w: float
kappa: float
m: int
mu: float
param_const: bool
phi_c: float
phi_ini: int
phi_ini_Stefan: int
rho_br: int
rho_i: int
rho_w: int

```

#### `src.spyice.parameters.results_params`

**class** `src.spyice.parameters.results_params.ResultsParams`

Bases: `object`

Class to store the results of the simulation.

##### Parameters

- **iter\_max** (*int*) – The maximum number of iterations.
- **nz** (*int*) – The number of depth levels.

##### Variables

- **error\_temperature** (*ndarray*) – An array of size `nz` to store the temperature errors.
- **error\_temperature\_sum** (*ndarray*) – An array of size `iter_max` to store the sum of temperature errors.
- **error\_temperature\_sum\_weighted** (*ndarray*) – An array of size `iter_max` to store the weighted sum of temperature errors.
- **temp\_grad** (*None*) – A placeholder for the temperature gradient.
- **thickness\_index\_total** (*ndarray*) – An array of size `iter_max` to store the total thickness index.
- **depth\_stefan\_all** (*ndarray*) – An array of size `iter_max` to store the depth Stefan values.

- **t\_stefan\_prev** (*ndarray*) – An array of size *nz* to store the previous Stefan temperature values.
- **t\_k\_prev** (*ndarray*) – An array of size *nz* to store the previous temperature values.
- **t\_stefan\_diff** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the differences in Stefan temperature values.
- **t\_k\_diff** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the differences in temperature values.
- **t\_stefan\_list** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the Stefan temperature values.
- **t\_k\_list** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the temperature values.
- **t\_k\_buffo\_list** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the buffered temperature values.
- **thickness\_list** (*ndarray*) – An array of size *iter\_max* to store the thickness values.
- **thickness\_list\_buffo** (*ndarray*) – An array of size *iter\_max* to store the buffered thickness values.
- **phi\_k\_list** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the phi values.
- **phi\_buffo\_list** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the buffered phi values.
- **s\_k\_list** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the *s* values.
- **s\_buffo\_list** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the buffered *s* values.
- **h\_k\_list** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the *h* values.
- **h\_solid\_list** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store the solid *h* values.
- **temp\_interface** (*ndarray*) – An array of size *iter\_max* to store the interface temperatures.
- **t\_k\_iter** (*list*) – A list to store the temperature values at each iteration.
- **phi\_k\_iter** (*list*) – A list to store the phi values at each iteration.
- **all\_phi\_iter** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store all phi values at each iteration.
- **t\_k\_iter\_all** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store all temperature values at each iteration.
- **phi\_k\_iter\_all** (*ndarray*) – A 2D array of size (*iter\_max*, *nz*) to store all phi values at each iteration.

- **all\_phi\_iter\_all** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all phi values at each iteration.
- **all\_t** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all temperature values.
- **all\_s** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all s values.
- **all\_phi** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all phi values.
- **all\_h** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all h values.
- **all\_h\_solid** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all solid h values.
- **all\_w** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all w values.
- **all\_thick** (*ndarray*) – An array of size iter\_max to store all thickness values.
- **all\_t\_passed** (*ndarray*) – An array of size iter\_max to store all passed temperature values.

**\_\_init\_\_**(*iter\_max, nz*)

#### Parameters

- **iter\_max** (*int*) – The maximum number of iterations.
- **nz** (*int*) – The number of depth levels.

#### Variables

- **error\_temperature** (*ndarray*) – An array of size nz to store the temperature errors.
- **error\_temperature\_sum** (*ndarray*) – An array of size iter\_max to store the sum of temperature errors.
- **error\_temperature\_sum\_weighted** (*ndarray*) – An array of size iter\_max to store the weighted sum of temperature errors.
- **temp\_grad** (*None*) – A placeholder for the temperature gradient.
- **thickness\_index\_total** (*ndarray*) – An array of size iter\_max to store the total thickness index.
- **depth\_stefan\_all** (*ndarray*) – An array of size iter\_max to store the depth Stefan values.
- **t\_stefan\_prev** (*ndarray*) – An array of size nz to store the previous Stefan temperature values.
- **t\_k\_prev** (*ndarray*) – An array of size nz to store the previous temperature values.
- **t\_stefan\_diff** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the differences in Stefan temperature values.

- **t\_k\_diff** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the differences in temperature values.
- **t\_stefan\_list** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the Stefan temperature values.
- **t\_k\_list** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the temperature values.
- **t\_k\_buffo\_list** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the buffered temperature values.
- **thickness\_list** (*ndarray*) – An array of size iter\_max to store the thickness values.
- **thickness\_list\_buffo** (*ndarray*) – An array of size iter\_max to store the buffered thickness values.
- **phi\_k\_list** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the phi values.
- **phi\_buffo\_list** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the buffered phi values.
- **s\_k\_list** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the s values.
- **s\_buffo\_list** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the buffered s values.
- **h\_k\_list** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the h values.
- **h\_solid\_list** (*ndarray*) – A 2D array of size (iter\_max, nz) to store the solid h values.
- **temp\_interface** (*ndarray*) – An array of size iter\_max to store the interface temperatures.
- **t\_k\_iter** (*list*) – A list to store the temperature values at each iteration.
- **phi\_k\_iter** (*list*) – A list to store the phi values at each iteration.
- **all\_phi\_iter** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all phi values at each iteration.
- **t\_k\_iter\_all** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all temperature values at each iteration.
- **phi\_k\_iter\_all** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all phi values at each iteration.
- **all\_phi\_iter\_all** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all phi values at each iteration.
- **all\_t** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all temperature values.
- **all\_s** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all s values.

- **all\_phi** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all phi values.
- **all\_h** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all h values.
- **all\_h\_solid** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all solid h values.
- **all\_w** (*ndarray*) – A 2D array of size (iter\_max, nz) to store all w values.
- **all\_thick** (*ndarray*) – An array of size iter\_max to store all thickness values.
- **all\_t\_passed** (*ndarray*) – An array of size iter\_max to store all passed temperature values.

**static store\_results**(*results\_dataclass, temp, s\_sw, phi, h, h\_solid, w, thickness, t\_passed, t*)

Stores the results of the simulation in the given results\_dataclass.

#### Parameters

- **results\_dataclass** – An instance of the ResultsDataClass where the results will be stored.
- **temp** – The temperature values to be stored.
- **s\_sw** – The s\_sw values to be stored.
- **phi** – The phi values to be stored.
- **h** – The h values to be stored.
- **h\_solid** – The h\_solid values to be stored.
- **w** – The w values to be stored.
- **thickness** – The thickness value to be stored.
- **t\_passed** – The t\_passed value to be stored.
- **t** – The t value to be stored.

#### Returns

The updated results\_dataclass with the stored results. - all\_t: The temperature values. - all\_s: The s\_sw values. - all\_phi: The phi values. - all\_h: The h values. - all\_h\_solid: The h\_solid values. - all\_w: The w values. - all\_thick: The thickness values. - all\_t\_passed: The t\_passed values.

**static store\_results\_for\_iter\_t**(*results\_dataclass, t, thickness\_index, t\_k, t\_stefan, s\_k, s\_k\_buffo, phi\_k, phi\_k\_buffo, h\_k, h\_solid, thickness, thickness\_buffo, thickness\_stefan, t\_k\_buffo, buffo=False*)

Stores the results for a given iteration 't' in the 'results\_dataclass'.

#### Parameters

- **results\_dataclass** – An instance of the results dataclass.

- **t** – The iteration number.
- **thickness\_index** – The index of the thickness.
- **t\_k** – The temperature values.
- **t\_stefan** – The Stefan temperature values.
- **s\_k** – The entropy values.
- **s\_k\_buffo** – The entropy values for buffo.
- **phi\_k** – The phase fraction values.
- **phi\_k\_buffo** – The phase fraction values for buffo.
- **h\_k** – The enthalpy values.
- **h\_solid** – The solid enthalpy values.
- **thickness** – The thickness values.
- **thickness\_buffo** – The thickness values for buffo.
- **thickness\_stefan** – The Stefan thickness values.
- **t\_k\_buffo** – The temperature values for buffo.
- **buffo** – A boolean indicating if buffo is enabled (default is False).

#### Returns

The updated 'results\_dataclass' with the stored results.

#### `src.spyice.parameters.user_input`

**class** `src.spyice.parameters.user_input.UserInput`

Bases: `object`

Represents the user input parameters for the model.

#### Variables

- **constants** (`RealConstants` (page ??) / `DebugConstants` (page ??)) – The type of constants to use.
- **max\_iterations** (`int`) – The maximum number of iterations.
- **is\_stefan** (`bool`) – Flag indicating whether Stefan condition is applied.
- **is\_buffo** (`bool`) – Flag indicating whether Buffo condition is applied.
- **liquidus\_relation\_type** (`str`) – The type of liquidus relation to use.
- **grid\_resolution\_dz** (`float`) – The grid resolution in the z-direction.
- **boundary\_condition\_type** (`str`) – The type of boundary condition to use.
- **temperature\_tolerance** (`float`) – The temperature tolerance.
- **salinity\_tolerance** (`float`) – The salinity tolerance.
- **liquid\_fraction\_tolerance** (`float`) – The liquid fraction tolerance.



- **initial\_temperature** (*str*) – The initial temperature profile.
- **initial\_salinity** (*str*) – The initial salinity profile.
- **initial\_liquid\_fraction** (*str*) – The initial liquid fraction profile.
- **output\_suffix** (*str*) – The suffix to be added to the output files.
- **temperature\_top\_type** (*str*) – The type of temperature condition at the top boundary.
- **phase\_type** (*int*) – The type of phase to consider.
- **grid\_timestep\_dt** (*float*) – The grid timestep.
- **dir\_output\_name** (*str*) – The name of the output directory.
- **critical\_liquid\_fraction** (*float*) – The critical liquid fraction.
- **boundary\_salinity** (*float*) – The boundary salinity (automatically calculated).
- **temperature\_melt** (*float*) – The temperature at which the material melts (automatically calculated).
- **boundary\_top\_temperature** (*float*) – The temperature at the top boundary (automatically calculated).
- **geometry\_type** (*int*) – The type of geometry.
- **counter\_limit** (*int*) – The counter limit.

#### **\_\_post\_init\_\_()**

Performs post-initialization tasks.

```
__init__(constants=RealConstants(param_const=False, phi_ini=1, phi_ini_Stefan=1,
    beta=0.0005836, kappa=1.37e-07, mu=0.00188, Ra_c=1.01,
    alpha=0.0015600000000000002, L=334774, D_s=2e-09, k_i=2.0, k_br=0.6,
    k_w=0.6, c_i=2000, c_br=3985, c_w=3985, rho_br=1028, rho_i=917,
    rho_w=1028, m=2, g=9.8, phi_c=0.06823, P_s=0.01, a_phi=5.9e-06,
    b_phi=169491.5254237288), config_data=<factory>, max_iterations=500,
    is_stefan=True, is_buffo=True, is_voller=False, is_salinity_equation=True,
    liquidus_relation_type='Normal', grid_resolution_dz=0.01,
    boundary_condition_type='Dirichlet', temperature_tolerance=0.01,
    salinity_tolerance=0.01, liquid_fraction_tolerance=0.001,
    initial_temperature='T(S)', initial_salinity='S34', initial_liquid_fraction='P1',
    output_suffix='nonconst_dens-mushfix', temperature_top_type='Stefan',
    phase_type=1, grid_timestep_dt=10,
    dir_output_name_hydra='Temperature_{S_IC}_{bc_condition}_{dz}_{dt}_{iter_max}_{cap_dens}',
    dir_output_name='Temperature_{S_IC}_{bc_condition}_{dz}_{dt}_{iter_max}_{cap_dens}',
    critical_liquid_fraction=0.1, counter_limit=100000)
```

#### Parameters

- **constants** (RealConstants (page ??) / DebugConstants (page ??))
- **config\_data** (DictConfig)
- **max\_iterations** (*int*)
- **is\_stefan** (*bool*)

- **is\_buffo** (*bool*)
- **is\_voller** (*bool*)
- **is\_salinity\_equation** (*bool*)
- **liquidus\_relation\_type** (*str*)
- **grid\_resolution\_dz** (*float*)
- **boundary\_condition\_type** (*str*)
- **temperature\_tolerance** (*float*)
- **salinity\_tolerance** (*float*)
- **liquid\_fraction\_tolerance** (*float*)
- **initial\_temperature** (*str*)
- **initial\_salinity** (*str*)
- **initial\_liquid\_fraction** (*str*)
- **output\_suffix** (*str*)
- **temperature\_top\_type** (*str*)
- **phase\_type** (*int*)
- **grid\_timestep\_dt** (*float*)
- **dir\_output\_name\_hydra** (*str*)
- **dir\_output\_name** (*str*)
- **critical\_liquid\_fraction** (*float*)
- **counter\_limit** (*int*)

Return type

None

**boundary\_condition\_type:** `str = 'Dirichlet'`

**boundary\_salinity:** `float`

**boundary\_top\_temperature:** `float`

**config\_data:** `DictConfig`

**constants:** *RealConstants* (page ??) | *DebugConstants* (page ??) =  
*RealConstants*(param\_const=False, phi\_ini=1, phi\_ini\_Stefan=1,  
beta=0.0005836, kappa=1.37e-07, mu=0.00188, Ra\_c=1.01,  
alpha=0.0015600000000000002, L=334774, D\_s=2e-09, k\_i=2.0, k\_br=0.6,  
k\_w=0.6, c\_i=2000, c\_br=3985, c\_w=3985, rho\_br=1028, rho\_i=917,  
rho\_w=1028, m=2, g=9.8, phi\_c=0.06823, P\_s=0.01, a\_phi=5.9e-06,  
b\_phi=169491.5254237288)

**counter\_limit:** `int = 100000`

**critical\_liquid\_fraction:** `float = 0.1`

```

dir_output_name: str =
'Temperature_{S_IC}_{bc_condition}_{dz}_{dt}_{iter_max}_{cap_dens}'

dir_output_name_hydra: str =
'Temperature_{S_IC}_{bc_condition}_{dz}_{dt}_{iter_max}_{cap_dens}'

geometry_type: int

grid_resolution_dz: float = 0.01

grid_timestep_dt: float = 10

initial_liquid_fraction: str = 'P1'

initial_salinity: str = 'S34'

initial_temperature: str = 'T(S)'

is_buffo: bool = True

is_salinity_equation: bool = True

is_stefan: bool = True

is_voller: bool = False

liquid_fraction_tolerance: float = 0.001

liquidus_relation_type: str = 'Normal'

max_iterations: int = 500

output_suffix: str = 'nonconst_dens-mushfix'

phase_type: int = 1

salinity_tolerance: float = 0.01

temperature_melt: float

temperature_tolerance: float = 0.01

temperature_top_type: str = 'Stefan'

```

`src.spyice.parameters.user_input.fourier_number_timestep(constants)`

Calculates the Fourier number for the given timestep.

#### Parameters

**constants** (*object*) – An object containing the required constants.

#### Returns

The calculated Fourier number.

#### Return type

float

## Module contents

`spyice.postprocess` package

## Submodules

`src.spyice.postprocess.analysis`

**class** `src.spyice.postprocess.analysis.Analysis`

Bases: `object`

Represents an analysis object that performs error analysis on temperature differences.

### Parameters

- **`t_k_diff`** (*float*) – The temperature difference in Kelvin.
- **`t_stefan_diff`** (*float*) – The temperature difference in Stefan-Boltzmann units.

**`__init__`**(*t\_k\_diff*, *t\_stefan\_diff*)

### Parameters

- **`t_k_diff`** (*float*) – The temperature difference in Kelvin.
- **`t_stefan_diff`** (*float*) – The temperature difference in Stefan-Boltzmann units.

**`calculate_errors`**(*field\_array*, *error\_norms\_object*)

Calculates the errors of a given field array using the provided error norms object.

### Parameters

- **`field_array`** (*numpy.ndarray*) – The field array to calculate the errors for.
- **`error_norms_object`** (`ErrorNorms` (page ??)) – The error norms object used to calculate the errors.

### Returns

A tuple containing the one-norm error, infinity-norm error, and two-norm error.

### Return type

tuple

**`error_analytical_numerical`**()

Calculates the errors of Numerical and Analytical using error norms one, two and infinity.

### Parameters

**`None`**

### Returns

**A tuple containing the errors calculated using different error norms:**

- **`T_k_Stefan_diff_L1norm`** (*float*): L1 norm of the difference between `T_k_list` and `T_Stefan_list`.

- **T\_k\_Stefan\_diff\_infnorm** (float): Infinity norm of the difference between **T\_k\_list** and **T\_Stefan\_list**.
- **T\_k\_Stefan\_diff\_L2norm** (float): L2 norm of the difference between **T\_k\_list** and **T\_Stefan\_list**.
- **T\_Stefan\_diff\_L1norm** (float): L1 norm of the difference between consecutive **T\_Stefan** values.
- **T\_Stefan\_diff\_infnorm** (float): Infinity norm of the difference between consecutive **T\_Stefan** values.
- **T\_Stefan\_diff\_L2norm** (float): L2 norm of the difference between consecutive **T\_Stefan** values.
- **T\_k\_diff\_infnorm** (float): Infinity norm of the difference between consecutive **T\_k** values.
- **T\_k\_diff\_L2norm** (float): L2 norm of the difference between consecutive **T\_k** values.
- **T\_k\_diff\_L1norm** (float): L1 norm of the difference between consecutive **T\_k** values.

**Return type**

tuple

**export\_residuals**(*residuals*, *temperature\_mushy*, *phi\_mushy*, *salinity\_mushy*, *output\_dir*)

Exports the residuals to a file.

**Parameters**

- **residuals** (*np.array*) – The residuals to export.
- **output\_dir** (*str*) – The output directory to save the residuals to.

**Returns**

None

**classmethod get\_error\_results**(*t\_k\_diff*, *t\_stefan\_diff*, *residual*, *temperature\_mushy*, *phi\_mushy*, *salinity\_mushy*, *output\_dir*)

Runs error analysis on the given temperature differences.

**Parameters**

- **cls** – The class object.
- **t\_k\_diff** (*array*) – The temperature difference for **k**.
- **t\_stefan\_diff** (*array*) – The temperature difference for **Stefan**.
- **residual** (*array*)
- **temperature\_mushy** (*array*)
- **phi\_mushy** (*array*)
- **salinity\_mushy** (*array*)
- **output\_dir** (*str*)

**Returns**

An instance of the `AnalysisData` class containing the error analysis results.

**Return type**

*AnalysisData* (page ??)

**set\_analysis()**

Sets up the analysis for the current object.

This method initializes the *error\_norms\_object* attribute with an instance of the *ErrorNorms* class, passing the *t\_k\_diff* and *t\_stefan\_diff* attributes as arguments. It then calculates the numerical analytical difference using the *numerical\_analytical\_diff* method of the *error\_norms\_object*.

**Parameters**

**None**

**Returns**

**None**

**static set\_dataclass(data\_to\_be\_converted, dataclass)**

Sets the values of a dataclass object using a dictionary.

**Parameters**

- **data\_to\_be\_converted** (*dict*) – A dictionary containing the values to be set.
- **dataclass** (*dataclass*) – The dataclass object to be updated.

**Returns**

The updated dataclass object.

**Return type**

*dataclass*

**store\_field\_errors(T\_k\_Stefan\_diff\_L1norm, T\_k\_Stefan\_diff\_infnorm, T\_k\_Stefan\_diff\_L2norm, T\_Stefan\_diff\_infnorm, T\_Stefan\_diff\_L2norm, T\_Stefan\_diff\_L1norm, T\_k\_diff\_infnorm, T\_k\_diff\_L2norm, T\_k\_diff\_L1norm)**

Stores the field errors.

**Parameters**

- **T\_k\_Stefan\_diff\_L1norm** (*float*) – The L1 norm of the temperature and concentration difference.
- **T\_k\_Stefan\_diff\_infnorm** (*float*) – The infinity norm of the temperature and concentration difference.
- **T\_k\_Stefan\_diff\_L2norm** (*float*) – The L2 norm of the temperature and concentration difference.
- **T\_Stefan\_diff\_infnorm** (*float*) – The infinity norm of the temperature difference.
- **T\_Stefan\_diff\_L2norm** (*float*) – The L2 norm of the temperature difference.

- **T\_Stefan\_diff\_L1norm** (*float*) – The L1 norm of the temperature difference.
- **T\_k\_diff\_infnorm** (*float*) – The infinity norm of the temperature and concentration difference.
- **T\_k\_diff\_L2norm** (*float*) – The L2 norm of the temperature and concentration difference.
- **T\_k\_diff\_L1norm** (*float*) – The L1 norm of the temperature and concentration difference.

**class** src.spyice.postprocess.analysis.**AnalysisData**

Bases: object

Represents the analysis data.

**Variables**

**all\_variables** (*dict*) – A dictionary containing all the variables.

**\_\_init\_\_** (*all\_variables*)

**Parameters**

**all\_variables** (*dict*)

**Return type**

None

**all\_variables:** dict

src.spyice.postprocess.visualise\_model

**class** src.spyice.postprocess.visualise\_model.**VisualiseModel**

Bases: object

A class for visualizing sea ice model results.

**Parameters**

- **user\_input\_dataclass** (UserInput (page ??)) – An instance of the UserInput class containing user input data.
- **results\_dataclass** (ResultsParams (page ??)) – An instance of the ResultsParams class containing results data.
- **error\_analysis\_dataclass** (Analysis (page ??)) – An instance of the Analysis class containing error analysis data.

**\_\_init\_\_** (*user\_input\_dataclass, results\_dataclass, error\_analysis\_dataclass*)

**Parameters**

- **user\_input\_dataclass** (UserInput (page ??)) – An instance of the UserInput class containing user input data.
- **results\_dataclass** (ResultsParams (page ??)) – An instance of the ResultsParams class containing results data.
- **error\_analysis\_dataclass** (Analysis (page ??)) – An instance of the Analysis class containing error analysis data.

**Return type**

None

**phi\_slope**(*iteration*)

Calculates the indices of the mushy regions based on the phi values.

**Parameters****iteration** (*int*) – The iteration number.**Returns**

The indices of the mushy regions.

**Return type**

numpy.ndarray

**plot\_H\_iter**(*h, t, s=None, param\_name='Temperature', unit='K', savefig=False*)**plot\_H\_iter\_all**(*savefig=False*)**plot\_H\_iter\_heatmap**(*h, p, temp\_all, d, t, param\_name='Temperature', unit='K', savefig=False*)**plot\_H\_iter\_heatmap\_mushy**(*h, p, temp\_all, t, param\_name='Temperature', unit='K', savefig=False*)**plot\_all\_phi\_mush**(*phi\_mush, t, savefig=False*)**plot\_depth\_over\_time**(*savefig=False*)

Plots the depth over time.

**Parameters****savefig** (*bool, optional*) – Whether to save the figure. Defaults to False.**plot\_depth\_over\_time\_heatmap**(*savefig=False*)

Plots the depth over time.

**Parameters****savefig** (*bool, optional*) – Whether to save the figure. Defaults to False.**plot\_error\_temp**(*zoom\_x, norm='inf', savefig=True*)

Plots the temperature errors using the specified norm.

**Parameters**

- **zoom\_x** (*int*) – The maximum value for the x-axis.
- **norm** (*str, optional*) – The norm to be used for plotting. Defaults to “inf”.
- **savefig** (*bool, optional*) – Whether to save the figure. Defaults to True.

**plot\_error\_temp\_diff**(*zoom\_x, savefig='True'*)

Plots the temperature differences between two consecutive iterations.

**Parameters**

- **zoom\_x** (*int*) – The maximum value for the x-axis.



- **savefig** (*str*, *optional*) – Indicates whether to save the figure or not. Defaults to “True”.

**plot\_liquidfraction\_heatmap**(*savefig=True*)

Plots the temperature heatmap.

Parameters

**savefig** (*bool*)

**plot\_response\_pt1\_pt2**(*tempmushPT1*, *tempmushPT2*, *savefig=False*)

**plot\_salinity\_heatmap**(*savefig=True*)

Plots the temperature heatmap.

Parameters

**savefig** (*bool*)

**plot\_temperature**(*z\_depth*, *savefig=True*, *Buffo\_matlab=False*)

Plots the temperature evolution at a given depth.

Parameters

- **z\_depth** (*float*) – The depth at which to plot the temperature evolution.
- **savefig** (*bool*, *optional*) – Whether to save the figure. Defaults to True.
- **Buffo\_matlab** (*bool*, *optional*) – Whether to include Buffo-matlab data in the plot. Defaults to False.

**plot\_temperature\_heatmap**(*savefig=True*)

Plots the temperature heatmap.

Parameters

**savefig** (*bool*)

**plot\_temperature\_heatmap\_as\_gif**( )

## Module contents

**spyice.preprocess** package

Submodules

**src.spyice.preprocess.geometry\_settings**

**class** **src.spyice.preprocess.geometry\_settings.GeometrySettings**

Bases: object

Set up model geometry with two parameters, an integer geom and a float dz.

Parameters

- **geom** (*int*) – 1 (is a test case scenario) or 2 (is according to W3 in Buffo et al. 2018)

- **dz** (*float*) – The parameter *dz* appears to be a float type. It is likely used to represent a specific value related to geometry or spatial calculations. If you need further assistance or have any specific questions about how to use this parameter in your code, feel free to ask!

`__init__(geom, dz)`

#### Parameters

- **geom** (*int*) – 1 (is a test case scenario) or 2 (is according to W3 in Buffo et al. 2018)
- **dz** (*float*) – The parameter *dz* appears to be a float type. It is likely used to represent a specific value related to geometry or spatial calculations. If you need further assistance or have any specific questions about how to use this parameter in your code, feel free to ask!

#### Return type

None

`src.spyice.preprocess.initial_boundary_conditions`

#### exception

`src.spyice.preprocess.initial_boundary_conditions.SalinityUnavailableError`

Bases: Exception

Exception raised when the S\_IC option is not available in initial conditions.

`src.spyice.preprocess.initial_boundary_conditions.boundary_condition(argument, t_passed, initial_salinity, **kwargs)`

Calculates the boundary conditions for temperature or salinity.

#### Parameters

- **argument** (*str*) – The argument specifying whether to calculate temperature or salinity.
- **t\_passed** (*float*) – The time passed.
- **initial\_salinity** (*float*) – The initial salinity value.
- **\*\*kwargs** – Additional keyword arguments.

#### Returns

A tuple containing the boundary conditions for temperature or salinity.

#### Return type

tuple

#### Raises

**None** –

`src.spyice.preprocess.initial_boundary_conditions.calculate_boundary_salinity(initial_salinity,`

Calculates the boundary salinity values based on the initial salinity.

**Parameters**

**initial\_salinity** (*str*) – The initial salinity value.

**Returns**

A tuple containing the bottom and top salinity values.

**Return type**

tuple

**Raises**

**SalinityException** – If the initial salinity value is not valid.

```
src.spyice.preprocess.initial_boundary_conditions.calculate_boundary_temperature(t_passed,
                                                                              ini-
                                                                              tial_salini
                                                                              kwargs)
```

Calculates the boundary temperature based on the given parameters.

**Parameters**

- **t\_passed** (*float*) – The time passed.
- **initial\_salinity** (*str*) – The initial salinity value.
- **kwargs** (*dict*) – Additional keyword arguments.

**Returns**

A tuple containing the bottom temperature and the top temperature.

**Return type**

tuple

```
src.spyice.preprocess.initial_boundary_conditions.compute_melting_temperature_from_salini
```

Computes the melting temperature from the given initial salinity.

**Parameters**

**initial\_salinity** (*float*) – The initial salinity value.

**Returns**

The computed melting temperature.

**Return type**

float

```
src.spyice.preprocess.initial_boundary_conditions.raise_salinity_exception(salinity_value)
```

Raises a custom exception if the salinity value is empty.

**Parameters**

**salinity\_value** (*str*) – The salinity value.

**Raises**

**SalinityUnavailableError** (page ??) – If the salinity value is empty.

```
src.spyice.preprocess.initial_boundary_conditions.set_boundary_temperature(t_passed,
                                                                              tem-
                                                                              per-
                                                                              a-
                                                                              ture_bottom,
                                                                              **kwargs)
```

Sets the boundary temperature based on the given parameters.

**Parameters**

- **t\_passed** (*float*) – The time passed.
- **temperature\_bottom** (*float*) – The bottom temperature.
- **\*\*kwargs** – Additional keyword arguments.

**Returns**

A tuple containing the top temperature and the bottom temperature.

**Return type**

tuple

```
src.spyice.preprocess.initial_boundary_conditions.set_inital_salinity(initial_salinity,  
                                                                    nz,  
                                                                    bound-  
                                                                    ary_salinity)
```

Sets the initial salinity values for each layer in the model.

**Parameters**

- **initial\_salinity** (*str*) – The type of initial salinity distribution.
- **nz** (*int*) – The number of layers in the model.
- **boundary\_salinity** (*float*) – The salinity value at the boundary.

**Returns**

An array of initial salinity values for each layer.

**Return type**

numpy.ndarray

**Raises**

**SalinityException** – If the initial\_salinity value is not recognized.

```
src.spyice.preprocess.initial_boundary_conditions.set_inital_temperature(initial_temperature,  
                                                                    nz,  
                                                                    bound-  
                                                                    ary_salinity,  
                                                                    bound-  
                                                                    ary_top_temperature)
```

Sets the initial temperature profile based on the given parameters.

**Parameters**

- **initial\_temperature** (*str*) – The type of initial temperature profile to set.
- **nz** (*int*) – The number of vertical grid points.
- **boundary\_salinity** (*float*) – The salinity at the boundary.
- **boundary\_top\_temperature** (*float*) – The temperature at the top boundary.

**Returns**

The initial temperature profile as a 1D numpy array.

**Return type**

numpy.ndarray

**Raises**

**None** –

```
src.spyice.preprocess.initial_boundary_conditions.set_initial_conditions(nz,
                                                                    bound-
                                                                    ary_salinity,
                                                                    ini-
                                                                    tial_temperature='T0',
                                                                    ini-
                                                                    tial_salinity='S1',
                                                                    ini-
                                                                    tial_liquid_fraction='P1',
                                                                    bound-
                                                                    ary_top_temperature=265.0)
```

Sets the initial conditions for the simulation.

**Parameters**

- **nz** (*int*) – Number of vertical grid points.
- **boundary\_salinity** (*float*) – Salinity value at the boundary.
- **initial\_temperature** (*str*, *optional*) – Initial temperature profile. Defaults to “T0”.
- **initial\_salinity** (*str*, *optional*) – Initial salinity profile. Defaults to “S1”.
- **initial\_liquid\_fraction** (*str*, *optional*) – Initial liquid fraction profile. Defaults to “P1”.
- **boundary\_top\_temperature** (*float*, *optional*) – Temperature value at the top boundary. Defaults to 265.0.

**Returns**

A tuple containing the temperature, salinity, liquid fraction, and upwind velocity arrays.

**Return type**

tuple

```
src.spyice.preprocess.initial_boundary_conditions.set_initial_liquidfraction(initial_liquid_f
                                                                    nz)
```

Sets the initial liquid fraction based on the given input.

**Parameters**

- **initial\_liquid\_fraction** (*str*) – The initial liquid fraction type.
- **nz** (*int*) – The number of grid points.

**Returns**

The array representing the initial liquid fraction.

**Return type**  
numpy.ndarray

**Raises**  
**None** –

### Examples

```
>>> set_initial_liquidfraction("P1", 10)
array([1., 1., 1., 1., 1., 1., 1., 1., 1., 1.])
>>> set_initial_liquidfraction("P_Stefan", 5)
array([0., 0., 0., 0., 0.])
>>> set_initial_liquidfraction("Po", 8)
array([0., 0., 0., 0., 0., 0., 0., 0., 0., 0.])
```

`src.spyice.preprocess.initial_boundary_conditions.t_w3(dt)`

Calculates the top boundary temperature and freeze date.

**Parameters**  
**dt** (*float*) – The time step.

**Returns**  
A tuple containing the top boundary temperature array and the freeze date.

**Return type**  
tuple

`src.spyice.preprocess.initial_boundary_conditions.temperature_gradient(phi, nz)`

Calculates the temperature gradient based on the given potential temperature profile and number of vertical levels.

**Parameters**

- **phi** (*list*) – The potential temperature profile.
- **nz** (*int*) – The number of vertical levels.

**Returns**  
The calculated temperature gradient.

**Return type**  
float

**Raises**  
**None** –

`src.spyice.preprocess.modify_initial_boundary`

**class** `src.spyice.preprocess.modify_initial_boundary.ModifyInitialBoundary`  
Bases: object

Defines functions for managing initial boundary conditions in a sea ice model.

**\_\_init\_\_()**

**Return type**

None

**bc\_neumann**(*phi\_k*, *nz*, *bc\_condition=None*)

Applies Neumann boundary condition to the given field.

**Parameters**

- **phi\_k** (*numpy.ndarray*) – The field to which the boundary condition is applied.
- **nz** (*int*) – The number of grid points in the vertical direction.
- **bc\_condition** (*str*, *optional*) – The type of boundary condition. Defaults to None.

**Returns**

None

**set\_boundary\_condition\_type**(*critical\_depth*, *bc\_type='Neumann'*)

Sets the boundary condition type for the given critical depth.

**Parameters**

- **critical\_depth** (*float*) – The critical depth value.
- **bc\_type** (*str*, *optional*) – The type of boundary condition. Defaults to “Neumann”.

**Raises**

**None** –

**Returns**

None

**src.spyice.preprocess.pre\_process**

**class** src.spyice.preprocess.pre\_process.**PreProcess**

Bases: *UserInput* (page ??), *GeometrySettings* (page ??), *ResultsParams* (page ??)

Class for preprocessing data before modeling.

Initialize the PreProcess class. :param config\_data: The configuration data object. :type config\_data: ConfigData :param output\_dir: The output directory. :type output\_dir: str

**Raises**

**None** –

**Returns**

None

**\_\_init\_\_**(*constants\_type*, *config\_data*, *output\_dir*)

Initialize the PreProcess class. :param config\_data: The configuration data object. :type config\_data: ConfigData :param output\_dir: The output directory. :type output\_dir: str

**Raises**

**None** –

**Returns**

None

**get\_userinput()**

Returns a `UserInput` object with the following attributes:

**Parameters**

**self** (`PreProcess` (page ??)) – The `PreProcess` instance.

**Returns**

**A `UserInput` object with the following attributes:**

- **constants**: The `constants` attribute.
- **grid\_timestep\_dt**: The `grid_timestep_dt` attribute.
- **initial\_salinity**: The `initial_salinity` attribute.
- **dir\_output\_name**: The `dir_output_name` attribute.
- **max\_iterations**: The `max_iterations` attribute.

**Return type**`UserInput` (page ??)**classmethod `get_variables(config, out_dir_final)`**

Retrieves variables and user input data after preprocessing.

**Parameters**

- **cls** – The class object.
- **config** – The configuration object.
- **out\_dir\_final** (*str*) – The output directory.

**Returns**

A tuple containing the filtered variables and user input data.

**Return type**`tuple[PreprocessData (page ??), UserInput (page ??)]`**preprocess()**

Preprocesses the data before running the simulation. This method sets up the initial conditions and boundary conditions for the simulation. It calculates the solid enthalpy and updates the enthalpy based on temperature, salinity, and liquid fraction. Finally, it initializes the ice thickness and prints a message indicating that the initial and boundary conditions have been applied.

**Args:**

None

**static `set_dataclass(data_to_be_converted, dataclass)`**

Sets the attributes of a dataclass object using a dictionary.

**Parameters**

- **data\_to\_be\_converted** (*dict*) – A dictionary containing the attribute names and values to be set.
- **dataclass** (*dataclass*) – The dataclass object to be modified.



**Returns**

The modified dataclass object with the attributes set.

**Return type**

dataclass

**set\_preprocess**(*config\_data*, *output\_dir*)

**class** `src.spyice.preprocess.pre_process.PreprocessData`

Bases: object

Class representing the preprocessing of data.

**Variables**

**is\_preprocessing** (*bool*) – Flag indicating if preprocessing is enabled.

**\_\_init\_\_**(*is\_preprocessing=True*)

**Parameters**

**is\_preprocessing** (*bool*)

**Return type**

None

**is\_preprocessing: bool = True**

`src.spyice.preprocess.pre_process.set_up_iter`(*iter\_max*, *grid\_timestep\_dt*)

Sets up the iteration parameters for the simulation. :param iter\_max: The maximum number of iterations. :type iter\_max: int :param grid\_timestep\_dt: The time step for the grid. :type grid\_timestep\_dt: float

**Returns**

Always returns 0.

**Return type**

int

`src.spyice.preprocess.store`

**Module contents**

`spyice.utils` package

**Submodules**

`src.spyice.utils.config_sort`

**class** `src.spyice.utils.config_sort.ConfigData`

Bases: object

Class representing configuration data.

**Parameters**

**setup** (*bool*) – Indicates if the setup is enabled or not.

**\_\_init\_\_**()

**Return type**

None

**setup = True**

**class** src.spylce.utils.config\_sort.**ConfigSort**

Bases: object

A class that provides methods for retrieving configuration parameters.

**Parameters**

**config** (*DictConfig*) – The configuration dictionary.

**Variables**

**config** (*DictConfig*) – The configuration dictionary.

**get\_config\_params(config**

*DictConfig*): Retrieves configuration parameters using the 'consts', 'dt', 'S\_IC', and 'iter\_max' keys.

**get\_ownconfig\_params(config)**

Retrieves configuration parameters using the 'constants', 'dt', 'S\_IC', and 'iter\_max' keys.

**getconfig\_dataclass(config, config\_type='default')**

Retrieves configuration parameters based on the specified config\_type.

**Parameters**

- **config** (*dataclass*)
- **config\_type** (*str*)

**Return type**

*dataclass*

**\_\_init\_\_**(*config*)

A class that provides methods for retrieving configuration parameters.

**Parameters**

**config** (*DictConfig*) – The configuration dictionary.

**Variables**

**config** (*DictConfig*) – The configuration dictionary.

**get\_config\_params(config**

*DictConfig*): Retrieves configuration parameters using the 'consts', 'dt', 'S\_IC', and 'iter\_max' keys.

**get\_ownconfig\_params(config)**

Retrieves configuration parameters using the 'constants', 'dt', 'S\_IC', and 'iter\_max' keys.

**getconfig\_dataclass(config, config\_type='default')**

Retrieves configuration parameters based on the specified config\_type.

**Parameters**

- **config** (*dataclass*)
- **config\_type** (*str*)

**Return type**

*dataclass*

**get\_config\_params(*config*)**

Get the configuration parameters from the given *config* dictionary.

**Parameters**

**config** (*DictConfig*) – The configuration dictionary.

**Returns**

None

**Raises**

**None** –

**get\_ownconfig\_params(*config*)**

Retrieves the parameters from the given configuration dictionary.

**Parameters**

**config** (*dict*) – The configuration dictionary.

**Returns**

None

**classmethod getconfig\_dataclass(*config*, *config\_type*='default')**

Retrieves configuration parameters based on the specified *config\_type*.

**Parameters**

- **config** (*dataclass*) – The configuration dictionary.
- **config\_type** (*str*) – The type of configuration “default” or “jupyter”. “jupyter” is used for Jupyter notebook configurations.

**Returns**

An instance of the ConfigData class.

**Return type**

*ConfigData* (page ??)

**src.spyice.utils.config\_sort.read\_omegaconfig(*config*, *parameter*)**

Reads the parameter from the OmegaConf configuration.

**Parameters**

**parameter** (*str*) – The parameter to read.

**Returns**

The value of the parameter.

**Return type**

str

**Raises**

**None** –

**src.spyice.utils.error\_norms****class src.spyice.utils.error\_norms.ErrorNorms**

Bases: object

Defines functions for calculating error norms between numerical and analytical values.

**Parameters**

- **numerical\_values** (*list*) – A list of numerical values.
- **analytical\_values** (*list*) – A list of analytical values.

**Variables**

- **numerical\_values** (*list*) – A list of numerical values.
- **analytical\_values** (*list*) – A list of analytical values.
- **iteration\_count** (*int*) – The count of iterations.

**\_\_init\_\_**(*numerical\_values, analytical\_values*)

**Parameters**

- **numerical\_values** (*list*) – A list of numerical values.
- **analytical\_values** (*list*) – A list of analytical values.

**Variables**

- **numerical\_values** (*list*) – A list of numerical values.
- **analytical\_values** (*list*) – A list of analytical values.
- **iteration\_count** (*int*) – The count of iterations.

**Return type**

None

**infinity\_norm**(*numerical\_analytical\_diff*)

Calculates the infinity norm of the given numerical-analytical difference.

**Parameters**

**numerical\_analytical\_diff** (*array-like*) – The numerical-analytical difference.

**Returns**

The infinity norm of the difference. If the analytical values are 1-dimensional, a single float value is returned. Otherwise, a numpy array containing the infinity norm for each iteration is returned.

**Return type**

float or numpy.ndarray

**numerical\_analytical\_diff**()

Calculates the absolute difference between the analytical values and the numerical values.

**Parameters**

**None**

**Returns**

The array containing the absolute differences between the analytical values and the numerical values.

**Return type**

numpy.ndarray

**one\_norm**(*numerical\_analytical\_diff*)

Calculates the one norm of the numerical-analytical difference.

**Parameters**

**numerical\_analytical\_diff** (*ndarray*) – The numerical-analytical difference.

**Returns**

The one norm of the numerical-analytical difference. If the analytical values are 1D, a float is returned. Otherwise, an ndarray is returned with the one norm for each iteration.

**Return type**

ndarray or float

**Raises**

**None** –

**two\_norm**(*numerical\_analytical\_diff*)

Calculates the two-norm error of the numerical-analytical difference.

**Parameters**

**numerical\_analytical\_diff** (*ndarray*) – The numerical-analytical difference.

**Returns**

The two-norm error. If the analytical values are 1D, a float is returned. If the analytical values are 2D, an ndarray is returned with the two-norm error for each iteration.

**Return type**

ndarray or float

### src.spyice.utils.helpers

src.spyice.utils.helpers.**export\_residuals**(*self, residuals, temperature\_mushy, phi\_mushy, salinity\_mushy, output\_dir*)

Exports the residuals to a file.

**Parameters**

- **residuals** (*np.array*) – The residuals to export.
- **output\_dir** (*str*) – The output directory to save the residuals to.

**Returns**

None

src.spyice.utils.helpers.**set\_dataclass**(*data\_to\_be\_converted, dataclass*)

Sets the values of a dataclass object using a dictionary.

**Parameters**

- **data\_to\_be\_converted** (*dict*) – The dictionary containing the data to be converted.
- **dataclass** (*dataclass*) – The dataclass object to be updated.

**Returns**

The updated dataclass object.

**Return type**

dataclass

`src.spyice.utils.helpers.t_total(t_passed, dt)`

Computes total time passed based on current time step `dt` and, total time of previous time step.

**Parameters**

- **t\_passed** (*float*) – The time that has already passed.
- **dt** (*float*) – The time increment.

**Returns**

The total time.

**Return type**

float

**src.spyice.utils.spyice\_exceptions module**

**exception** `src.spyice.utils.spyice_exceptions.ConvergenceError`

Bases: Exception

Exception raised when convergence is not reached.

`__init__(message='Convergence not reached.')`

**exception** `src.spyice.utils.spyice_exceptions.InvalidPhaseError`

Bases: Exception

Custom exception class for SeaIceModel errors.

**src.spyice.utils.spyice\_logger**

**class** `src.spyice.utils.spyice_logger.SpyiceLogger`

Bases: StringIO

A class that replaces the print function with a logger object.

**Parameters**

- **logger** – The logger object used for logging.
- **\*args** – Additional positional arguments.
- **\*\*kwargs** – Additional keyword arguments.

`__init__(logger, *args, **kwargs)`

**Parameters**

- **logger** – The logger object used for logging.
- **\*args** – Additional positional arguments.
- **\*\*kwargs** – Additional keyword arguments.

**flush()**

Flush write buffers, if applicable.

This is not implemented for read-only and non-blocking streams.

**write(message)**

Writes the given message to the logger if it is not empty.

**Parameters**

**message** (*str*) – The message to be written to the logger.

**Returns**

None

**Module contents****3.1.2 Submodules****src.spyice.coefficients**

**src.spyice.coefficients.update\_coefficients**(*argument, x\_initial, w, phi, nz, salinity\_initial*)

Updates of coefficients required to solve the Advection Reaction Diffusion Equation for each time step for temperature or salinity

**Parameters**

- **argument** (*str*) – Either ‘temperature’ or ‘salinity’
- **x\_initial** (*float*) – Initial value for salinity or temperature [ppt] or [K]
- **w** (*float*) – Brine velocity [ms-1]
- **phi** (*float*) – Liquid fraction [-]
- **nz** (*int*) – Number of computational nodes

**Returns**

**A tuple containing the following coefficients:**

- **a** (numpy.ndarray): ‘temperature’: heat capacity | ‘salinity’: liquid fraction
- **b** (numpy.ndarray): Brine velocity [ms-1]
- **c** (numpy.ndarray): ‘temperature’: thermal conductivity | ‘salinity’: salt diffusivity
- **d** (numpy.ndarray): ‘temperature’: latent heat | ‘salinity’: factor to determine salinity increase due to liquid fraction decrease

**Return type**

tuple

**src.spyice.main\_process****class** src.spyice.main\_process.**MainProcess**

Bases: object

Main class to run the model.

**Parameters**

- **config** – The configuration object.
- **hyd\_output\_dir** (*Path* / *str*) – The directory path for the hydraulic output. Defaults to the ‘outputs’ directory in the current working directory.
- **project\_path** (*Path* / *str*) – The project path. Defaults to the current working directory.

```
__init__(config,  
         hyd_output_dir=WindowsPath('C:/Users/sneha/Documents/MBDHiWi/MBDHiwi/spyicedir/spyice'),  
         project_path='C:\\Users\\sneha\\Documents\\MBDHiWi\\MBDHiwi\\spyicedir\\spyicedir\\')
```

**Parameters**

- **config** – The configuration object.
- **hyd\_output\_dir** (*Path* / *str*) – The directory path for the hydraulic output. Defaults to the ‘outputs’ directory in the current working directory.
- **project\_path** (*Path* / *str*) – The project path. Defaults to the current working directory.

**plot\_model**(*userinput\_data*, *results\_data*, *analysis\_data*)

Plots various visualizations of the model.

**Parameters**

- **userinput\_data** (*UserInputData*) – The user input data.
- **results\_data** (*ResultsData*) – The results data.
- **analysis\_data** (*AnalysisData* (page ??)) – The error analysis data.

**Returns**

None

**Raises****None** –**run\_model**()

Runs the model using the provided configuration and output directory.

**Parameters****None****Returns**

None

**Return type**

None



**src.spyice.rhs**

**src.spyice.rhs.apply\_boundary\_condition**(*argument, x\_initial, source, factor1, factor3, a, delta\_upwind, w, nz, t\_passed, salinity\_initial, \_temperature\_top, is\_stefan, is\_buffo=False, is\_voller=False, bc\_neumann=None*)

Creates the right hand side of the matrix equation considering source terms.

**Parameters**

- **argument** (*str*) – Either ‘salinity’ for salt equation or ‘temperature’ for temperature equation.
- **x\_initial** (*float*) – Value of X at the last time step.
- **source** (*float*) – Source term.
- **factor1** (*float*) – Factor 1.
- **factor3** (*float*) – Factor 3.
- **a** (*float*) – A parameter.
- **delta\_upwind** (*float*) – Difference of ice volume fraction between this and the last time step.
- **w** (*float*) – W parameter.
- **nz** (*int*) – Number of computational nodes.
- **t\_passed** (*float*) – Time passed in seconds.
- **salinity\_initial** (*float*) – Initial salinity value.
- **\_temperature\_top** (*float*) – Top temperature value.
- **is\_stefan** (*bool*) – Indicates if Stefan condition is used.
- **is\_buffo** (*bool, optional*) – Indicates if Buffo condition is used. Defaults to False.
- **bc\_neumann** (*float, optional*) – Neumann boundary condition. Defaults to None.
- **float** – The right hand side of the equation.

**src.spyice.rhs.correct\_for\_brine\_movement**(*argument, x\_initial, w, t\_passed, nz, salinity\_initial, top\_temp*)

Corrects for brine movement based on the given arguments.

**Parameters**

- **argument** (*str*) – The argument for correction, either “salinity” or “temperature”.
- **x\_initial** (*numpy.ndarray*) – The initial values of x.
- **w** (*numpy.ndarray*) – The values of w.
- **t\_passed** (*float*) – The time passed.
- **nz** (*int*) – The number of elements.
- **salinity\_initial** (*float*) – The initial salinity value.

- **top\_temp** (*float*) – The top temperature value.

**Returns**

The corrected values of x.

**Return type**

numpy.ndarray

**Raises**

**None** –

**src.spyice.statevariables**

```
src.spyice.statevariables.compute_error_for_convergence(temperature_calculated,  
                                                         temperature_previous,  
                                                         salinity_calculated,  
                                                         salinity_previous, liq-  
                                                         uid_fraction_calculated,  
                                                         liquid_fraction_previous,  
                                                         voller=False, **kwargs)
```

Computes the errors for convergence between the calculated and previous values of temperature, salinity, and liquid fraction.

**Parameters**

- **temperature\_calculated** (*numpy.ndarray*) – Array of calculated temperature values.
- **temperature\_previous** (*numpy.ndarray*) – Array of previous temperature values.
- **salinity\_calculated** (*numpy.ndarray*) – Array of calculated salinity values.
- **salinity\_previous** (*numpy.ndarray*) – Array of previous salinity values.
- **liquid\_fraction\_calculated** (*numpy.ndarray*) – Array of calculated liquid fraction values.
- **liquid\_fraction\_previous** (*numpy.ndarray*) – Array of previous liquid fraction values.

**Returns**

**A tuple containing the following error values:**

- **temperature\_error\_max** (*float*): Maximum temperature error for convergence check.
- **temperature\_error\_all** (*numpy.ndarray*): Full temperature error for convergence check.
- **salinity\_error\_max** (*float*): Maximum salinity error for convergence check.
- **salinity\_err\_all** (*numpy.ndarray*): Full salinity error for convergence check.

- `liquid_fraction_error_max` (float): Maximum liquid fraction error for convergence check.
- `liquid_fraction_error_all` (numpy.ndarray): Full liquid fraction error for convergence check.

#### Return type

tuple

```
src.spyice.statevariables.overwrite_statevariables(temperature_calculated,
                                                    salinity_calculated,
                                                    liquid_fraction_calculated,
                                                    t_k_lhs_A_matrix=None,
                                                    temp_factor3=None)
```

Overwrites the state variables with the calculated values. :param `temperature_calculated`: The calculated temperature. :type `temperature_calculated`: float  
:param `salinity_calculated`: The calculated salinity. :type `salinity_calculated`: float  
:param `liquid_fraction_calculated`: The calculated liquid fraction. :type `liquid_fraction_calculated`: float

#### Returns

A tuple containing the following state variables:

- `temperature_new` (float): The current temperature for iteration.
- `salinity_new` (float): The current salinity for iteration.
- `liquid_fraction_new` (float): The current liquid fraction for iteration.

#### Return type

tuple

```
src.spyice.statevariables.reset_error_for_while_loop(temperature_tolerance,
                                                    salinity_tolerance,
                                                    liquid_fraction_tolerance)
```

Resets the error values for a while loop based on the given tolerances.

#### Parameters

- **`temperature_tolerance`** (*float*) – The tolerance for temperature error.
- **`salinity_tolerance`** (*float*) – The tolerance for salinity error.
- **`liquid_fraction_tolerance`** (*float*) – The tolerance for liquid fraction error.

#### Returns

A tuple containing the reset error values for temperature, salinity, and liquid fraction.

#### Return type

tuple

```
src.spyice.statevariables.set_initial_statevariables(temperature, salinity,
                                                    liquid_fraction)
```

Initializes the state variables for the given temperature, salinity, and liquid fraction.

**Parameters**

- **temperature** (*float*) – The initial temperature.
- **salinity** (*float*) – The initial salinity.
- **liquid\_fraction** (*float*) – The initial liquid fraction.

**Returns**

**A tuple containing the initialized state variables:**

- `temperature_initial` (*float*): The initial temperature.
- `temperature_new` (*float*): The current temperature for iteration.
- `temperature_previous` (*float*): The previous temperature for initialization.
- `salinity_initial` (*float*): The initial salinity.
- `salinity_new` (*float*): The current salinity for iteration.
- `salinity_previous` (*float*): The previous salinity for initialization.
- `phi_initial` (*float*): The initial liquid fraction.
- `liquid_fraction_new` (*float*): The current liquid fraction for iteration.
- `phi_prev` (*float*): The previous liquid fraction for initialization.

**Return type**

tuple

```
src.spyice.statevariables.set_statevariables(temperature_calculated,  
                                             salinity_calculated,  
                                             liquid_fraction_calculated)
```

Set the state variables for temperature, salinity, and liquid fraction.

**Parameters**

- **temperature\_calculated** (*float*) – The calculated temperature.
- **salinity\_calculated** (*float*) – The calculated salinity.
- **liquid\_fraction\_calculated** (*float*) – The calculated liquid fraction.

**Returns**

**A tuple containing the following state variables:**

- `temperature_initial` (*float*): The initial temperature.
- `temperature_previous` (*float*): The previous temperature.
- `salinity_initial` (*float*): The initial salinity.
- `salinity_previous` (*float*): The previous salinity.
- `phi_initial` (*float*): The initial liquid fraction.
- `phi_prev` (*float*): The previous liquid fraction.

**Return type**

tuple

```
src.spyice.statevariables.voller_residual_scheme(temperature_calculated,
                                                temperature_previous,
                                                liquid_fraction_calculated,
                                                liquid_fraction_previous, kwargs)
```

Calculate the residual sum for the Voller residual scheme. This function computes the residual sum based on the calculated and previous temperature and liquid fraction values, along with additional parameters provided in the kwargs dictionary. The residual sum is used to assess the convergence of the numerical scheme. (RHS - LHS matrix)  
Parameters: ——— temperature\_calculated : array-like

The calculated temperature values at the current iteration.

**temperature\_previous**

[array-like] The temperature values from the previous iteration.

**liquid\_fraction\_calculated**

[array-like] The calculated liquid fraction values at the current iteration.

**liquid\_fraction\_previous**

[array-like] The liquid fraction values from the previous iteration.

**kwargs**

[dict] A dictionary containing additional parameters required for the computation. Expected keys are: - "A\_matrix": The coefficient matrix for the temperature calculation. - "phi\_initial": The initial liquid fraction values. - "t\_initial": The initial temperature values. - "temp\_factor3": A factor used in the residual calculation.

**Returns:****residual\_sum**

[float] The sum of the absolute values of the residuals.

**src.spyice.update\_physical\_values**

```
src.spyice.update_physical_values.H_function(_self, _temperature, _enthalpy_s1)
```

Calculates the value of H function.

**Parameters**

- **\_self** (*float*) – The value of self.
- **\_temperature** (*float*) – The temperature.
- **\_enthalpy\_s1** (*float*) – The enthalpy.

**Returns**

The calculated value of H function.

**Return type**

float

```
src.spyice.update_physical_values.H_function_derivate(_x, _enthalpy_s1)
```

Calculates the derivative of the H function.

**Parameters**

- **\_x** (*float*) – The value of x.
- **\_enthalpy\_s1** (*float*) – The value of enthalpy\_s1.

**Returns**

The derivative of the H function.

**Return type**

float

`src.spyice.update_physical_values.H_newton_iteration(_temperature, _enthalpy_s1)`

Performs Newton iteration to find the root of the H\_function.

**Parameters**

- **\_temperature** (*float*) – The temperature value.
- **\_enthalpy\_s1** (*float*) – The enthalpy value.

**Returns**

The root of the H\_function.

**Return type**

float

**Raises**

**None** –

`src.spyice.update_physical_values.calculate_liquidus_temperature(_salinity)`

`src.spyice.update_physical_values.calculate_melting_temperature_from_salinity(_salinity, _temperature_melt=27, _liquid_relation=`

Calculates the melting temperature of seawater based on salinity.

**Parameters**

- **\_salinity** (*numpy.ndarray*) – Array of salinity values.
- **\_temperature\_melt** (*float, optional*) – Melting temperature. Defaults to `_temperature_melt`.
- **\_liquid\_relation** (*str, optional*) – Liquid relation type. Must be either “Normal” or “Frezchem”. Defaults to “Normal”.

**Returns**

Array of melting temperature values.

**Return type**

*numpy.ndarray*

**Raises**

**TypeError** – If `_liquid_relation` is not “Normal” or “Frezchem”.

`src.spyice.update_physical_values.physical_control_for_infinite_values(_phi)`

Calculates the control values for infinite phi values.

**Parameters**

**\_phi** (*numpy.ndarray*) – The input array of phi values.

**Returns**

The control values for the given phi values.

**Return type**

*numpy.ndarray*

```
src.spyice.update_physical_values.phi_func(_enthalpy_k1, _enthalpy_s1)
```

Calculates the phi value based on the given enthalpy values.

**Parameters**

- **\_enthalpy\_k1** (*float*) – The enthalpy value for k1.
- **\_enthalpy\_s1** (*float*) – The enthalpy value for s1.

**Returns**

The calculated phi value.

**Return type**

*float*

```
src.spyice.update_physical_values.update_enthalpy(_temperature, _salinity,  
                                                    _liquid_fraction, _nz,  
                                                    _method='likebuffo')
```

Updates the enthalpy based on the given parameters.

**Parameters**

- **\_temperature** (*float*) – The temperature value.
- **\_salinity** (*float*) – The salinity value.
- **\_liquid\_fraction** (*float*) – The liquid fraction value.
- **\_nz** (*int*) – The nz value.
- **\_method** (*str*, *optional*) – The method used for calculating enthalpy. Defaults to “likebuffo”.

**Returns**

The updated enthalpy value.

**Return type**

*float*

**Raises**

**TypeError** – If the given method is not available.

```
src.spyice.update_physical_values.update_enthalpy_solid_state(_salinity, _nz,  
                                                                _liq_rel='Normal',  
                                                                _tempera-  
                                                                ture_melt=271.189559549752)
```

Updates the enthalpy in the solid state based on the given parameters. :param *\_salinity*: The salinity value. :type *\_salinity*: *float* :param *\_nz*: The nz value. :type *\_nz*: *int* :param *\_liq\_rel*: The liquid relation. Defaults to “Normal”. :type *\_liq\_rel*: *str*, *optional* :param *\_temperature\_melt*: The melting temperature. Defaults to *\_temperature\_melt*. :type *\_temperature\_melt*: *float*, *optional*

**Returns**

The updated enthalpy in the solid state.

**Return type**

float

```
src.spyice.update_physical_values.update_liquid_fraction_buffo(_temperature,  
                                                                _salinity,  
                                                                _liquid_fraction,  
                                                                _enthalpy,  
                                                                _enthalpy_solid,  
                                                                _nz,  
                                                                _is_stefan=False,  
                                                                _method='likebuffo',  
                                                                _pt2_system=False)
```

Updates the liquid fraction based on temperature, salinity, enthalpy, and other parameters.

**Parameters**

- **\_temperature** (*float*) – The temperature value.
- **\_salinity** (*float*) – The salinity value.
- **\_liquid\_fraction** (*float*) – The liquid fraction value.
- **\_enthalpy** (*float*) – The enthalpy value.
- **\_enthalpy\_solid** (*float*) – The solid enthalpy value.
- **\_nz** (*int*) – The number of vertical grid points.
- **\_is\_stefan** (*bool, optional*) – Whether to use Stefan condition. Defaults to False.
- **\_method** (*str, optional*) – The method to use. Defaults to “likebuffo”.

**Returns**

A tuple containing the updated liquid fraction values.

**Return type**

tuple

**Raises**

**AssertionError** – If the liquid fraction has a non-physical value.

```
src.spyice.update_physical_values.update_liquid_fraction_mixture_with_enthalpy_equation(_
```

Update the liquid fraction mixture using under-relaxation. :param \_temperature: The current temperature. :type \_temperature: float :param \_salinity: The salinity of the mixture. :type \_salinity: float :param \_liquid\_fraction\_previous: The previous liquid frac-



tion. :type `_liquid_fraction_previous`: float :param `under_relaxation_factor`: The under-relaxation factor to be applied. :type `under_relaxation_factor`: float :param `_is_stefan`: Flag to indicate if the Stefan condition should be applied. Defaults to False. :type `_is_stefan`: bool, optional

#### Returns

The updated liquid fraction mixture.

#### Return type

float

`src.splyce.update_physical_values.update_liquid_fraction_mixture_with_under_relaxation(_t`

Update the liquid fraction mixture using under-relaxation. :param `_temperature`: The current temperature. :type `_temperature`: float :param `_salinity`: The salinity of the mixture. :type `_salinity`: float :param `_liquid_fraction_previous`: The previous liquid fraction. :type `_liquid_fraction_previous`: float :param `under_relaxation_factor`: The under-relaxation factor to be applied. :type `under_relaxation_factor`: float :param `_is_stefan`: Flag to indicate if the Stefan condition should be applied. Defaults to False. :type `_is_stefan`: bool, optional

#### Returns

The updated liquid fraction mixture.

#### Return type

float

`src.splyce.update_physical_values.update_liquid_fraction_voller_continuous_thermal_proper`

Updates the liquid fraction based on temperature, salinity, enthalpy, and other parameters.

#### Parameters

- **`_temperature`** (*float*) – The temperature value.
- **`_salinity`** (*float*) – The salinity value.
- **`_liquid_fraction`** (*float*) – The liquid fraction value.

- **\_enthalpy** (*float*) – The enthalpy value.
- **\_enthalpy\_solid** (*float*) – The solid enthalpy value.
- **\_nz** (*int*) – The number of vertical grid points.
- **\_is\_stefan** (*bool, optional*) – Whether to use Stefan condition. Defaults to False.
- **\_method** (*str, optional*) – The method to use. Defaults to “like-buffo”.

**Returns**

A tuple containing the updated liquid fraction values.

**Return type**

tuple

**Raises**

**AssertionError** – If the liquid fraction has a non-physical value.

```
src.spylce.update_physical_values.update_liquid_fraction_voller_under_relaxation(_temperature,  
_salinity,  
_temperature,  
_parameter,  
_temperature_initial,  
_liquid_fraction,  
_temperature,  
_parameter,  
_parameter,  
_temperature_previous,  
_salinity_previous,  
_nz,  
temp_factor_3,  
t_k_A_LH,  
t_k_A_LH,  
under_relaxation_factor,  
_is_stefan,  
_pt2_system,  
_non-constant_physical)
```

Update the liquid fraction using the Voller under-relaxation method. :param - *\_temperature*: The temperature value. :type - *\_temperature*: float :param - *\_salinity*: The salinity value. :type - *\_salinity*: float :param - *\_liquid\_fraction*: The liquid fraction value. :type - *\_liquid\_fraction*: float :param - *\_nz*: The number of vertical grid points. :type - *\_nz*: int :param - *temp\_factor\_3*: The temperature factor. :type - *temp\_factor\_3*: float :param - *under\_relaxation\_factor*: The under-relaxation factor. :type - *under\_relaxation\_factor*: float :param - *\_is\_stefan*: Whether to use Stefan condition. Defaults to False. :type - *\_is\_stefan*: bool, optional

**Returns**

The updated liquid fraction values.

**Return type**

- `numpy.ndarray`

**Raises**

- **AssertionError** – If the liquid fraction has a non-physical value

```
src.spyice.update_physical_values.update_state_variables(preprocess_data_object,
                                                         t_prev, s_prev, phi_prev,
                                                         buffo, stefan, voller,
                                                         t_initial, s_initial,
                                                         phi_initial,
                                                         t_k_A_LHS_matrix_prev,
                                                         temp_factor_3,
                                                         source_term,
                                                         _is_salinity_equation=False)
```

Update the state variables Temperature, Salinity, Liquid Fraction, Enthalpy and Enthalpy of solid based on the given parameters.

**Parameters**

- **preprocess\_data\_object** (`PreprocessData` (page ??)) – The preprocess data object.
- **t\_km1** (`numpy.ndarray`) – The previous temperature values.
- **s\_km1** (`numpy.ndarray`) – The previous salinity values.
- **phi\_km1** (`numpy.ndarray`) – The previous liquid fraction values.
- **buffo** (`bool`) – The buffo flag.
- **stefan** (`bool`) – The stefan flag.
- **t\_initial** (`numpy.ndarray`) – The initial temperature values.
- **s\_initial** (`numpy.ndarray`) – The initial salinity values.
- **phi\_initial** (`numpy.ndarray`) – The initial liquid fraction values.
- **source\_term** (`numpy.ndarray`) – The source term values.

**Methods called:**

Updates the state variables based on the given parameters.

- `update_enthalpy`: Update the enthalpy of the system with previous temperature, salinity, liquid fraction.
- `update_enthalpy_solid_state`: Update the enthalpy of the solid state with previous salinity using the liquidus relation.
- `update_liquid_fraction`: Update the liquid fraction of the system with previous temperature, salinity, enthalpy, solid enthalpy.
- `update_temperature_and_salinity`: Update the temperature and salinity of the system with updated liquid fraction and previous temperature, salin-

ity, initial temperature, initial salinity, initial liquid fraction, source term, buffo, stefan.

### Returns

A tuple containing the following updated values:

- `h_k` (`numpy.ndarray`): The updated enthalpy values.
- `h_solid` (`numpy.ndarray`): The updated solid enthalpy values.
- `phi_k` (`numpy.ndarray`): The updated liquid fraction values.
- `t_k` (`numpy.ndarray`): The updated temperature values.
- `s_k` (`numpy.ndarray`): The updated salinity values.

### Return type

tuple

### Parameters

**`preprocess_data_object`** (`PreprocessData` (page ??))

```
src.spyice.update_physical_values.update_temperature_and_salinity(preprocess_data_object,  
                                                                    t_prev,  
                                                                    s_prev,  
                                                                    phi_k,  
                                                                    t_initial,  
                                                                    s_initial,  
                                                                    phi_initial,  
                                                                    source_term,  
                                                                    buffo, stefan,  
                                                                    voller,  
                                                                    _is_salinity_equation=False,  
                                                                    _noncon-  
                                                                    stant_physical_properties=Fa  
                                                                    t_k_A_LHS_matrix_prev=No
```

Update the temperature and salinity based on the given parameters.

### Parameters

- **`preprocess_data_object`** (`PreprocessData` (page ??)) – The preprocess data object.
- **`t_prev`** (`numpy.ndarray`) – The previous temperature values.
- **`s_prev`** (`numpy.ndarray`) – The previous salinity values.
- **`phi_k`** (`numpy.ndarray`) – The liquid fraction values.
- **`t_initial`** (`numpy.ndarray`) – The initial temperature values.
- **`s_initial`** (`numpy.ndarray`) – The initial salinity values.
- **`phi_initial`** (`numpy.ndarray`) – The initial liquid fraction values.
- **`source_term`** (`numpy.ndarray`) – The source term values.
- **`buffo`** (`bool`) – The buffo flag.

- **stefan** (*bool*) – The stefan flag.
- **\_is\_salinity\_equation** (*bool, optional*) – Whether to consider the salinity equation. Defaults to False.
- **voller** (*bool*)
- **\_nonconstant\_physical\_properties** (*bool*)
- **t\_k\_A\_LHS\_matrix\_prev** (*np.ndarray*)

**Returns**

A tuple containing the updated temperature and salinity values. a\_p: The main diagonal of the matrix A for the Voller scheme liquid fraction update

**Return type**

tuple

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