## Installation Guide

1. What type of installation should I choose? ........................................ 1
2. Using the Python-interpreter bundled to Thermo-Calc .............................. 1
   1.2.1 Limitations ........................................................................ 1
   1.2.2 Step 1: Install an IDE (Integrated Development Environment) ............ 2
   1.2.3 Step 2: Configure PyCharm to use the bundled Python-interpreter .......... 2
   1.2.4 Step 3: Run a TC-Python Example ....................................... 3
       1.2.4.1 Open the TC-Python Project in PyCharm ......................... 3
3. Installing TC-Python into the Python-interpreter of your choice ................. 4
   1.3.1 Step 1: Install a Python Distribution ................................... 4
       1.3.1.1 Install Anaconda ...................................................... 4
   1.3.2 Step 2: Install Thermo-Calc and the TC-Python SDK ...................... 4
   1.3.3 Step 3: Install TC-Python ............................................... 4
   1.3.4 Step 4: Install an IDE (Integrated Development Environment) ........... 5
   1.3.5 Step 5: Open the IDE and Run a TC-Python Example ..................... 5
       1.3.5.1 Open the TC-Python Project in PyCharm ....................... 5
       1.3.5.2 Fixing potential issues with the environment ...................... 6
4. Updating to a newer version ................................................................ 6

## Mac OS: Setting Environment Variables

## Architecture overview

1. TCPython .................................................................................. 11
2. SystemBuilder and System ......................................................... 12
3. Calculation .............................................................................. 13
   3.3.1 Single equilibrium calculations ....................................... 13
   3.3.2 Batch equilibrium calculations ........................................ 14
   3.3.3 Precipitation calculations ............................................ 15
   3.3.4 Scheil calculations ....................................................... 15
   3.3.5 Property diagram calculations ....................................... 15
   3.3.6 Phase diagram calculations ........................................... 16
   3.3.7 Diffusion calculations .................................................. 17
   3.3.8 Property Model calculations .......................................... 17
   3.3.9 Material to Material calculations .................................... 18
4. Result ....................................................................................... 18
   3.4.1 DiffusionResult .......................................................... 19
5. Property Model Framework ........................................................ 20
   3.5.1 Property models vs. TC-Python ....................................... 20
   3.5.2 Architecture .............................................................. 20
   3.5.3 Property Model directory .............................................. 21
# 4 Best Practices

4.1 Re-use of the single equilibrium calculation state ................................................. 23
4.2 Re-use and saving of results .................................................................................. 24
4.3 All TC-Python objects are non-copyable .............................................................. 25
4.4 Python Virtual Environments ............................................................................... 25
4.5 Using `TCPython()` efficiently ........................................................................... 25
4.6 Parallel calculations ............................................................................................ 26
4.7 Handling crashes of the calculation engine ............................................................ 27
4.8 Using TC-Python within a Jupyter Notebook or the Python console .................... 28
4.9 Property Model Framework ................................................................................. 29
   4.9.1 Debugging Property Model code ................................................................ 29
   4.9.2 Developing Property Models in several files ................................................ 31
      4.9.2.1 `side-by-side` modules ...................................................................... 31
      4.9.2.2 `common` modules .......................................................................... 32
   4.9.3 Alternative Python for Property Models ....................................................... 33
      4.9.3.1 Default bundled Python interpreter .................................................. 33
      4.9.3.2 Configuring another Python interpreter ............................................ 33

# 5 API Reference

5.1 Calculations ........................................................................................................... 35
   5.1.1 Module “single_equilibrium” ........................................................................ 35
   5.1.2 Module “batch_equilibrium” ......................................................................... 48
   5.1.3 Module “precipitation” .................................................................................. 52
   5.1.4 Module “scheil” .......................................................................................... 74
   5.1.5 Module “step_or_map_diagrams” .................................................................. 85
   5.1.6 Module “diffusion” ..................................................................................... 112
   5.1.7 Module “propertymodel” ............................................................................ 150
   5.1.8 Module “material_to_material” .................................................................. 155
   5.2 Module “system” .............................................................................................. 178
   5.3 Module “entities” ............................................................................................. 186
   5.4 Module “server” ............................................................................................. 190
   5.5 Module “quantity_factory” ............................................................................. 195
   5.6 Module “utils” ............................................................................................... 207
   5.7 Module “propertymodel_sdk” ......................................................................... 209
   5.8 Module “exceptions” ....................................................................................... 225
   5.9 Module “abstract_base” .................................................................................. 227

# 6 Troubleshooting

6.1 Diagnostics script ............................................................................................... 235
6.2 “No module named tc_python” error on first usage .............................................. 237
6.3 “pip install” fails with “Failed to establish a new network connection” or similar ... 238

Python Module Index

239
This guide helps you to get a working TC-Python API installation.

There is a PDF guide included with your installation. In the Thermo-Calc menu, select Help → Manuals Folder. Then double-click to open the Software Development Kits (SDKs) folder.

Note: A license is required to run TC-Python.

1.1 What type of installation should I choose?

There are two possibilities to install TC-Python:

1. Using the Python-interpreter bundled to Thermo-Calc: This interpreter has TC-Python preinstalled together with some popular Python-packages. This is the recommended option for new users to TC-Python, but it is limited to the preinstalled packages.

2. Installing TC-Python into the Python-interpreter of your choice: This is the recommended option for any more advanced usage and provides full flexibility.

1.2 Using the Python-interpreter bundled to Thermo-Calc

Note: A Python-interpreter is bundled to Thermo-Calc beginning with version 2021a.

1.2.1 Limitations

The bundled Python 3.7.2 interpreter is containing the following major packages:

<table>
<thead>
<tr>
<th>Package</th>
<th>Version</th>
</tr>
</thead>
<tbody>
<tr>
<td>matplotlib</td>
<td>3.3.2</td>
</tr>
<tr>
<td>numpy</td>
<td>1.19.2</td>
</tr>
<tr>
<td>scikit-learn</td>
<td>0.23.2</td>
</tr>
<tr>
<td>scipy</td>
<td>1.5.2</td>
</tr>
<tr>
<td>TC-Python</td>
<td>2022a</td>
</tr>
</tbody>
</table>
Please contact the Thermo-Calc support if you think that further packages might be useful in future releases.

**Note:** The following TC-Python examples are requiring additional packages that are not available in the bundled Python-interpreter, they can therefore not be run:

- `pyex_M_01_Input_from_file.py`
- `pyex_M_02_Ouput_to_file.py`

**Warning:** The Python-interpreter bundled to Thermo-Calc is also used for running the property models in Thermo-Calc. Any changes to the interpreter packages can therefore break Thermo-Calc and should be avoided. If the installation gets broken, it can be fixed by reinstalling Thermo-Calc after having removed it.

### 1.2.2 Step 1: Install an IDE (Integrated Development Environment)

Any editor can be used to write the Python code, but an IDE is recommended, e.g. PyCharm. These instructions are based on the use of PyCharm.

Use of an IDE will give you access to code completion, which is of great help when you use the API as it will give you the available methods on the objects you are working with.

2. Click to choose your OS and then click Download. You can use the Community version of PyCharm.
3. Follow the instructions. It is recommended you keep all the defaults.

**Note:** For Mac installations, you also need to set some environment variables as described below in Mac OS: Setting Environment Variables.

### 1.2.3 Step 2: Configure PyCharm to use the bundled Python-interpreter

Open PyCharm and configure the interpreter:

1. Go the menu File→Settings.
2. Navigate in the tree to Project: YourProjectName and choose Project Interpreter.
3. Click on the settings symbol close to the Project Interpreter dropdown menu and choose Add.
4. Now choose System Interpreter and add the bundled Thermo-Calc Python 3 interpreter. It is located in different places depending on the operating system:
## Operating system

<table>
<thead>
<tr>
<th>Operating system</th>
<th>Path to the bundled Python-interpreter</th>
</tr>
</thead>
<tbody>
<tr>
<td>Windows</td>
<td>C:\Program Files\Thermo-Calc\2022a\python\python.exe</td>
</tr>
<tr>
<td>Linux</td>
<td>/home/UserName/Thermo-Calc/2022a/python/bin/python3</td>
</tr>
<tr>
<td>MacOS</td>
<td>/Applications/Thermo-Calc-2022a.app/Contents/Resources/python/bin/python3</td>
</tr>
</tbody>
</table>

5. Select your added interpreter and confirm.

### 1.2.4 Step 3: Run a TC-Python Example

Now you are ready to start working with TC-Python.

It is recommended that you open one or more of the included examples to both check that the installation has worked and to start familiarizing yourself with the code.

#### 1.2.4.1 Open the TC-Python Project in PyCharm

When you first open the TC-Python project and examples, it can take a few moments for the Pycharm IDE to index before some of the options are available.

1. Open PyCharm and then choose File→Open. The first time you open the project you will need to navigate to the path of the TC-Python installation:

<table>
<thead>
<tr>
<th>Operating system</th>
<th>Path to the TC-Python folder</th>
</tr>
</thead>
<tbody>
<tr>
<td>Windows</td>
<td>C:\Users\UserName\Documents\Thermo-Calc\2022a\SDK\TC-Python</td>
</tr>
<tr>
<td>Linux</td>
<td>/home/UserName/Thermo-Calc/2022a/SDK/TC-Python</td>
</tr>
<tr>
<td>MacOS</td>
<td>/Users/Shared/Thermo-Calc/2022a/SDK/TC-Python</td>
</tr>
</tbody>
</table>

2. Click on the Examples folder and then click OK.

3. From any subfolder:
   - Double-click to open an example file to examine the code.
   - Right-click an example and choose Run.

**Note:** If you are not following the recommended approach and create a new project (File→New Project...), you need to consider that by default the options to choose the interpreter are hidden within the Create Project window. So click on Project Interpreter: New Virtual Environment and in most cases choose your System Interpreter containing the Python bundled to Thermo-Calc instead of the default New Virtual Environment.
1.3 Installing TC-Python into the Python-interpreter of your choice

1.3.1 Step 1: Install a Python Distribution

If you already have a Python distribution installation, version 3.5 or higher, skip this step.

These instructions are based on using the Anaconda platform for the Python distribution. Install version 3.5 or higher to be able to work with TC-Python, although it is recommended that you use the most recent version.

1.3.1.1 Install Anaconda

2. Click to choose your OS (operating system) and then click Download. Follow the instructions. It is recommended you keep all the defaults.

1.3.2 Step 2: Install Thermo-Calc and the TC-Python SDK

Note: TC-Python is available starting with Thermo-Calc version 2018a.

1. Install Thermo-Calc
2. When the installation is complete, open the TC-Python folder that includes the *.whl file needed for the next step. There is also an file:Examples folder with Python files you can use in the IDE to understand and work with TC-Python.

1.3.3 Step 3: Install TC-Python

On Windows, it is recommended that you use the Python distribution prompt (i.e. Anaconda, ...), especially if you have other Python installations. Do not use Virtual Environments unless you have a good reason for that.

1. Open the command line. For example, in Anaconda on a Windows OS, go to Start → Anaconda → Anaconda Prompt.
2. At the command line, enter the following. Make sure there are no spaces at the end of the string or in the folder name or it will not run:

```
pip install <path to the TC-Python folder>/TC_Python-<version>-py3-none-any.whl
```

Tip: Note that on Linux depending on the interpreter usually pip3 is used.

<table>
<thead>
<tr>
<th>Operating system</th>
<th>Path to the TC-Python folder</th>
</tr>
</thead>
<tbody>
<tr>
<td>Windows</td>
<td>C:\Users\UserName\Documents\Thermo-Calc\2022a\SDK\TC-Python</td>
</tr>
<tr>
<td>Linux</td>
<td>/home/UserName/Thermo-Calc/2022a/SDK/TC-Python</td>
</tr>
<tr>
<td>MacOS</td>
<td>/Users/Shared/Thermo-Calc/2022a/SDK/TC-Python</td>
</tr>
</tbody>
</table>
3. Press <Enter>. When the process is completed, there is a confirmation that TC-Python is installed.

**Note:** If your computer is located behind a proxy-server, the default `pip`-command will fail with a network connection error. In that case you need to install the dependencies of TC-Python in a special configuration:

```
pip install -proxy user:password@proxy_ip:port py4j jproperties
```

See “pip install” fails with “Failed to establish a new network connection” or similar for detailed information.

### 1.3.4 Step 4: Install an IDE (Integrated Development Environment)

Any editor can be used to write the Python code, but an IDE is recommended, e.g. PyCharm. These instructions are based on the use of PyCharm.

Use of an IDE will give you access to code completion, which is of great help when you use the API as it will give you the available methods on the objects you are working with.

2. Click to choose your OS and then click **Download**. You can use the **Community** version of PyCharm.
3. Follow the instructions. It is recommended you keep all the defaults.

**Note:** For Mac installations, you also need to set some environment variables as described below in *Mac OS: Setting Environment Variables*.

### 1.3.5 Step 5: Open the IDE and Run a TC-Python Example

After you complete all the software installations, you are ready to open the IDE to start working with TC-Python.

It is recommended that you open one or more of the included examples to both check that the installation has worked and to start familiarizing yourself with the code.

#### 1.3.5.1 Open the TC-Python Project in PyCharm

When you first open the TC-Python project and examples, it can take a few moments for the Pycharm IDE to index before some of the options are available.

1. Open PyCharm and then choose **File**→**Open**. The first time you open the project you will need to navigate to the path of the TC-Python installation.

<table>
<thead>
<tr>
<th>Operating system</th>
<th>Path to the TC-Python folder</th>
</tr>
</thead>
<tbody>
<tr>
<td>Windows</td>
<td>C:\Users\UserName\Documents\Thermo-Calc\2022a\SDK\TC-Python</td>
</tr>
<tr>
<td>Linux</td>
<td>/home/UserName/Thermo-Calc/2022a/SDK/TC-Python</td>
</tr>
<tr>
<td>MacOS</td>
<td>/Users/Shared/Thermo-Calc/2022a/SDK/TC-Python</td>
</tr>
</tbody>
</table>
2. Click on the Examples folder and then click OK.
3. From any subfolder:
   • Double-click to open an example file to examine the code.
   • Right-click an example and choose Run.

1.3.5.2 Fixing potential issues with the environment

In most cases you should run TC-Python within your global Python 3 interpreter and not use Virtual Environments unless you have a good reason to do so. A common problem on first usage of TC-Python is the error message “No module named tc_python”. You can resolve this and other problems with the interpreter settings as follows:

1. Go the menu File → Settings.
2. Navigate in the tree to Project: YourProjectName and choose Project Interpreter.
3. Click on the settings symbol close to the Project Interpreter dropdown menu and choose Add.
4. Now choose System Interpreter and add your existing Python 3 interpreter.
5. Select your added interpreter and confirm.

Note: If you are not following the recommended approach and create a new project (File → New Project...), you need to consider that by default the options to choose the interpreter are hidden within the Create Project window. So click on Project Interpreter: New Virtual Environment and in most cases choose your System Interpreter instead of the default New Virtual Environment.

Note: If you really need to use a Virtual Environment, please consider the hints given in the Python Virtual Environments chapter.

1.4 Updating to a newer version

When updating to a newer version of Thermo-Calc, you need to also install the latest version of TC-Python. This is not necessary if are using the bundled Python-interpreter that has it automatically installed. It is not sufficient to run the installer of Thermo-Calc:

```
pip install <path to the TC-Python folder>/TC_Python-<version>-py3-none-any.whl
```

Tip: Note that on Linux depending on the interpreter usually pip3 is used.

In case of problems you may wish to uninstall the previous version of TC-Python in advance:

```
pip uninstall TC-Python
pip install <path to the TC-Python folder>/TC_Python-<version>-py3-none-any.whl
```
However, that should normally not be required.

You can check the currently installed version of TC-Python by running:

```
pip show TC-Python
```
MAC OS: SETTING ENVIRONMENT VARIABLES

In order to use TC-Python on Mac you need to set some environment variables.

TC22A_HOME=/Applications/Thermo-Calc-2022a.app/Contents/Resources

If you use a license server:

LSHOST=<name-of-the-license-server>

If you have a node-locked license:

LSHOST=NO-NET
LSERVRC=/Users/Shared/Thermo-Calc/lservrc

In PyCharm, you can add environment variables in the configurations.

Select Run→Edit Configurations to open the Run/Debug Configurations window. Choose Templates and then Python. Enter the environment variable(s) by clicking the button to the right of the Environment Variables text field. Now the environment variables(s) will be set for each new configuration by default.

Note: Existing configurations need to be removed and recreated to obtain the environment variables in them.

The same way for configuring the environment variables can be used on other operating systems as if necessary.
TC-Python contains classes of these types:

- **TCPython** – this is where you start with general settings.
- **SystemBuilder** and **System** – where you choose database and elements etc.
- **Calculation** – where you choose and configure the calculation.
- **Result** – where you get the results from a calculation you have run.

### 3.1 TCPython

This is the starting point for all TC-Python usage.

You can think of this as the start of a “wizard”.

You use it to select databases and elements. That will take you to the next step in the wizard, where you configure the system.

**Example:**

```python
from tc_python import *

with TCPython() as start:
    start.select_database_and_elements(...
    # e.t.c
    # after with clause

# or like this
with TCPython():
    SetUp().select_database_and_elements(...
    # e.t.c
    # after with clause
```

**Tip:** If you use TC-Python from Jupyter Lab / Notebook, you should use TC-Python slightly different to be able to use multiple cells. See *Using TC-Python within a Jupyter Notebook or the Python console* for details.

**Note:** When your python script runs a row like this:

```python
with TCPython() as start:
```
a process running a calculation server starts. Your code, via TC-Python, uses socket communication to send and receive messages to and from that server.

When your Python script has run as far as this row:

```python
# after with clause
```

the calculation server automatically shuts down, and all temporary files are deleted. It is important to ensure that this happens by structuring your Python code using a `with()` clause as in the above example.

**Note:** To re-use results from previous calculations, set a folder where TC-Python saves results, and looks for previous results.

This is done with the function `set_cache_folder()`.

```python
from tc_python import *
with TCPython() as start:
    start.set_cache_folder("cache")
```

This folder can be a network folder and shared by many users. If a previous TC-Python calculation has run with the same cache_folder and EXACTLY the same system and calculation settings, the calculation is not re-run. Instead the result is automatically loaded from disk.

It is also possible to explicitly save and load results.

```python
from tc_python import *
with TCPython() as start:
    #... diffusion calculation (could be any calculation type)
    calculation_result.save_to_disk('path to folder')
    #...
    loaded_result = start.load_result_from_disk().diffusion('path to folder')
```

### 3.2 SystemBuilder and System

A **SystemBuilder** is returned when you have selected your database and elements in **TCPython**.

The **SystemBuilder** lets you further specify your system, for example the phases that should be part of your system.

**Example:**

```python
from tc_python import *
with TCPython() as start:
    start.select_database_and_elements("ALDEMO", ["Al", "Sc"])  
    # e.t.c
```

When all configuration is done, you call `get_system()` which returns an instance of a **System** class. The **System** class is fixed and cannot be changed. If you later want to change the database, elements or something else, change the **SystemBuilder** and call `get_system()` again, or create a new **SystemBuilder** and call `get_system()`.

From the **System** you can create one or more calculations, which is the next step in the “wizard”.
Note: You can use the same System object to create several calculations.

3.3 Calculation

The best way to see how a calculation can be used is in the TC-Python examples included with the Thermo-Calc installation.

Some calculations have many settings. Default values are used where it is applicable, and are overridden if you specify something different.

When you have configured your calculation you call calculate() to start the actual calculation. That returns a Result, which is the next step.

3.3.1 Single equilibrium calculations

In single equilibrium calculations you need to specify the correct number of conditions, depending on how many elements your System contains.

You do that by calling set_condition().

An important difference from other calculations is that single equilibrium calculations have two functions to get result values.

The calculate() method, which gives a SingleEquilibriumTempResult, is used to get actual values. This result is “temporary”, meaning that if you run other calculations or rerun the current one, the resulting object no longer gives values corresponding to the first calculation.

This is different from how other calculations work. If you want a Result that you can use after running other calculations, you need to call calculate_with_state(), which returns a SingleEquilibriumResult.

Note: calculate() is the recommended function and works in almost all situations. Also it has much better performance than calculate_with_state().

Example:

```python
from tc_python import *

with TCPython() as start:
    gibbs_energy = {
        start.
            select_database_and_elements("FEDEMO", ["Fe", "Cr", "C"]).
            get_system().
            with_single_equilibrium_calculation().
                set_condition(ThermodynamicQuantity.temperature(), 2000.0).
                set_condition(ThermodynamicQuantity.mole_fraction_of_a_component("Cr
                    →"), 0.1).
                set_condition(ThermodynamicQuantity.mole_fraction_of_a_component("C"),
                    → 0.01).
                calculate().
                get_value_of("G")
            }
```
3.3.2 Batch equilibrium calculations

Batch equilibrium calculations are used when you want to do many single equilibrium calculations and it is known from the beginning which result values are required from the equilibrium. This is a vectorized type of calculation that can reduce the overhead from Python and TC-Python similar to the approach used in *numpy*-functions for example.

**Tip:** The performance of batch equilibrium calculations can be significantly better than looping and using single equilibrium calculations if the actual Thermo-Calc calculation is fast. There is little advantage if the Thermo-Calc equilibrium calculations take a long time (typically for large systems and databases).

Example:

```python
from tc_python import *

with TCPython() as start:
    calculation = (start
        .set_cache_folder(os.path.basename(__file__) + "_cache")
        .select_database_and_elements("NIDEMO", ["Ni", "Al", "Cr")
        .get_system()
        .with_batch_equilibrium_calculation()
        .set_condition("T", 800.1)
        .set_condition("X(Al)", 1E-2)
        .set_condition("X(Cr)", 1E-2)
        .disable_global_minimization()
    )

list_of_x_A1 = [1, 2, 3, 4, 5, 6, 7, 8, 9, 10]
list_of_x_Cr = [3, 5, 7, 9, 11, 13, 15]
lists_of_conditions = []
for x_A1 in list_of_x_A1:
    for x_Cr in list_of_x_Cr:
        lists_of_conditions.append(["X(Al)", x_A1 / 100],
                                   ["X(Cr)", x_Cr / 100])
    calculation.set_conditions_for_equilibria(lists_of_conditions)

results = calculation.calculate(["BM", "VM"])

masses = results.get_values_of("BM")
volumes = results.get_values_of("VM")

print(masses)
print(volumes)
```
3.3.3 Precipitation calculations

All that can be configured in the Precipitation Calculator in Graphical Mode can also be done here in this calculation. However, you must at least enter a matrix phase, a precipitate phase, temperature, simulation time and compositions.

Example:

```python
from tc_python import *

with TCPython() as start:
    precipitation_curve = (    start.
        select_thermodynamic_and_kinetic_databases_with_elements("ALDEMO", "MALDEMO", ["Al", "Sc"]).
        get_system().
        with_isothermal_precipitation_calculation().
            set_composition("Sc", 0.18).
            set_temperature(623.15).
            set_simulation_time(1e5).
            with_matrix_phase(MatrixPhase("FCC_A1").
                add_precipitate_phase(PrecipitatePhase("AL3SC")).
            )
        .calculate()
    )
```

3.3.4 Scheil calculations

All Scheil calculations available in Graphical Mode or Console Mode can also be done here in this calculation. The minimum you need to specify are the elements and compositions. Everything else is set to a default value.

Example:

```python
from tc_python import *

with TCPython() as start:
    temperature_vs_mole_fraction_of_solid = (        start.
        select_database_and_elements("FEDEMO", ["Fe", "C"]).
        get_system().
        with_scheil_calculation().
            set_composition("C", 0.3).
            calculate().
            get_values_of(ScheilQuantity.temperature(),
                ScheilQuantity.mole_fraction_of_all_solid_phases())
    )
```

3.3.5 Property diagram calculations

For the property diagram (step) calculation, everything that you can configure in the Equilibrium Calculator when choosing Property diagram in Graphical Mode can also be configured in this calculation. In Console Mode the property diagram is created using the Step command. The minimum you need to specify are elements, conditions and the calculation axis. Everything else is set to default values, if you do not specify otherwise.

Example:
from tc_python import *

with TCPython() as start:
    property_diagram = {
        start.
        select_database_and_elements("FEDEMO", ["Fe", "C"]).
        get_system().
        with_property_diagram_calculation().
        with_axis(CalculationAxis(ThermodynamicQuantity.temperature())).
        set_min(500).
        set_max(3000)).
        set_condition(ThermodynamicQuantity.mole_fraction_of_a_component("C"),
                      0.01).
        calculate().
        get_values_grouped_by_stable_phases_of(ThermodynamicQuantity.
                                              temperature(),
                                              ThermodynamicQuantity.volume_)
        →fraction_of_a_phase("ALL")
    }

3.3.6 Phase diagram calculations

For the phase diagram (map) calculation, everything that you can configure in the Equilibrium Calculator when choosing Phase diagram in Graphical Mode can also be configured in this calculation. In Console Mode the phase diagram is created using the Map command. The minimum you need to specify are elements, conditions and two calculation axes. Everything else is set to default values, if you do not specify otherwise.

Example:

from tc_python import *

with TCPython() as start:
    phase_diagram = {
        start.
        select_database_and_elements("FEDEMO", ["Fe", "C"]).
        get_system().
        with_phase_diagram_calculation().
        with_first_axis(CalculationAxis(ThermodynamicQuantity.temperature())).
        set_min(500).
        set_max(3000)).
        with_second_axis(CalculationAxis(ThermodynamicQuantity.mole_fraction_of_a_component("C")),
                         set_min(0).
                         set_max(1)).
        set_condition(ThermodynamicQuantity.mole_fraction_of_a_component("C"),
                      0.01).
        calculate().
        get_values_grouped_by_stable_phases_of(ThermodynamicQuantity.mass_)
        →fraction_of_a_component("C"),
                      ThermodynamicQuantity.
        →temperature())
    }
3.3.7 Diffusion calculations

For diffusion calculations, everything that you can configure in the Diffusion Calculator can also be configured in this calculation. The minimum you need to specify are elements, temperature, simulation time, a region with a grid and width, a phase and an initial composition.

Example:

```python
from tc_python import *

with TCPython() as start:
    diffusion_result = {
        start.
        select_thermodynamic_and_kinetic_databases_with_elements("FEDEMO",  
        "MFEDEMO", ["Fe", "Ni"]).
        set_system().
        with_isothermal_diffusion_calculation().
        set_temperature(1400.0).
        set_simulation_time(108000.0).
        add_region(Region("Austenite").
            set_width(1E-4).
            with_grid(CalculatedGrid.linear().set_no_of_points(50)).
            with_composition_profile(CompositionProfile().
                add("Ni", ElementProfile.linear(10.0, 50.0))
            ).
        add_phase("FCC_A1")).
        calculate()
    }

distance, ni_fraction = diffusion_result.get_mass_fraction_of_component_at_time(
    "Ni", 108000.0)
```

3.3.8 Property Model calculations

For Property Model calculations, everything that you can configure in the Property Model Calculator in Graphical Mode can also be configured in this calculation. The minimum you need to specify are elements, composition and which Property Model you want to use.

Example:

```python
from tc_python import *

with TCPython() as start:
    property_model = {
        start.
        select_database_and_elements("FEDEMO", ["Fe", "C"]).
        get_system().
        with_property_model_calculation("Driving force").
        set_composition("C", 1.0).
        set_argument("precipitate", "GRAPHITE")

        print("Available arguments: /": format(property_model.get_arguments()))
        result = property_model.calculate()

        print("Available result quantities: /": format(result.get_result_quantities()))
        driving_force = result.get_value_of("normalizedDrivingForce")
    }
```
3.3.9 Material to Material calculations

Material to Material calculations are generally regular single equilibrium, property diagram or phase diagram calculations but they are specialised to handle the mixture of two materials A and B. Everything that you can configure in the Material to Material Calculator in Graphical Mode can also be configured in this calculation. The minimum required configuration is shown below for a Property diagram calculation for varying amount of material B. The other calculators (single fraction of material B and phase diagram calculations) are configured in a similar way.

Example:

```python
define tc_python import *
with TCPython() as start:
    material_to_material_property_diagram = (  
        start.
            select_database_and_elements("FEDEMO", ["Fe", "Cr", "Ni", "C"]).  
            get_system().  
            with_material_to_material().  
            with_property_diagram_calculation().  
            set_material_a("Cr": 10.0, "Ni": 15.0, "Fe").  
            set_material_b("Cr": 15.0, "Ni": 10.0, "Fe").  
            set_activities("C": 0.1).  
            with_constant_condition(ConstantCondition.temperature(800 + 273.15)).  
            with_axis(MaterialToMaterialCalculationAxis.fraction_of_material_b(from_  
                --fraction=0.0,  
                --fraction=1.0,  
                --fraction=0.5))  
        )  
    result = material_to_material_property_diagram.calculate()  
    data = result.get_values_grouped_by_quantity_of(MATERIAL_B_FRACTION,  
        ThermodynamicQuantity.volume_  
        --fraction_of_a_phase(ALL_PHASES))  
    for group in data.values():  
        fractions_of_b = group.x  
        volume_fractions_of_phase = group.y  
        phase_name = group.label
```

3.4 Result

All calculations have a method called calculate() that starts the calculations and when finished, returns a Result. The Result classes have very different methods, depending on the type of calculation.

The Result is used to get numerical values from a calculation that has run.

The Result can be saved to disk by the method save_to_disk().

Previously saved results can be loaded by the method load_result_from_disk() on the SetUp class.

Example:
# code above sets up the calculation
r = calculation.calculate()
time, meanRadius = r.get_mean_radius_of("AL3SC")

The **Result** objects are completely independent from calculations done before or after they are created. The objects return valid values corresponding to the calculation they were created from, for their lifetime. The only exception is if you call `calculate()` and not `calculate_with_state()` on a single equilibrium calculation.

As in the following example you can mix different calculations and results, and use old results after another calculation has run.

**Example:**

```python
# ...
# some code to set up a single equilibrium calculation
# ...
single_eq_result = single_eq_calculation.calculate_with_state()
# ...
# some code to set up a precipitation calculation
# ...
prec_result = precipitation_calculation.calculate()
# ...
# some code to set up a Scheil calculation
# ...
scheil_result = scheil_calculations.calculate()
# now it is possible to get results from the single equilibrium calculation,
# without having to re-run it (because it has been calculated with saving of the_
# state)
gibbs = single_eq_result.get_value_of("G")
```

### 3.4.1 DiffusionResult

The DiffusionResult class, that is returned when calling `calculate()` on any DiffusionCalculation, has the possibility to create a ContinuedDiffusionCalculation, in addition to the “normal” functionality for results. This makes it possible to run a diffusion calculation and then, depending on the result, change some settings and continue.

**Example:**

```python
# ...
# some code to set up a Diffusion calculation
# ...
first_diffusion_result = diffusion_calculation.calculate()
# ...
continued_calculation = first_diffusion_result.with_continued_calculation()
continued_calculation.set_simulation_time(110000.0)
continued_calculation.with_left_boundary_condition(BoundaryCondition.mixed_zero_flux_#
and_activity().set_activity_for_element('C', 1.0))
second_result = continued_calculation.calculate()
```

(continues on next page)
3.5 Property Model Framework

The **Python Property Model SDK** extends the Thermo-Calc software to enable you to create your own Property Models. A **Property Model** is a Python-based calculation that can use any TC-Python functionality (including diffusion and precipitation calculations) but is usable through the Graphical User Interface (UI) of Thermo-Calc in a more simple way. It is typically used to model material properties but by no means limited to that. Examples of Property Models provided by Thermo-Calc include Martensite and Pearlite formation in steel.

The Property Model Framework uses standard Python 3 beginning with Thermo-Calc 2021a and can access all TC-Python functionality and any Python package including **numpy**, **scipy**, **tensorflow**, etc. The actual calculation code is nearly identical, regardless if called from within a Property Model or from standard Python.

This is a complete rewrite of the original version of the framework that was based on Jython 2.7 and therefore had a number of limitations. **Property models written with the old Property Model Framework before Thermo-Calc 2021a are not compatible with the new framework.** However, the migration should be relatively easy because the syntax was changed as little as possible.

3.5.1 Property models vs. TC-Python

The main difference between a **Property Model** and regular **TC-Python code** is that a Property Model is directly integrated into the UI of Thermo-Calc via a plugin architecture while TC-Python code can only be accessed by programs and scripts written in Python.

The user should develop a Property Model if the functionality needs to be available from the Thermo-Calc UI, especially if it should be applied by other users not familiar to programming languages. Otherwise it is preferable to implement the functionality directly in a TC-Python program. If required, Property Models can as well be accessed from within TC-Python.

3.5.2 Architecture

Every Property Model needs to contain a class that implements the interface `tc_python.propertymodel_sdk.PropertyModel`. There are naming conventions that must to be fulfilled: the file name is required to follow the the pattern `XYPythonModel.py` and the name of the class needs to match this. Additionally the file must be placed in a directory named `XYPython` within the Property Model directory. The content of the placeholder `XY` can be freely chosen.

A simple complete Property Model, saved in a file called `SimplePythonModel.py` in the directory `SimplePython`, looks like this:

```python
from tc_python import *

class SimplePythonModel(PropertyModel):
    def provide_model_category(self) -> List[str]:
        return ["Demo"]
```

(continues on next page)
def provide_model_name(self) -> str:
    return "My Demo Model"

def provide_model_description(self) -> str:
    return "This is a demo model."

def provide_ui_panel_components(self) -> List[UIComponent]:
    return [UIBooleanComponent("CHECKBOX", "Should this be checked?", "Simple checkbox", setting=False)]

def provide_calculation_result_quantities(self) -> List[ResultQuantity]:
    return [create_general_quantity("RESULT", "A result")]

def evaluate_model(self, context: CalculationContext):
    if context.get_ui_boolean_value("CHECKBOX"):
        self.logger.info("The checkbox is checked")

        composition_as_mass_fraction = context.get_mass_fractions()
        temp_in_k = context.get_temperature()
        calc = context.system.with_single_equilibrium_calculation()

        context.set_result_quantity_value="RESULT", 5.0)  # the value would normally have been calculated

The basic building blocks of the Property Model API are:

- `tc_python.propertymodel_sdk.ResultQuantity`: Defines a calculation result of a Property Model that will be provided to the UI after each model evaluation
- `tc_python.propertymodel_sdk.CalculationContext`: Provides access to the data from the UI (such as the entered composition and temperature) and to the current TC-Python system object which is the entrypoint for using TC-Python from within the Property Model
- `tc_python.propertymodel_sdk.UIComponent`: These are the UI-components that create the user interface of the Property Model within the model panel of the Thermo-Calc application UI. Different components are available (for example checkboxes, text fields and lists).

### 3.5.3 Property Model directory

The Property Model py-files need to be located within subdirectories of the Property Model directory, e.g. PropertyModelDir/XYPython/XYPythonModel.py. The default Property Model directory can be changed in the menu Tools -> Options in the graphical user interface.

<table>
<thead>
<tr>
<th>Operating system</th>
<th>Default Property Model directory</th>
</tr>
</thead>
<tbody>
<tr>
<td>Windows</td>
<td>C:\Users\UserName\Documents\Thermo-Calc\2022a\PropertyModels</td>
</tr>
<tr>
<td>Linux</td>
<td>/home/UserName/Thermo-Calc/2022a/PropertyModels</td>
</tr>
<tr>
<td>MacOS</td>
<td>/Users/Shared/Thermo-Calc/2022a/PropertyModels</td>
</tr>
</tbody>
</table>
4.1 Re-use of the single equilibrium calculation state

The Thermo-Calc core keeps an internal state containing the data from previously performed calculations (such as composition of sublattices, previously formed phases, ...). This will be used for start values of future calculations (if not explicitly overwritten) and can strongly influence their convergence and calculation time. It can be useful to save and restore later the core-state in advanced use cases, these include:

- Improving the convergence speed in case of very complicated equilibria if a similar equilibrium had been calculated already before. Similarity refers here primarily to composition, temperature and entered phase set. This case can occur for example with the Nickel-database TCNi.

- Convenient and fast switching between states that have changed a lot (for example regarding suspended phases, numerical settings, ...)

The mechanism of saving and restoring the state is called bookmarking and is controlled with the two methods `tc_python.single_equilibrium.SingleEquilibriumCalculation.bookmark_state()` and `tc_python.single_equilibrium.SingleEquilibriumCalculation.set_state_to_bookmark()`. The following short example demonstrates how to switch between two different states easily in practice:

```python
from tc_python import *

with TCPython() as session:
    calc = (session.
        select_database_and_elements("FEDEMO", ["Fe", "C"]).
        get_system().
        with_single_equilibrium_calculation().
        set_condition(ThermodynamicQuantity.temperature(), 2000.0).
        set_condition("X(C)", 0.01))
    
    calc.calculate()
    bookmark_temp_condition = calc.bookmark_state()
    calc.set_phase_to_fixed("BCC", 0.5)
    calc.remove_condition(ThermodynamicQuantity.temperature())
    bookmark_fixed_phase_condition = calc.bookmark_state()
    
    result_temp = calc.set_state_to_bookmark(bookmark_temp_condition)
    print("Conditions do contain temperature: {}\n".format(result_temp.get_conditions()))
    # this calculation had already been performed
    print("Stable phases (do not contain BCC): {}\n".format(result_temp.get_stable_phases()))
```

(continues on next page)
result_fixed_phase = calc.set_state_toBookmark(bookmark_fixed_phase_condition)
print("Conditions do not contain temperature: \{}").format(result_fixed_phase.get_conditions())
# this calculation had **not yet** been performed
print("Stable phases (do contain BCC): \{}").format(calc.calculate().get_stable_phases())

4.2 Re-use and saving of results

Before a calculation is run in TC-Python, a check is made to see if the exact same calculation has run before, and if that is the case, the result from the calculation can be loaded from disk instead of being re-calculated.

This functionality is always enabled within a script running TC-Python, but you can make it work the same way when re-running a script, or even when running a completely different script.

To use results from previous calculations, set a folder where TC-Python saves results, and looks for previous results. This is controlled by the method `tc_python.server.SetUp.set_cache_folder()`.

```python
from tc_python import *

with TCPython() as start:
    start.set_cache_folder("cache")
```

This folder can be a network folder and shared by many users. The calculation is not re-run if there is a previous TC-Python calculation with the same cache folder and exactly the same settings; the result is instead loaded from disk.

Another possibility is to explicitly save the result to disk and reload it later:

```python
from tc_python import *

with TCPython() as start:
    # ... the system and calculator are set up and the calculation is performed
    result = calculator.calculate()
    result.save_to_disk("./result_dir")
```

You can then load the result again in another session:

```python
from tc_python import *

with TCPython() as start:
    result = SetUp().load_result_from_disk().diffusion("./result_dir")
    x, frac = result.get_mole_fraction_of_component_at_time("Cr", 1000.0)
```
4.3 All TC-Python objects are non-copyable

*Never create a copy* of an instance of a class in TC-Python, neither by using the Python built-in function `deepcopy()` nor in any other way. All classes in TC-Python are proxies for classes in the underlying calculation server and normally hold references to result files. A copied class object in Python would consequently point to the same classes and result files in the calculation server.

Instead of making a copy, always create a new instance:

```python
from tc_python import *

with TCPython() as start:
    system = start.select_database_and_elements("FEDEMO", ["Fe", "Cr"])
    calculator = system.with_single_equilibrium_calculation()

    # *do not* copy the 'calculator' object, create another one instead
    calculator_2 = system.with_single_equilibrium_calculation()

    # now you can use both calculators for different calculations ...
```

4.4 Python Virtual Environments

A Python installation can have several virtual environments. You can think of a virtual environment as a collection of third party packages that you have access to in your Python scripts. `tc_python` is such a package.

To run TC-Python, you need to install it into the same virtual environment as your Python scripts are running in. If your scripts fail on `import tc_python`, you need to execute the following command in the terminal of the same Python environment as your script is running in:

```
pip install TC_Python-<version>-py3-none-any.whl
```

If you use the PyCharm IDE, you should do that within the `Terminal` built into the IDE. This `Terminal` runs automatically within your actual (virtual) environment.

To prevent confusion, it is recommend in most cases to install TC-Python within your global interpreter, for example by running the `pip install` command within your default Anaconda prompt.

4.5 Using `with TCPython()` efficiently

Normally you should call `with TCPython()` only once within each process.

**Note:** When leaving the `with`-clause, the Java backend engine process is stopped and all temporary data is deleted. Finally when entering the next `with`-clause a new Java process is started. This can take several seconds.

If appropriate, it is safe to run `with TCPython()` in a loop. **Due to the time it takes this only makes sense if the calculation time per iteration is longer than a minute.**

To prevent calling `with TCPython()` multiple times and cleaning up temporary data, you can use the following pattern. **Example:**
from tc_python import *

# ...

def calculation(calculator):
    # you could also pass the 'session' or 'system' object if more appropriate
    calculator.set_condition("W(Cr)", 0.1)
    # further configuration ...

    result = calculator.calculate()
    # ...
    result.invalidate()  # if the temporary data needs to be cleaned up immediately

if __name__ == '__main__':
    with TCPython() as session:
        system = session.select_database_and_elements("FEDEMO", ["Fe", "Cr"])
        calculator = system.with_single_equilibrium_calculation()

        for i in range(50):
            calculation(calculator)

4.6 Parallel calculations

It is possible to perform parallel calculations with TC-Python using multi-processing.

Note: Please note that multi-threading is not suitable for parallelization of computationally intensive tasks in Python. Additionally the Thermo-Calc core is not thread-safe. Using suitable Python-frameworks it is also possible to dispatch the calculations on different computers of a cluster.

A general pattern that can be applied is shown below. This code snippet shows how to perform single equilibrium calculations for different compositions in parallel. In the same way all other calculators of Thermo-Calc can be used or combined. For performance reasons in a real application, probably numpy arrays instead of Python arrays should be used.

Example:

import concurrent.futures

from tc_python import *

def do_perform(parameters):
    # this function runs within an own process
    with TCPython() as start:
        elements = ["Fe", "Cr", "Ni", "C"]
        calculation = (start.select_database_and_elements("FEDEMO", elements).get_system().
                       with_single_equilibrium_calculation().
                       set_condition("T", 1100).
                       set_condition("W(C)", 0.1 / 100).
                       set_condition("W(Ni)", 2.0 / 100))

(continues on next page)
phase_fractions = []
cr_contents = range(parameters["cr_min"],
                     parameters["cr_max"],
                     parameters["delta_cr"])

for cr in cr_contents:
    result = (calculation.
              set_condition("W(Cr)", cr / 100).
              calculate())

    phase_fractions.append(result.get_value_of("NPM(BCC_A2)"))

return phase_fractions

if __name__ == "__main__":
    parameters = [
        {"index": 0, "cr_min": 10, "cr_max": 15, "delta_cr": 1},
        {"index": 1, "cr_min": 15, "cr_max": 20, "delta_cr": 1}
    ]

    bcc_phase_fraction = []
    num_processes = 2

    with concurrent.futures.ProcessPoolExecutor(num_processes) as executor:
        for result_from_process in zip(parameters, executor.map(do_perform, parameters)):
            # params can be used to identify the process and its parameters
            params, phase_fractions_from_process = result_from_process
            bcc_phase_fraction.extend(phase_fractions_from_process)

            # use the result in `bcc_phase_fraction`, for example for plotting

4.7 Handling crashes of the calculation engine

In some cases the Thermo-Calc calculation engine can crash. If batch calculations are performed, this brings down the complete batch. To handle this situation there is an exception you can use.

UnrecoverableCalculationException

That exception is thrown if the calculation server enters a state where no further calculations are possible. You should catch that exception outside of the with TCPython() clause and continue within a new with-clause.

Example:

from tc_python import *

for temperature in range(900, 1100, 10):
    try:
        with TCPython() as start:
            diffusion_result = (start.
                                 select_thermodynamic_and_kinetic_databases_with_elements("FEDEMO", "MFEDEMO", ["Fe", "Ni"]).

(continues on next page)
get_system().
with_isothermal_diffusion_calculation().
    set_temperature(temperature).
    set_simulation_time(108000.0).
    add_region(Region("Austenite").
        set_width(1E-4).
        with_grid(CalculatedGrid.linear().set_no_of_points(50)).
        with_composition_profile(CompositionProfile().
            add("Ni", ElementProfile.linear(10.0, 50.0))
        ).
    add_phase("FCC_A1")).
    calculate()}

    distance, ni_fraction = diffusion_result.get_mass_fraction_of_component_at_time("Ni", 108000.0)
    print(ni_fraction)

except UnrecoverableCalculationException as e:
    print('Could not calculate. Continuing with next...')

4.8 Using TC-Python within a Jupyter Notebook or the Python console

TC-Python can also be used from within an interactive Jupyter Notebook and a Python console as well as similar products. The main difference from a regular Python program is that it is not recommended to use a with-clause to manage the TC-Python resources. That is only possible within a single Jupyter Notebook cell. Instead the standalone functions `tc_python.server.start_api_server()` and `tc_python.server.stop_api_server()` should be used for manually managing the resources.

**Note:** The resources of TC-Python are primarily the Java-process running on the backend side that performs the actual calculations and the temporary-directory of TC-Python that can grow to a large size over time, especially if precipitation calculations are performed. If a with-clause is used, these resources are automatically cleared after use.

You need to make sure that you execute the two functions `tc_python.server.start_api_server()` and `tc_python.server.stop_api_server()` exactly once within the Jupyter Notebook session. If not stopping TC-Python, extra Java-processes might be present and the temporary disk-space is not cleared. However, these issues can be resolved manually.

The temporary directories of TC-Python are named, for example, `TC_TMP474758848895385507` that has a random ID. The temporary directory on different operating systems varies according to the pattern shown in the table.

<table>
<thead>
<tr>
<th>Operating system</th>
<th>Temporary directory</th>
</tr>
</thead>
<tbody>
<tr>
<td>Windows</td>
<td>C:\Users{UserName}\AppData\Local\Temp\TC_TMP4747588488953835507</td>
</tr>
<tr>
<td>MacOS</td>
<td>/var/folders/g7/7du8lti_b7mm84n184fn3k910000lg/T/TC_TMP4747588488953835507</td>
</tr>
<tr>
<td>Linux</td>
<td>/tmp/TC_TMP4747588488953835507</td>
</tr>
</tbody>
</table>

In a Jupyter Notebook some features of an IDE such as auto-completion (TAB-key), available method lookup (press . and then TAB) and parameter lookup (set the cursor within the method-parenthesis and press SHIFT + TAB or SHIFT + TAB + TAB for the whole docstring) are also available.
Example using TC-Python with a Jupyter Notebook:

```python
In [1]: from tc_python import *
In [2]: start_api_server()
In [3]: system = Setup().select_database_and_elements("FREDERICK", ["Fe", "Ni", "Cr"]).set_system()
   calc = system_with_single_equilibrium_calculation()
In [4]: temp = 825 # in K
   Ni_conc = 10.0 # in wt-%
   Cr_conc = 8.0 # in wt-%
   calc.
   \set_condition(ThermodynamicQuantity, temperature(), temp).
   \set_condition(ThermodynamicQuantity, mass_fractions_of_a_component("Ni"), Ni_conc / 100).
   \set_condition(ThermodynamicQuantity, mass_fractions_of_a_component("Cr"), Cr_conc / 100)
   result = calc.calculate()
In [5]: result.get_value_of(ThermodynamicQuantity, mole_fractions_of_a_phase("FeC_Al"))
Out[5]: 0.3345580340424432
In [6]: stop_api_server()
```

4.9 Property Model Framework

4.9.1 Debugging Property Model code

You can debug property models while running them from Thermo-Calc.

- Start Thermo-Calc and create a Property Model calculator.
- Select the model you want to debug and check the debug checkbox in the lower right corner of the Python code tab.

Now the model that you want to debug has been updated with code needed to connect with Thermo-Calc.

- Start debugging the model in the IDE of your choice.

**Note:** You must use a Python interpreter where TC-Python is installed.

In PyCharm it looks like this:
Note: When your IDE and Thermo-Calc have successfully connected, you will see this in the Thermo-Calc log:

```
10:34:42,170 INFO Waiting for developer(!) to start Python process in debugger...

10:34:42,171 INFO Connected successfully to the Python process for the model 'DrivingForcePythonModel' in DEBUG mode
```

You can stop the debug session in your IDE, change the model code, and start debugging again. The changes you made will take effect in Thermo-Calc without the need to restart. If you for instance changed the method evaluate_model(), the change will take effect the next time you press **Perform**.

It is also possible to start the models from TC-Python. The workflow is exactly the same as described above, except instead of starting Thermo-Calc graphical user interface, you start a Python script and use the parameter debug_model=True when selecting your model.

```
from tc_python import *

with TCPython() as start:
    property_model = (start.
        select_database_and_elements("FEDEMO", ["Fe", "C"]).
        get_system().
        with_property_model_calculation("my own Driving Force", debug_model=True).
        set_composition("C", 1.0).
    )
    property_model.calculate()

...```
4.9.2 Developing Property Models in several files

You can split your Property Model code in several .py files, and there are two ways of doing that:

- *side-by-side* modules
- *common* modules

*Side-by-side* modules are Python files located in the same folder as the Property Model.
*Common* modules are Python files located in a folder outside of the Property Model folder, which makes it possible to share them with several models as a common library.

4.9.2.1 *side-by-side* modules

You are required to:

- Add a __init__.py file to your Property Model folder
- Add all imports of *side-by-side* modules in your main Property Model Python file also to the __init__.py file

Example:

CriticalTemperaturesPythonModel.py (The main Property Model file):

```python
from CriticalTemperaturesPython import CriticalTemperatures
from tc_python import *
import numpy as np

class CriticalTemperaturesPythonModel(PropertyModel):
...
```

__init__.py:

```python
from CriticalTemperaturesPython.critical_temperatures_library import
→CriticalTemperatures
```

If you are using PyCharm, the package name of the Property Model might be highlighted as an error, in this case you can mark the Property Model directory (i.e. the root of the present model directory) by right-clicking on it in the project window of PyCharm and marking it as Sources Root:

```python
from tc_python import *
import numpy as np
from scipy import optimize
from enum import Enum

class CriticalTemperatures(object):
...
```

4.9. Property Model Framework
Note: Modules installed in the Python interpreter such as numpy, scipy, etc can be imported as normal. This only concerns files imported as side-by-side modules.

4.9.2.2 common modules

*common* modules work very similar to side-by-side modules except the import statements are done in the “main” __init__.py file in *Property Model directory*.

You are required to:

- Add a __init__.py file to your property model folder.
- Add all imports of *common* modules in your main property model python file also to both the __init__.py file in *Property Model directory* AND the __init__.py of the property model.

Example:

CriticalTemperaturesPythonModel.py (The main Property Model file):

```python
from PropertyModels import Martensite
from tc_python import *

class CriticalTemperaturesPythonModel(PropertyModel):
...
```

__init__.py: (The init file located in the property model folder)

```python
from PropertyModels import Martensite
```

__init__.py: (The init file located in *Property Model directory*)

```python
from PropertyModels.common.martensite_library import Martensite
```

The file critical_temperatures_library.py should in this example be located in a folder called *common* in the *Property Model directory*.

critical_temperatures_library.py:

```python
from tc_python import *
import numpy as np
from scipy import optimize
from enum import Enum

class CriticalTemperatures(object):
...
```

Note: *common* modules don’t have to be located in folder called *common*. It can be any name, as long as the imports match the folder name.

Note: In this example the *Property Model directory* is called ‘PropertyModels’. If you use a different directory for your property model, your imports have to match that.
4.9.3 Alternative Python for Property Models

4.9.3.1 Default bundled Python interpreter

Thermo-Calc is by default using a Python 3.7.2 interpreter bundled to the software for running the property models. It is containing the following major packages:

<table>
<thead>
<tr>
<th>Package</th>
<th>Version</th>
</tr>
</thead>
<tbody>
<tr>
<td>matplotlib</td>
<td>3.3.2</td>
</tr>
<tr>
<td>numpy</td>
<td>1.19.2</td>
</tr>
<tr>
<td>scikit-learn</td>
<td>0.23.2</td>
</tr>
<tr>
<td>scipy</td>
<td>1.5.2</td>
</tr>
<tr>
<td>TC-Python</td>
<td>2022a</td>
</tr>
</tbody>
</table>

**Warning:** Any changes to the interpreter packages can therefore break Thermo-Calc and should be avoided. If the installation gets broken, it can be fixed by reinstalling Thermo-Calc after having removed it.

Please contact the Thermo-Calc support if you think that further packages might be useful in future releases. If these packages are insufficient for you, it is possible to use another Python-interpreter: Configuring another Python interpreter.

The interpreter is located in different places depending on the platform:

<table>
<thead>
<tr>
<th>Operating system</th>
<th>Path to the bundled Python-interpreter</th>
</tr>
</thead>
<tbody>
<tr>
<td>Windows</td>
<td>C:\Program Files\Thermo-Calc\2022a\python\python.exe</td>
</tr>
<tr>
<td>Linux</td>
<td>/home/UserName/Thermo-Calc/2022a/python/bin/python3</td>
</tr>
<tr>
<td>MacOS</td>
<td>/Applications/Thermo-Calc-2022a.app/Contents/Resources/python/bin/python3</td>
</tr>
</tbody>
</table>

4.9.3.2 Configuring another Python interpreter

If you require additional Python-packages or prefer to use your own interpreter installed on your system, you can change the interpreter used by Thermo-Calc to run the property models. Select **Tools→Options** in the Thermo-Calc GUI and modify the path to that of your Python 3 interpreter of choice:
5.1 Calculations

5.1.1 Module “single_equilibrium”

class tc_python.single_equilibrium.AbstractSingleEquilibriumCalculation(calculator)
    Bases: tc_python.abstract_base.AbstractCalculation

Abstract configuration required for a single equilibrium calculation.

Note: This is an abstract class that cannot be used directly.

    abstract calculate() → tc_python.single_equilibrium.SingleEquilibriumTempResult

def disable_global_minimization()
    Turns the global minimization completely off.

    Returns This SingleEquilibriumCalculation object

def enable_global_minimization()
    Turns the global minimization on (using the default settings).

    Returns This SingleEquilibriumCalculation object

def get_components() → List[str]
    Returns a list of components in the system (including all components auto-selected by the database(s)).

    Returns The components

def get_gibbs_energy_addition_for(phase: str) → float
    Used to get the additional energy term (always being a constant) of a given phase. The value given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

    It is not composition-, temperature- or pressure-dependent.

    Parameters phase – Specify the name of the (stoichiometric or solution) phase with the addition

    Returns Gibbs energy addition to G per mole formula unit.

def get_system_data() → tc_python.abstract_base.SystemData
    Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using with_system_modifications().
Run a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

**Parameters**
- `command` – The Thermo-Calc Console Mode command

**Returns**
This `SingleEquilibriumCalculation` object

**Note:** It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

**Warning:** As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).

Sets the specified component to the status ENTERED, that is the default state.

**Parameters**
- `component` – The component name or `ALL_COMPONENTS`

**Returns**
This `SingleEquilibriumCalculation` object

Sets the specified component to the status SUSPENDED, i.e. it is ignored in the calculation.

**Parameters**
- `reset_conditions` – if ‘True’ also remove composition conditions for the component if they are defined
- `component` – The component name or `ALL_COMPONENTS`

**Returns**
This `SingleEquilibriumCalculation` object

Used to specify the additional energy term (always being a constant) of a given phase. The value (`gibbs_energy`) given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters**
- `phase` – Specify the name of the (stoichiometric or solution) phase with the addition
- `gibbs_energy` – Addition to G per mole formula unit

**Returns**
This `SingleEquilibriumCalculation` object

Sets the phase to the status DORMANT, necessary for calculating the driving force to form the specified phase.

**Parameters**
- `phase` – The phase name or `ALL_PHASES` for all phases

**Returns**
This `SingleEquilibriumCalculation` object
**set_phase_to_entered** *(phase: str, amount: float = 1.0)*

Sets the phase to the status ENTERED, that is the default state.

**Parameters**

- **phase** – The phase name or ALL_PHASES for all phases
- **amount** – The phase fraction (between 0.0 and 1.0)

**Returns** This **SingleEquilibriumCalculation** object

**set_phase_to_fixed** *(phase: str, amount: float)*

Sets the phase to the status FIXED, i.e. it is guaranteed to have the specified phase fraction after the calculation.

**Parameters**

- **phase** – The phase name
- **amount** – The fixed phase fraction (between 0.0 and 1.0)

**Returns** This **SingleEquilibriumCalculation** object

**set_phase_to_suspended** *(phase: str)*

Sets the phase to the status SUSPENDED, i.e. it is ignored in the calculation.

**Parameters**

- **phase** – The phase name or ALL_PHASES for all phases

**Returns** This **SingleEquilibriumCalculation** object

**with_options** *(options: tc_python.single_equilibrium.SingleEquilibriumOptions)*

Sets the simulation options.

**Parameters**

- **options** – The simulation options

**Returns** This **SingleEquilibriumCalculation** object

**with_reference_state** *(component: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)*

The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.

All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

**Parameters**

- **component** – The name of the element must be given.
- **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.
- **temperature** – The Temperature (in K) for the reference state. Or CURRENT_TEMPERATURE which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.

- **pressure** – The Pressure (in Pa) for the reference state.

**Returns** This `SingleEquilibriumCalculation` object

with_system_modifications(system_modifications: tc_python.abstract_base.SystemModifications) Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

**Note:** This is only possible if the system has been read from unencrypted (i.e. user) databases loaded as a *.tdb*-file.

**Parameters** system_modifications – The system modification to be performed

**Returns** This `SingleEquilibriumCalculation` object

**class** tc_python.single_equilibrium.SingleEquilibriumCalculation(calculator)

**Bases:** tc_python.single_equilibrium.AbstractSingleEquilibriumCalculation

Configuration for a single equilibrium calculation.

**Note:** Specify the conditions and possibly other settings, the calculation is performed with `calculate()`.

**bookmark_state**(bookmark_id: str = '') → str Puts a “bookmark” on the current calculation-state of the calculator allowing the program to return to this state later as needed.

By bookmarking a state, you can simplify the convergence of equilibria when they strongly depend on the starting conditions (i.e. the state). Also use it to improve performance by running a calculation, then bookmarking it, and later returning to it for other equilibria whose conditions are “close” to the bookmarked equilibrium.

This method is used in combination with the method `set_state_to_bookmark()`.

**Parameters** bookmark_id – The bookmark id. If omitted a generated id is used and returned

**Returns** The bookmark id

**calculate**() → tc_python.single_equilibrium.SingleEquilibriumTempResult Performs the calculation and provides a temporary result object that is only valid until something gets changed in the calculation state. The method `calculate()` is the default approach and should be used in most cases.

**Returns** A new `SingleEquilibriumTempResult` object which can be used to get specific values from the calculated result. It is undefined behavior to use that object after the state of the calculation has been changed.

**Warning:** If the result object should be valid for the whole program lifetime, use `calculate_with_state()` instead.

**calculate_with_state**() → tc_python.single_equilibrium.SingleEquilibriumResult Performs the calculation and provides a result object that reflects the present state of the calculation during the whole lifetime of the object.
Note: Because this method has performance and temporary disk space overhead (i.e. it is resource heavy), only use it when it is necessary to access the result object after the state is changed. In most cases you should use the method `calculate()`.

Returns A new `SingleEquilibriumResult` object which can be used later at any time to get specific values from the calculated result.

disable_global_minimization()
Turns the global minimization completely off.

Returns This `SingleEquilibriumCalculation` object
enable_global_minimization()
Turns the global minimization on (using the default settings).

Returns This `SingleEquilibriumCalculation` object
get_components() → List[str]
Returns a list of components in the system (including all components auto-selected by the database(s)).

Returns The components
get_gibbs_energy_addition_for(phase: str) → float
Used to get the additional energy term (always being a constant) of a given phase. The value given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

Parameters phase – Specify the name of the (stoichiometric or solution) phase with the addition

Returns Gibbs energy addition to G per mole formula unit.

get_interfacial_energy(matrix_phase: str, precipitate_phases: List[str], zero_volume_elements: List[str] = ['C', 'N']) → Dict[str, float]
Estimates the interfacial energy between a matrix phase and a precipitate phase using thermodynamic data from a CALPHAD database. The approximation model is based on Becker’s bond energy approach.

Default: elements with no contribution to volume are C and N.

Parameters

• matrix_phase – The matrix phase.

• precipitate_phases – The list of precipitate phases for which interfacial energy between them and the matrix phase is to be calculated.

• zero_volume_elements – The elements that are assumed to not contribute to the volume.

Returns A dictionary containing interfacial energy per precipitate phase.

get_system_data() → tc_python.abstract_base.SystemData
Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using `with_system_modifications()`.

Note: Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb-file.
Returns The system data

**remove_all_conditions()**
Removes all set conditions.

**Returns** This `SingleEquilibriumCalculation` object

**remove_condition(quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str])**
Removes the specified condition.

**Parameters**
- **quantity** – the thermodynamic quantity to set as condition; a Console Mode syntax string can be used as an alternative (for example “X(Cr)”)

**Returns** This `SingleEquilibriumCalculation` object

**run_poly_command(command: str)**
Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

**Parameters**
- **command** – The Thermo-Calc Console Mode command

**Returns** This `SingleEquilibriumCalculation` object

---

**Note:** It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

---

**Warning:** As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).

**set_component_to_entered(component: str)**
Sets the specified component to the status ENTERED, that is the default state.

**Parameters**
- **component** – The component name or `ALL_COMPONENTS`

**Returns** This `SingleEquilibriumCalculation` object

**set_component_to_suspended(component: str, reset_conditions: bool = False)**
Sets the specified component to the status SUSPENDED, i.e. it is ignored in the calculation.

**Parameters**
- **reset_conditions** – if ‘True’ also remove composition conditions for the component if they are defined
- **component** – The component name or `ALL_COMPONENTS`

**Returns** This `SingleEquilibriumCalculation` object

**set_condition(quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], value: float)**
Sets the specified condition.

**Parameters**
- **quantity** – The thermodynamic quantity to set as condition; a Console Mode syntax string can be used as an alternative (for example “X(Cr)”)
- **value** – The value of the condition

**Returns** This `SingleEquilibriumCalculation` object
set_gibbs_energy_addition_for(phase: str, gibbs_energy: float)

Used to specify the additional energy term (always being a constant) of a given phase. The value
(gibbs_energy) given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent
a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

Parameters

- phase – Specify the name of the (stoichiometric or solution) phase with the addition
- gibbs_energy – Addition to G per mole formula unit

Returns This SingleEquilibriumCalculation object

set_phase_to_dormant(phase: str)

Sets the phase to the status DORMANT, necessary for calculating the driving force to form the specified phase.

Parameters phase – The phase name or ALL_PHASES for all phases

Returns This SingleEquilibriumCalculation object

set_phase_to_entered(phase: str, amount: float = 1.0)

Sets the phase to the status ENTERED, that is the default state.

Parameters

- phase – The phase name or ALL_PHASES for all phases
- amount – The phase fraction (between 0.0 and 1.0)

Returns This SingleEquilibriumCalculation object

set_phase_to_fixed(phase: str, amount: float)

Sets the phase to the status FIXED, i.e. it is guaranteed to have the specified phase fraction after the calculation.

Parameters

- phase – The phase name
- amount – The fixed phase fraction (between 0.0 and 1.0)

Returns This SingleEquilibriumCalculation object

set_phase_to_suspended(phase: str)

Sets the phase to the status SUSPENDED, i.e. it is ignored in the calculation.

Parameters phase – The phase name or ALL_PHASES for all phases

Returns This SingleEquilibriumCalculation object

set_state_to_bookmark(bookmark_id: str) \(\rightarrow\) tc_python.single_equilibrium.SingleEquilibriumTempResult

Resets the calculation state to a previously bookmarked state.

After calling this method, the calculation behaves exactly as it would after the bookmarked calculation ran.

This method is used in combination with the method bookmark_state().

Parameters bookmark_id – The bookmark id of the state to return to.

Returns A new SingleEquilibriumTempResult object which can be used to get specific values from the calculated result. It is undefined behavior to use that object after the state of the calculation has been changed.
**with_options** *(options: tc_python.single_equilibrium.SingleEquilibriumOptions)*

Sets the simulation options.

Parameters **options** – The simulation options

Returns This `SingleEquilibriumCalculation` object

**with_reference_state** *(component: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)*

The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.

All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

Parameters

- **component** – The name of the element must be given.
- **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.
- **temperature** – The Temperature (in K) for the reference state. Or `CURRENT_TEMPERATURE` which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.
- **pressure** – The Pressure (in Pa) for the reference state.

Returns This `SingleEquilibriumCalculation` object

**with_system_modifications** *(system_modifications: tc_python.abstract_base.SystemModifications)*

Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

Parameters **system_modifications** – The system modification to be performed

Returns This `SingleEquilibriumCalculation` object

---

**Note:** This is only possible if the system has been read from unencrypted (i.e. `user`) databases loaded as a `*.tdb`-file.

---

**Parameters** **system_modifications** – The system modification to be performed

Returns This `SingleEquilibriumCalculation` object

---

**class** `tc_python.single_equilibrium.SingleEquilibriumOptions`

Bases: `object`

General simulation conditions for the thermodynamic calculations.

---

Chapter 5. API Reference
disable_approximate_driving_force_for_metastable_phases()
Disables the approximation of the driving force for metastable phases.

Default: Enabled

Note: When enabled, the metastable phases are included in all iterations. However, these may not have reached their most favorable composition and thus their driving forces may be only approximate.

If it is important that these driving forces are correct, use disable_approximate_driving_force_for_metastable_phases() to force the calculation to converge for the metastable phases.

Returns This SingleEquilibriumOptions object

disable_control_step_size_during_minimization()
Disables stepsize control during minimization (non-global).

Default: Enabled

Returns This SingleEquilibriumOptions object

disable_force_positive_definite_phase_hessian()
Disables forcing of positive definite phase Hessian. This determines how the minimum of an equilibrium state in a normal minimization procedure (non-global) is reached. For details, search the Thermo-Calc documentation for “Hessian minimization”.

Default: Enabled

Returns This SingleEquilibriumOptions object

enable_approximate_driving_force_for_metastable_phases()
Enables the approximation of the driving force for metastable phases.

Default: Enabled

Note: When enabled, the metastable phases are included in all iterations. However, these may not have reached their most favorable composition and thus their driving forces may be only approximate.

If it is important that these driving forces are correct, use disable_approximate_driving_force_for_metastable_phases() to force the calculation to converge for the metastable phases.

Returns This SingleEquilibriumOptions object

enable_control_step_size_during_minimization()
Enables stepsize control during normal minimization (non-global).

Default: Enabled

Returns This SingleEquilibriumOptions object

enable_force_positive_definite_phase_hessian()
Enables forcing of positive definite phase Hessian. This determines how the minimum of an equilibrium state in a normal minimization procedure (non-global) is reached. For details, search the Thermo-Calc documentation for “Hessian minimization”.

Default: Enabled
Returns This `SingleEquilibriumOptions` object

**set_global_minimization_max_grid_points** (*max_grid_points: int = 2000*)
Sets the maximum number of grid points in global minimization. Only applicable if global minimization is actually used.

Default: 2000 points

Parameters `max_grid_points` – The maximum number of grid points

Returns This `SingleEquilibriumOptions` object

**set_max_no_of_iterations** (*max_no_of_iterations: int = 500*)
Set the maximum number of iterations.

Default: max. 500 iterations

Note: As some models give computation times of more than 1 CPU second/iteration, this number is also used to check the CPU time and the calculation stops if 500 CPU seconds/iterations are used.

Parameters `max_no_of_iterations` – The max. number of iterations

Returns This `SingleEquilibriumOptions` object

**set_required_accuracy** (*accuracy: float = 1e-06*)
Sets the required relative accuracy.

Default: 1.0E-6

Note: This is a relative accuracy, and the program requires that the relative difference in each variable must be lower than this value before it has converged. A larger value normally means fewer iterations but less accurate solutions. The value should be at least one order of magnitude larger than the machine precision.

Parameters `accuracy` – The required relative accuracy

Returns This `SingleEquilibriumOptions` object

**set_smallest_fraction** (*smallest_fraction: float = 1e-12*)
Sets the smallest fraction for constituents that are unstable.

It is normally only in the gas phase that you can find such low fractions.

The default value for the smallest site-fractions is 1E-12 for all phases except for IDEAL phase with one sublattice site (such as the GAS mixture phase in many databases) for which the default value is always as 1E-30.

Parameters `smallest_fraction` – The smallest fraction for constituents that are unstable

Returns This `SingleEquilibriumOptions` object

**class** `tc_python.single_equilibrium.SingleEquilibriumResult` (*result*)
Bases: `tc_python.abstract_base.AbstractResult`

Result of a single equilibrium calculation, it can be evaluated using a Quantity or Console Mode syntax.

**change_pressure** (*pressure: float*)
Change the pressure and re-evaluate the results from the equilibrium without minimizing Gibbs energy, i.e. with higher performance. The properties are calculated at the new pressure using the phase amount,
temperature and composition of phases from the initial equilibrium. Use \texttt{get\_value\_of()} to obtain
them.

\textbf{Parameters pressure} – The pressure [Pa]

\textbf{Returns} This \texttt{SingleEquilibriumCalculation} object

\textbf{change\_temperature(temperature: float)}

Change the temperature and re-evaluate the results from the equilibrium without minimizing Gibbs energy,
\textit{i.e.} with high performance. The properties are calculated at the new temperature using the phase amount,
pressure and composition of phases from the initial equilibrium. Use \texttt{get\_value\_of()} to obtain them.

\textbf{Note:} This is typically used when calculating room temperature properties (e.g. density) for a material
when it is assumed that the equilibrium phase amount and composition freeze-in at a higher temperature
during cooling.

\textbf{Parameters temperature} – The temperature [K]

\textbf{Returns} This \texttt{SingleEquilibriumCalculation} object

\textbf{get\_components()} \rightarrow \texttt{List[str]}

Ret\texttt{urns} the names of the components selected in the system (including any components auto-selected by
the database(s)).

\textbf{Returns} The names of the selected components

\textbf{get\_conditions()} \rightarrow \texttt{List[str]}

Ret\texttt{urns} the conditions.

\textbf{Returns} The selected conditions

\textbf{get\_phases()} \rightarrow \texttt{List[str]}

Ret\texttt{urns} the phases present in the system due to its configuration. It also contains all phases that
have been automatically added during the calculation, this is the difference to the method \texttt{System.get\_phases\_in\_system()}.

\textbf{Returns} The names of the phases in the system including automatically added phases

\textbf{get\_stable\_phases()} \rightarrow \texttt{List[str]}

Ret\texttt{urns} the stable phases (i.e. the phases present in the current equilibrium).

\textbf{Returns} The names of the stable phases

\textbf{get\_value\_of(quantity: Union[tc\_python.quantity\_factory.ThermodynamicQuantity, str])} \rightarrow \texttt{float}

Ret\texttt{urns} a value from a single equilibrium calculation.

\textbf{Parameters quantity} – The thermodynamic quantity to get the value of; a Console Mode
syntax strings can be used as an alternative (for example “NPM(FCC_A1)"

\textbf{Returns} The requested value

\textbf{run\_poly\_command(command: str)}

Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine. This
affects only the state of the result object.

\textbf{Parameters command} – The Thermo-Calc Console Mode command

\textbf{Returns} This \texttt{SingleEquilibriumCalculation} object
**Note:** It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

**Warning:** As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).

```python
save_to_disk(path: str)
```

Saves the result to disk. Note that the result is a folder, containing potentially many files. The result can later be loaded with `load_result_from_disk()`.

**Parameters**
- `path` – the path to the folder you want the result to be saved in. It can be relative or absolute.

**Returns**
- This `SingleEquilibriumResult` object

```python
class tc_python.single_equilibrium.SingleEquilibriumTempResult(result)
```

**Bases:** `tc_python.abstract_base.AbstractResult`

Result of a single equilibrium calculation that is only valid until something gets changed in the calculation state. It can be evaluated using a Quantity or Console Mode syntax.

**Warning:** Note that it is undefined behavior to use that object after something has been changed in the state of the calculation, this will result in an `InvalidResultStateException` exception being raised.

```python
change_pressure(pressure: float)
```

Change the pressure and re-evaluate the results from the equilibrium without minimizing Gibbs energy, i.e. with higher performance. The properties are calculated at the new pressure using the phase amount, temperature and composition of phases from the initial equilibrium. Use `get_value_of()` to obtain them.

**Parameters**
- `pressure` – The pressure [Pa]

**Returns**
- This `SingleEquilibriumCalculation` object

```python
change_temperature(temperature: float)
```

Change the temperature and re-evaluate the results from the equilibrium without minimizing Gibbs energy, i.e. with high performance. The properties are calculated at the new temperature using the phase amount, pressure and composition of phases from the initial equilibrium. Use `get_value_of()` to obtain them.

**Note:** This is typically used when calculating room temperature properties (e.g. density) for a material when it is assumed that the equilibrium phase amount and composition freeze-in at a higher temperature during cooling.

**Parameters**
- `temperature` – The temperature [K]

**Returns**
- This `SingleEquilibriumCalculation` object

```python
get_components() \rightarrow List[str]
```

Returns the names of the components selected in the system (including any components auto-selected by the database(s)).

**Returns**
- The names of the selected components
Raises **InvalidResultStateException** – If something has been changed in the state of the calculation since that result object has been created

**get_conditions** () → List[str]
Returns the conditions.

**Returns** List containing the selected conditions

Raises **InvalidResultStateException** – If something has been changed in the state of the calculation since that result object has been created

**get_phases** () → List[str]
Returns the phases present in the system due to its configuration. It also contains all phases that have been automatically added during the calculation, this is the difference to the method **System.get_phases_in_system**.

**Returns** The names of the phases in the system including automatically added phases

Raises **InvalidResultStateException** – If something has been changed in the state of the calculation since that result object has been created

**get_stable_phases** () → List[str]
Returns the stable phases (i.e. the phases present in the current equilibrium).

**Returns** The names of the stable phases

Raises **InvalidResultStateException** – If something has been changed in the state of the calculation since that result object has been created

**get_value_of** *(quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str]) → float*
Returns a value from a single equilibrium calculation.

**Parameters** **quantity** – The thermodynamic quantity to get the value of; a Console Mode syntax strings can be used as an alternative (for example “NPM(FCC_A1)"

**Returns** The requested value

Raises **InvalidResultStateException** – If something has been changed in the state of the calculation since that result object has been created

**run_poly_command** *(command: str)*
Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

**Parameters** **command** – The Thermo-Calc Console Mode command

**Returns** This **SingleEquilibriumCalculation** object

**Note:** It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

**Warning:** As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).
5.1.2 Module “batch_equilibrium”

class tc_python.batch_equilibrium.BatchEquilibriumCalculation(calculator)
    Bases: tc_python.abstract_base.AbstractCalculation

Configuration for a series of single equilibrium calculations performed in a vectorized fashion.

**Note:** Specify the conditions and call `calculate()`.

**Tip:** The performance of batch equilibrium calculations can be significantly better than looping and using `SingleEquilibriumCalculation` if the actual Thermo-Calc calculation is fast. There is little advantage if the Thermo-Calc equilibrium calculations take a long time (typically for large systems and databases).

calculate(quantities: List[Union[tc_python.quantity_factory.ThermodynamicQuantity, str]], logging_frequency: int = 10) → tc_python.batch_equilibrium.BatchEquilibriumResult

Runs the batch equilibrium calculation. The calculated `BatchEquilibriumResult` can then be queried for the values of the quantities specified.

Example:

```python
gg quantities = ['G', 'X(BCC)']
```

**Parameters** `logging_frequency` – Determines how often logging should be done.

**Returns** `A BatchEquilibriumResult` which later can be used to get specific values from the calculated result.

disable_global_minimization()

Returns This `BatchEquilibriumCalculation` object

enable_global_minimization()

Returns This `BatchEquilibriumCalculation` object

get_components() → List[str]

Returns a list of components in the system (including all components auto-selected by the database(s)).

**Returns** The components

get_gibbs_energy_addition_for(phase: str) → float

Used to get the additional energy term (always being a constant) of a given phase. The value given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters** `phase` – Specify the name of the (stoichiometric or solution) phase with the addition

**Returns** Gibbs energy addition to G per mole formula unit.

get_system_data() → tc_python.abstract_base.SystemData

Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using `with_system_modifications()`.
Note: Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb-file.

Returns The system data

remove_all_conditions()  
Removes all set conditions.

Returns This BatchEquilibriumCalculation object

remove_condition(quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str])  
Removes the specified condition.

Parameters quantity – the thermodynamic quantity to set as condition; a Console Mode syntax string can be used as an alternative (for example “X(Cr)”)

Returns This BatchEquilibriumCalculation object

run_poly_command(command: str)  
Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

Parameters command – The Thermo-Calc Console Mode command

Returns This BatchEquilibriumCalculation object

Note: It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

Warning: As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).

set_component_to_entered(component: str)  
Sets the specified component to the status ENTERED, that is the default state.

Parameters component – The component name or ALL_COMPONENTS

Returns This BatchEquilibriumCalculation object

set_component_to_suspended(component: str, reset_conditions: bool = False)  
Sets the specified component to the status SUSPENDED, i.e. it is ignored in the calculation.

Parameters

• reset_conditions – if ‘True’ also remove composition conditions for the component if they are defined

• component – The component name or ALL_COMPONENTS

Returns This BatchEquilibriumCalculation object

set_condition(quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], value: float)  
Sets the specified condition.

Parameters

• quantity – The thermodynamic quantity to set as condition; a Console Mode syntax string can be used as an alternative (for example “X(Cr)”)

5.1. Calculations 49
• **value** – The value of the condition

**Returns**  This `BatchEquilibriumCalculation` object

**set_conditions_for_equilibria**(equilibria: List[List[Union[tc_python.quantity_factory.ThermodynamicQuantity, str], float]])

Set the conditions of the equilibria to be calculated.

This is done by sending a list of equilibria at once.

Each equilibrium itself is a list of conditions that will be changed for that equilibrium.

A condition is described by a tuple containing:

1. A Console Mode syntax string or a `ThermodynamicQuantity` instance,
2. A float value specifying the value of the condition.

Example:

```python
>>> [[('T', 800), ('X(Cr)', 0.1)], [('T', 850), ('X(Cr)', 0.11)]
```

You can use `ThermodynamicQuantity` instead of a Console Mode syntax string when specifying type of condition.

Example:

```python
>>> [[[ThermodynamicQuantity.temperature(), 800), (ThermodynamicQuantity.mole_fraction_of_a_component('Cr'), 0.1)], [(ThermodynamicQuantity.temperature(), 850), (ThermodynamicQuantity.mole_fraction_of_a_component('Cr'), 0.15)]
```

**Parameters**  **equilibria** – The list of equilibria

**Returns**  This `BatchEquilibriumCalculation` object

**set_gibbs_energy_addition_for**(phase: str, gibbs_energy: float)

Used to specify the additional energy term (always being a constant) of a given phase. The value (`gibbs_energy`) given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters**

- **phase** – Specify the name of the (stoichiometric or solution) phase with the addition
- **gibbs_energy** – Addition to G per mole formula unit

**Returns**  This `BatchEquilibriumCalculation` object

**set_phase_to_dormant**(phase: str)

Sets the phase to the status DORMANT, necessary for calculating the driving force to form the specified phase.

**Parameters**  **phase** – The phase name or `ALL_PHASES` for all phases

**Returns**  This `BatchEquilibriumCalculation` object

**set_phase_to_entered**(phase: str, amount: float = 1.0)

Sets the phase to the status ENTERED, that is the default state.

**Parameters**

- **phase** – The phase name or `ALL_PHASES` for all phases
• **amount** – The phase fraction (between 0.0 and 1.0)

**Returns** This *BatchEquilibriumCalculation* object

**set_phase_to_fixed**(phase: str, amount: float)
Sets the phase to the status FIXED, i.e. it is guaranteed to have the specified phase fraction after the calculation.

**Parameters**

• **phase** – The phase name

• **amount** – The fixed phase fraction (between 0.0 and 1.0)

**Returns** This *BatchEquilibriumCalculation* object

**set_phase_to_suspended**(phase: str)
Sets the phase to the status SUSPENDED, i.e. it is ignored in the calculation.

**Parameters**

• **phase** – The phase name or *ALL_PHASES* for all phases

**Returns** This *BatchEquilibriumCalculation* object

**with_options**(options: tc_python.single_equilibrium.SingleEquilibriumOptions)
Sets the simulation options.

**Parameters**

• **options** – The simulation options

**Returns** This *BatchEquilibriumCalculation* object

**with_reference_state**(component: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)
The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.

All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

**Parameters**

• **component** – The name of the element must be given.

• **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.

• **temperature** – The Temperature (in K) for the reference state. Or CURRENT_TEMPERATURE which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.

• **pressure** – The Pressure (in Pa) for the reference state.
Returns This BatchEquilibriumCalculation object

with_system_modifications (system_modifications: tc_python.abstract_base.SystemModifications)
Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

Note: This is only possible if the system has been read from unencrypted (i.e. *user*) databases loaded as a *.tdb*-file.

Parameters system_modifications – The system modification to be performed

Returns This BatchEquilibriumCalculation object

class tc_python.batch_equilibrium.BatchEquilibriumResult (result)
Bases: object

Result of a batch equilibrium calculation. This can be used to query for specific values.

get_values_of (quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str]) → List[float]
Returns values from a batch equilibrium calculation.

Warning: The quantity must be one of the quantities specified for the BatchEquilibriumCalculation object that created the result object.

Example:

```python
>>> batch_result = batch_calculation.calculate(quantities = ['G', 'X(BCC)'])
>>> batch_result.get_values_of('G')
```

Parameters quantity – the thermodynamic quantity to get the value of; a Console Mode syntax strings can be used as an alternative (for example “NPM(FCC_A1)"

invalidate ()
Invalidates the object and frees the disk space used by it.

Note: This is only required if the disk space occupied by the object needs to be released during the calculation. No data can be retrieved from the object afterwards.

5.1.3 Module “precipitation”

class tc_python.precipitation.FixedGrainSize (grain_radius: float = 0.0001)
Bases: tc_python.precipitation.GrainGrowthModel

set_grain_aspect_ratio (grain_aspect_ratio: float = 1.0)
Enter a numerical value. Default: 1.0.

Parameters grain_aspect_ratio – The grain aspect ratio [-]

class tc_python.precipitation.GrainGrowth (grain_size_distribution: tc_python.precipitation.GrainSizeDistribution)
Bases: tc_python.precipitation.GrainGrowthModel
disable_zener_pinning()

Disable Zener pinning to ignore the particle pinning effect on the grain growth. Zener pinning is by default disabled when no grain size distribution is defined, i.e. a single constant grain size is used. The setting is by default enabled when a grain size distribution is defined.

**Returns** This GrainSizeDistribution object

enable_zener_pinning()

Enable Zener pinning to simulate the particle pinning effect on the grain growth. The setting is by default enabled when a grain size distribution is defined.

**Returns** This GrainSizeDistribution object

set_grain_boundary_energy(energy: float = 0.5)

Set the energy of the grain boundary.

**Parameters** energy – The grain boundary energy [J/m2]

**Returns** This GrainSizeDistribution object

set_grain_boundary_mobility_activation_energy(activation_energy: float = 242000.0)

Set the grain boundary mobility activation energy where the mobility is defined by an Arrhenius type of equation.

**Parameters** activation_energy – The mobility activation energy [J/mol]

**Returns** This GrainSizeDistribution object

set_grain_boundary_mobility_pre_factor(pre_factor: float = 0.004)

Set the grain boundary mobility prefactor where the mobility is defined by an Arrhenius type of equation.

**Parameters** pre_factor – The grain boundary mobility pre factor [m^4/(J s)]

**Returns** This GrainSizeDistribution object

class tc_python.precipitation.GrainGrowthModel

Factory class providing objects representing a grain growth model.

classmethod fixed_grain_size(grain_radius: float = 0.0001)

Fixed grain radius size. **Default**: 1.0E-4 m

**Parameters** grain_radius – The grain radius / size [m]

classmethod grain_growth(grain_size_distribution: tc_python.precipitation.GrainSizeDistribution)

Sets the initial grain size distribution for the matrix. **Default**: If the initial grain size distribution is not explicitly provided, a constant average grains size will be used and no grain growth evaluated during the simulation.

**Tip:** Use this option if you want to study the further evolution of an existing microstructure.

**Parameters** grain_size_distribution – grain size distribution

class tc_python.precipitation.GrainSizeDistribution

Represents the grain size distribution at a certain time.

add_radius_and_number_density(radius: float, number_density: float)

Adds a radius and number density pair to the grain size distribution.

5.1. Calculations 53
Parameters

- **radius** – The radius [m]
- **number_density** – The number of grains per unit volume per unit length [m^-4]

Returns This GrainSizeDistribution object

class tc_python.precipitation.GrowthRateModel(value)
Bases: enum.Enum

Choice of the used growth rate model for a precipitate.

The most efficient model is the Simplified model, which is the default and applicable to most alloy systems under the assumption that either the supersaturation is small, or the alloying elements have comparable diffusivity. If all alloying elements are substitutional but they have remarkable diffusivity difference, e.g. in Al-Zr system, or if the diffusivity is strongly composition-dependent, the General model is preferred. If the supersaturation is high, and meanwhile there are fast-diffusing interstitial elements such as C, the Advanced model is more appropriate to capture the NPLE mechanism.

ADVANCED = 3

The advanced model has been proposed by Chen, Jeppsson, and Ågren (CJA) (2008) and calculates the velocity of a moving phase interface in multicomponent systems by identifying the operating tie-line from the solution of the flux-balance equations. This model can treat both high supersaturation and cross-diffusion rigorously. Spontaneous transitions between different modes (LE and NPLE) of phase transformation can be captured without any ad-hoc treatment.

Note: Since it is not always possible to solve the flux-balance equations and it takes time, usage of a less rigorous but simple and efficient model is preferred if possible.

GENERAL = 5

The general model is based on the Morral-Purdy model, which follows the same quasi-steady state approximation as the Simplified model, but improves it by taking the cross-diffusion into account.

NPLE = 11

The Non-Partitioning Local Equilibrium (NPLE) growth rate model is only available for alloy systems where Fe is the major element and at least one interstitial element partitions into the precipitate phase. This model is specifically designed to deal with the fast diffusion of interstitial elements (C, N, etc.) in Fe alloys. Based on the Simplified growth model, it still holds a local equilibrium condition at the migrating interface. It chooses a tie-line under NPLE condition so that the u-fractions of all substitutional elements and minor interstitial elements in the precipitate phase are the same as those in the far-field matrix phase (i.e. the overall instantaneous matrix composition).

PARA_EQ = 10

The para-equilibrium model is only available for alloy systems where Fe is the major element and C is the only interstitial element, which also partitions into the precipitate phase. The interstitial elements, e.g. C, N, etc., usually have remarkably faster diffusion rate than the substitutional elements. Meanwhile, they are assumed to have negligible volume contribution, and as a result the composition variables are replaced by u-fractions when interstitial elements are included in the system. This model is specifically designed to address the fast diffusion of C in Fe alloys. Based on the Simplified growth rate model it holds a para-equilibrium condition at the migrating interface. Contrary to the regular ortho-equilibrium condition state that assumes that all alloying elements are in equilibrium at the interface, the para-equilibrium assumes only equilibrium for C. The substitutional elements are immobile and thus have the same compositions (u-fractions) across the interface.

SIMPLIFIED = 2

The simplified model is based on the advanced model but avoids the difficulty of finding the operating tie-line and uses instead the tie-line across the bulk composition. This is the default growth rate model.
class `tc_python.precipitation.MatrixPhase(matrix_phase_name: str)`

Bases: `object`

The matrix phase in a precipitation calculation

`add_precipitate_phase(precipitate_phase: tc_python.precipitation.PrecipitatePhase)`

Adds a precipitate phase.

**Parameters**

- `precipitate_phase` – The precipitate phase

`set_dislocation_density(dislocation_density: float = 5000000000000.0)`

Enter a numerical value. **Default**: 5.0E12 m^-2.

**Parameters**

- `dislocation_density` – The dislocation density [m^-2]

`set_grain_aspect_ratio(grain_aspect_ratio: float = 1.0)`

Enter a numerical value. **Default**: 1.0.

**Note:** Deprecated in version 2022a: Use `with_grain_growth_model()` instead. This method will be removed in release 2023a.

**Parameters**

- `grain_aspect_ratio` – The grain aspect ratio [-]

`set_grain_radius(grain_radius: float = 0.0001)`

Sets grain radius / size. **Default**: 1.0E-4 m

**Note:** Deprecated in version 2022a: Use `with_grain_growth_model()` instead. This method will be removed in release 2023a.

**Parameters**

- `grain_radius` – The grain radius / size [m]

`set_mobility_enhancement_activation_energy(mobility_enhancement_activation_energy: float = 0.0)`

A value that adds to the activation energy of mobility data from the database. **Default**: 0.0 J/mol

**Parameters**

- `mobility_enhancement_activation_energy` – The value that adds to the activation energy of mobility data from the database [J/mol].

`set_mobility_enhancement_prefactor(mobility_enhancement_prefactor: float = 1.0)`

A parameter that multiplies to the mobility data from database. **Default**: 1.0

**Parameters**

- `mobility_enhancement_prefactor` – The mobility enhancement factor [-]

`set_molar_volume(volume: float)`

Sets the molar volume of the phase.

**Default:** If not set, the molar volume is taken from the thermodynamic database (or set to 7.0e-6 m^3/mol if the database contains no molar volume information).

**Parameters**

- `volume` – The molar volume [m^3/mol]

`with_elastic_properties_cubic(c11: float, c12: float, c44: float)`

Sets the elastic properties to “cubic” and specifies the elastic stiffness tensor components. **Default**: if not chosen, the default is DISREGARD

**Parameters**

- `c11` – The stiffness tensor component c11 [GPa]
• **c12** – The stiffness tensor component c12 [GPa]

• **c44** – The stiffness tensor component c44 [GPa]

**with_elastic_properties_disregard()**
Set to disregard to ignore the elastic properties. **Default**: This is the default option

**with_elastic_properties_isotropic**(shear_modulus: float, poisson_ratio: float)
Sets elastic properties to isotropic. **Default**: if not chosen, the default is DISREGARD

**Parameters**

- **shear_modulus** – The shear modulus [GPa]
- **poisson_ratio** – The Poisson’s ratio [-]

**with_grain_growth_model**(grain_growth_model: tc_python.precipitation.GrainGrowthModel)
Sets the model for grain growth. Either fixed size or with a starting distribution

**Default**: Fixed grain radius size 1.0E-4 m

**Parameters**

- **grain_growth_model** – the grain growth model

---

**class** tc_python.precipitation.NumericalParameters

**Bases**: object

Numerical parameters

**set_max_overall_volume_change** (max_overall_volume_change: float = 0.001)
This defines the maximum absolute (not ratio) change of the volume fraction allowed during one time step.

**Default**: 0.001

**Parameters**

- **max_overall_volume_change** – The maximum absolute (not ratio) change of the volume fraction allowed during one time step [-]

**set_max_radius_points_per_magnitude** (max_radius_points_per_magnitude: float = 200.0)
Sets the maximum number of grid points over one order of magnitude in radius.

**Default**: 200.0

**Parameters**

- **max_radius_points_per_magnitude** – The maximum number of grid points over one order of magnitude in radius [-]

**set_max_rel_change_critical_radius** (max_rel_change_critical_radius: float = 0.1)
Used to place a constraint on how fast the critical radius can vary, and thus put a limit on time step.

**Default**: 0.1

**Parameters**

- **max_rel_change_critical_radius** – The maximum relative change of the critical radius [-]

**set_max_rel_change_nucleation_rate_log** (max_rel_change_nucleation_rate_log: float = 0.5)
This parameter ensures accuracy for the evolution of effective nucleation rate.

**Default**: 0.5

**Parameters**

- **max_rel_change_nucleation_rate_log** – The maximum logarithmic relative change of the nucleation rate [-]

**set_max_rel_radius_change** (max_rel_radius_change: float = 0.01)
The maximum value allowed for relative radius change in one time step.

**Default**: 0.01

**Parameters**

- **max_rel_radius_change** – The maximum relative radius change in one time step [-]

**set_max_rel_solute_composition_change** (max_rel_solute_composition_change: float = 0.01)
Set a limit on the time step by controlling solute depletion or saturation, especially at isothermal stage.

**Default**: 0.01
Parameters `max_rel_solute_composition_change` – The limit for the relative solute composition change [-]

```python
set_max_time_step(max_time_step: float = 0.1)
```

The maximum time step allowed for time integration as fraction of the simulation time. **Default:** 0.1

Parameters `max_time_step` – The maximum time step as fraction of the simulation time [-]

```python
set_max_time_step_during_heating(max_time_step_during_heating: float = 1.0)
```

The upper limit of the time step that has been enforced in the heating stages. **Default:** 1.0 s

Parameters `max_time_step_during_heating` – The maximum time step during heating [s]

```python
set_max_volume_fraction_dissolve_time_step(max_volume_fraction_dissolve_time_step: float = 0.01)
```

Sets the maximum volume fraction of subcritical particles allowed to dissolve in one time step. **Default:** 0.01

Parameters `max_volume_fraction_dissolve_time_step` – The maximum volume fraction of subcritical particles allowed to dissolve in one time step [-]

```python
set_min_radius_nucleus_as_particle(min_radius_nucleus_as_particle: float = 5e-10)
```

The cut-off lower limit of precipitate radius. **Default:** 5.0E-10 m

Parameters `min_radius_nucleus_as_particle` – The minimum radius of a nucleus to be considered as a particle [m]

```python
set_min_radius_points_per_magnitude(min_radius_points_per_magnitude: float = 100.0)
```

Sets the minimum number of grid points over one order of magnitude in radius. **Default:** 100.0

Parameters `min_radius_points_per_magnitude` – The minimum number of grid points over one order of magnitude in radius [-]

```python
set_radius_points_per_magnitude(radius_points_per_magnitude: float = 150.0)
```

Sets the number of grid points over one order of magnitude in radius. **Default:** 150.0

Parameters `radius_points_per_magnitude` – The number of grid points over one order of magnitude in radius [-]

```python
set_rel_radius_change_class_collision(rel_radius_change_class_collision: float = 0.5)
```

Sets the relative radius change for avoiding class collision. **Default:** 0.5

Parameters `rel_radius_change_class_collision` – The relative radius change for avoiding class collision [-]

class tc_python.precipitation.ParticleSizeDistribution
    Bases: object

Represents the state of a microstructure evolution at a certain time including its particle size distribution, composition and overall phase fraction.

```python
add_radius_and_number_density(radius: float, number_density: float)
```

Adds a radius and number density pair to the particle size distribution.

Parameters

- radius – The radius [m]
- number_density – The number of particles per unit volume per unit length [m^-4]

Returns This `ParticleSizeDistribution` object

```python
set_initial_composition(element_name: str, composition_value: float)
```

Sets the initial precipitate composition.
Parameters

• `element_name` – The name of the element
• `composition_value` – The composition value [composition unit defined for the calculation]

Returns This `ParticleSizeDistribution` object

Sets the type of the phase fraction or percentage. **Default**: By default volume fraction is used.

Parameters `volume_fraction_of_phase_type_enum` – Specifies if volume percent or fraction is used

Returns This `ParticleSizeDistribution` object

`set_volume_fraction_of_phase_value(value: float)`
Sets the overall volume fraction of the phase (unit based on the setting of `set_volume_fraction_of_phase_type()`).

Parameters `value` – The volume fraction 0.0 - 1.0 or percent value 0 - 100

Returns This `ParticleSizeDistribution` object

```python
class tc_python.precipitation.PrecipitateElasticProperties
Bases: object
```

Represents the elastic transformation strain of a certain precipitate class.

**Note:** This class is only relevant if the option `TransformationStrainCalculationOption.USER_DEFINED` has been chosen using `PrecipitatePhase.set_transformation_strain_calculation_option()`. The elastic strain can only be considered for non-spherical precipitates.

```python
set_e11(e11: float)
Sets the elastic strain tensor component e11. **Default**: 0.0

Parameters `e11` – The elastic strain tensor component e11

Returns This `PrecipitateElasticProperties` object

set_e12(e12: float)
Sets the strain tensor component e12. **Default**: 0.0

Parameters `e12` – The elastic strain tensor component e12

Returns This `PrecipitateElasticProperties` object

set_e13(e13: float)
Sets the elastic strain tensor component e13. **Default**: 0.0

Parameters `e13` – The elastic strain tensor component e13

Returns This `PrecipitateElasticProperties` object

set_e22(e22: float)
Sets the elastic strain tensor component e22. **Default**: 0.0

Parameters `e22` – The elastic strain tensor component e22

Returns This `PrecipitateElasticProperties` object

set_e23(e23: float)
Sets the elastic strain tensor component e23. **Default**: 0.0
Parameters `e23` – The elastic strain tensor component e23

Returns This `PrecipitateElasticProperties` object

```python
set_e33(e33: float)
```
Sets the elastic strain tensor component e33. Default: 0.0

Parameters `e33` – The elastic strain tensor component e33

Returns This `PrecipitateElasticProperties` object

class tc_python.precipitation.PrecipitateMorphology(value)

Bases: `enum.Enum`

Available precipitate morphologies.

- **CUBOID = 3**
  - Cuboidal precipitates, only available for bulk nucleation.

- **NEEDLE = 1**
  - Needle-like precipitates, only available for bulk nucleation.

- **PLATE = 2**
  - Plate-like precipitates, only available for bulk nucleation.

- **SPHERE = 0**
  - Spherical precipitates, this is the default morphology.

```python
class tc_python.precipitation.PrecipitatePhase(precipitate_phase_name: str)
```

Bases: `object`

Represents a certain precipitate class (i.e. a group of precipitates with the same phase and settings).

```python
disable_calculate_aspect_ratio_from_elastic_energy()
```
Disables the automatic calculation of the aspect ratio from the elastic energy of the phase.

Returns This `PrecipitatePhase` object

Note: If you use this method, you are required to set the aspect ratio explicitly using the method `set_aspect_ratio_value()`.

**Default:** This is the default setting (with an aspect ratio of 1.0).

```python
disable_driving_force_approximation()
```
Disables driving force approximation for this precipitate class. **Default:** Driving force approximation is disabled.

Returns This `PrecipitatePhase` object

```python
enable_calculate_aspect_ratio_from_elastic_energy()
```
Enables the automatic calculation of the aspect ratio from the elastic energy of the phase. **Default:** The aspect ratio is set to a value of 1.0.

Returns This `PrecipitatePhase` object

```python
enable_driving_force_approximation()
```
Enables driving force approximation for this precipitate class. This approximation is often required when simulating precipitation of multiple particles that use the same phase description. E.g. simultaneous precipitation of a Metal-Carbide(MC) and Metal-Nitride(MN) if configured as different composition sets of the same phase FCC_A1. **Default:** Driving force approximation is disabled.

Returns This `PrecipitatePhase` object

5.1. Calculations
**Tip:** Use this if simulations with several compositions sets of the same phase cause problems.

**set_alias** *(alias: str)*  
Sets an alias string that can later be used to get values from a calculated result. Typically used when having the same phase for several precipitates, but with different nucleation sites. For example two precipitates of the phase M7C3 with nucleation sites in ‘Bulk’ and at ‘Dislocations’. The alias can be used instead of the phase name when retrieving simulated results.  

**Parameters** alias – The alias string for this class of precipitates  
**Returns** This *PrecipitatePhase* object  

**Note:** Typically used when having using the same precipitate phase, but with different settings in the same calculation.

**set_aspect_ratio_value** *(aspect_ratio_value: float)*  
Sets the aspect ratio of the phase. **Default:** An aspect ratio of 1.0.  

**Parameters** aspect_ratio_value – The aspect ratio value  
**Returns** This *PrecipitatePhase* object  

**Note:** Only relevant if *disable_calculate_aspect_ratio_from_elastic_energy()* is used (which is the default).

**set_gibbs_energy_addition** *(gibbs_energy_addition: float)*  
Sets a Gibbs energy addition to the Gibbs energy of the phase. **Default:** 0.0 J/mol  

**Parameters** gibbs_energy_addition – The Gibbs energy addition [J/mol]  
**Returns** This *PrecipitatePhase* object  

**set_interfacial_energy** *(interfacial_energy: float)*  
Sets the interfacial energy. **Default:** If the interfacial energy is not set, it is automatically calculated using a broken-bond model.  

**Parameters** interfacial_energy – The interfacial energy [J/m^2]  
**Returns** This *PrecipitatePhase* object  

**Note:** The calculation of the interfacial energy using a broken-bond model is based on the assumption of an interface between a bcc- and a fcc-crystal structure with (110) and (111) lattice planes regardless of the actual phases.

**set_interfacial_energy_estimation_prefactor** *(interfacial_energy_estimation_prefactor: float)*  
Sets the interfacial energy prefactor. **Default:** Prefactor of 1.0 (only relevant if the interfacial energy is automatically calculated).  

**Parameters** interfacial_energy_estimation_prefactor – The prefactor for the calculated interfacial energy  
**Returns** This *PrecipitatePhase* object
Note: The interfacial energy prefactor is an amplification factor for the automatically calculated interfacial energy. Example: $\text{interfacial\_energy\_estimation\_prefactor} = 2.5 \Rightarrow 2.5 \times \text{calculated interfacial energy}$

```python
set_molar_volume(volume: float)
```
Sets the molar volume of the precipitate phase. Default: The molar volume obtained from the database. If no molar volume information is present in the database, a value of 7.0e-6 m$^3$/mol is used.

Parameters

- **volume** – The molar volume [m$^3$/mol]

Returns

This `PrecipitatePhase` object

```python
set_nucleation_at_dislocations(number_density: -1)
```
Activates nucleation at dislocations for this class of precipitates. Calling the method overrides any other nucleation setting for this class of precipitates. Default: If not set, by default bulk nucleation is chosen.

Parameters

- **number_density** – Number density of nucleation sites. If not set, the value is calculated based on the matrix settings (grain size, dislocation density) [m$^{-3}$].

Returns

This `PrecipitatePhase` object

```python
set_nucleation_at_grain_boundaries(wetting_angle: float = 90.0, number_density: float = -1)
```
Activates nucleation at grain boundaries for this class of precipitates. Calling the method overrides any other nucleation setting for this class of precipitates. Default: If not set, by default bulk nucleation is chosen.

Parameters

- **wetting_angle** – If not set, a default value of 90 degrees is used
- **number_density** – Number density of nucleation sites. If not set, the value is calculated based on the matrix settings (grain size) [m$^{-3}$].

Returns

This `PrecipitatePhase` object

```python
set_nucleation_at_grain_corners(wetting_angle: float = 90, number_density: float = -1)
```
Activates nucleation at grain corners for this class of precipitates. Calling the method overrides any other nucleation setting for this class of precipitates. Default: If not set, by default bulk nucleation is chosen.

Parameters

- **wetting_angle** – If not set, a default value of 90 degrees is used
- **number_density** – Number density of nucleation sites. If not set, the value is calculated based on the matrix settings (grain size) [m$^{-3}$].

Returns

This `PrecipitatePhase` object

```python
set_nucleation_at_grain_edges(wetting_angle: 90, number_density: -1)
```
Activates nucleation at the grain edges for this class of precipitates. Calling the method overrides any other nucleation setting for this class of precipitates. Default: If not set, by default bulk nucleation is chosen.

Parameters

- **wetting_angle** – If not set, a default value of 90 degrees is used
- **number_density** – Number density of nucleation sites. If not set, the value is calculated based on the matrix settings (grain size) [m$^{-3}$].

Returns

This `PrecipitatePhase` object

```python
set_nucleation_in_bulk(number_density: float = -1.0)
```
Activates nucleation in the bulk for this class of precipitates. Calling the method overrides any other
nucleation setting for this class of precipitates. **Default**: This is the default setting (with an automatically calculated number density).

**Parameters**

- **number_density** – Number density of nucleation sites. If not set, the value is calculated based on the matrix settings (molar volume) [m^-3]

**Returns**

This PrecipitatePhase object

**set_phase_boundary_mobility**(phase_boundary_mobility: float)

Sets the phase boundary mobility. **Default**: 10.0 m^4/(Js).

**Parameters**

- **phase_boundary_mobility** – The phase boundary mobility [m^4/(Js)]

**Returns**

This PrecipitatePhase object

**set_precipitate_morphology**(precipitate_morphology_enum: tc_python.precipitation.PrecipitateMorphology)

Sets the precipitate morphology. **Default**: PrecipitateMorphology.SPHERE

**Parameters**

- **precipitate_morphology_enum** – The precipitate morphology

**Returns**

This Precipitate_morphology object

**set_transformation_strain_calculation_option**(transformation_strain_calculation_option_enum: tc_python.precipitation.TransformationStrainCalculationOption)

Sets the transformation strain calculation option. **Default**: TransformationStrainCalculationOption.DISREGARD.

**Parameters**

- **transformation_strain_calculation_option_enum** – The chosen option

**Returns**

This PrecipitatePhase object

**with_elastic_properties**(elastic_properties: tc_python.precipitation.PrecipitateElasticProperties)

Sets the elastic properties. **Default**: The elastic transformation strain is disregarded by default.

**Parameters**

- **elastic_properties** – The elastic properties object

**Returns**

This PrecipitatePhase object

**Note:** This method has only an effect if the option TransformationStrainCalculationOption.USER_DEFINED is chosen using the method set_transformation_strain_calculation_option().

**with_growth_rate_model**(growth_rate_model_enum: tc_python.precipitation.GrowthRateModel)

Sets the growth rate model for the class of precipitates. **Default**: GrowthRateModel.SIMPLIFIED

**Parameters**

- **growth_rate_model_enum** – The growth rate model

**Returns**

This PrecipitatePhase object

**with_particle_size_distribution**(particle_size_distribution: tc_python.precipitation.ParticleSizeDistribution)

Sets the initial particle size distribution for this class of precipitates. **Default**: If the initial particle size distribution is not explicitly provided, the simulation will start from a supersaturated matrix.

**Parameters**

- **particle_size_distribution** – The initial particle size distribution object

**Returns**

This PrecipitatePhase object

**Tip:** Use this option if you want to study the further evolution of an existing microstructure.
class tc_python.precipitation.PrecipitationCCTCalculation(calculation)
Bases: tc_python.abstract_base.AbstractCalculation

Configuration for a Continuous-Cooling-Time (CCT) precipitation calculation.

calculate() \rightarrow tc_python.precipitation.PrecipitationCalculationTTorCCTResult
Runs the CCT diagram calculation.

Returns A PrecipitationCalculationTTorCCTResult which later can be used to
get specific values from the calculated result

get_system_data() \rightarrow tc_python.abstract_base.SystemData
Retruns the content of the database for the currently loaded system. This can be used to modify the param-
eters and functions and to change the current system by using with_system_modifications().

Note: Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb-file.

Returns The system data

set_composition(element_name: str, value: float)
Sets the composition of the elements. The unit for the composition can be changed using
set_composition_unit(). Default: Mole percent (CompositionUnit.MOLE_PERCENT)

Parameters
  • element_name – The element
  • value – The composition (fraction or percent depending on the composition unit)

Returns This PrecipitationCCTCalculation object

set_composition_unit(unit_enum: tc_python.utils.CompositionUnit)
Sets the composition unit. Default: Mole percent (CompositionUnit.MOLE_PERCENT).

Parameters unit_enum – The new composition unit

Returns This PrecipitationCCTCalculation object

set_cooling_rates(cooling_rates: List[float])
Sets all cooling rates for which the CCT diagram should be calculated.

Parameters cooling_rates – A list of cooling rates [K/s]

Returns This PrecipitationCCTCalculation object

set_max_temperature(max_temperature: float)
Sets maximum temperature of the CCT diagram.

Parameters max_temperature – the maximum temperature [K]

Returns This PrecipitationCCTCalculation object

set_min_temperature(min_temperature: float)
Sets the minimum temperature of the CCT diagram.

Parameters min_temperature – the minimum temperature [K]

Returns This PrecipitationCCTCalculation object

stop_at_volume_fraction_of_phase(stop_criterion_value: float)
Sets the stop criterion as a volume fraction of the phase. This setting is applied to all phases.
Parameters **stop_criterion_value** – the volume fraction of the phase (a value between 0 and 1)

Returns This PrecipitationCCTCalculation object

**with_matrix_phase** *(matrix_phase: tc_python.precipitation.MatrixPhase)*

Sets the matrix phase.

Parameters **matrix_phase** – The matrix phase

Returns This PrecipitationCCTCalculation object

**with_numerical_parameters** *(numerical_parameters: tc_python.precipitation.NumericalParameters)*

Sets the numerical parameters. If not specified, reasonable defaults are be used.

Parameters **numerical_parameters** – The parameters

Returns This PrecipitationCCTCalculation object

**with_system_modifications** *(system_modifications: tc_python.abstract_base.SystemModifications)*

Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

Note: This is only possible if the system has been read from unencrypted (i.e. user) databases loaded as a * .tdb*-file.

Parameters **system_modifications** – The system modification to be performed

Returns This PrecipitationCCTCalculation object

class tc_python.precipitation.PrecipitationCalculationResult *(result)*

Bases: tc_python.abstract_base.AbstractResult

Result of a precipitation calculation. This can be used to query for specific values.

**save_to_disk** *(path: str)*

Saves the result to disc. Note that a result is a folder, containing potentially many files. The result can later be loaded with load_result_from_disk()

Parameters **path** – the path to the folder you want the result to be saved in. It can be relative or absolute.

Returns this PrecipitationCalculationResult object

class tc_python.precipitation.PrecipitationCalculationSingleResult *(result)*

Bases: tc_python.precipitation.PrecipitationCalculationCalculationResult

Result of a isothermal or non-isothermal precipitation calculation. This can be used to query for specific values.

Search the Thermo-Calc help for definitions of the axis variables, e.g. search isothermal variables or non-isothermal variables.

**get_aspect_ratio_distribution_for_particle_length_of** *(precipitate_id: str, time: float) → [typing.List[float], typing.List[float]]*

Returns the aspect ratio distribution of a precipitate in dependency of its mean particle length at a certain time.

Only available if the morphology is set to PrecipitateMorphology.NEEDLE or PrecipitateMorphology.PLATE.
Parameters

- `time` – The time [s]
- `precipitate_id` – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (mean particle length [m], aspect ratio)

`get_aspect_ratio_distribution_for_radius_of` (`precipitate_id: str, time: float`) → [typing.List[float], typing.List[float]]

Returns the aspect ratio distribution of a precipitate in dependency of its mean radius at a certain time.

Only available if the morphology is set to `PrecipitateMorphology.NEEDLE` or `PrecipitateMorphology.PLATE`.

Parameters

- `time` – The time [s]
- `precipitate_id` – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (mean radius [m], aspect ratio)

`get_critical_radius_of` (`precipitate_id: str`) → [typing.List[float], typing.List[float]]

Returns the critical radius of a precipitate in dependency of the time.

Parameters `precipitate_id` – The id of a precipitate can either be phase name or alias

Returns A tuple of two lists of floats (time [s], critical radius [m])

`get_cubic_factor_distribution_for_particle_length_of` (`precipitate_id: str, time: float`) → [typing.List[float], typing.List[float]]

Returns the cubic factor distribution of a precipitate in dependency of its mean particle length at a certain time.

Only available if the morphology is set to `PrecipitateMorphology.CUBOID`.

Parameters

- `time` – The time in seconds
- `precipitate_id` – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (particle length [m], cubic factor)

`get_cubic_factor_distribution_for_radius_of` (`precipitate_id: str, time: float`) → [typing.List[float], typing.List[float]]

Returns the cubic factor distribution of a precipitate in dependency of its mean radius at a certain time.

Only available if the morphology is set to `PrecipitateMorphology.CUBOID`.

Parameters

- `time` – The time [s]
- `precipitate_id` – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (radius [m], cubic factor)

`get_driving_force_of` (`precipitate_id: str`) → [typing.List[float], typing.List[float]]

Returns the (by R * T) normalized driving force of a precipitate in dependency of the time.

Parameters `precipitate_id` – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (time [s], normalized driving force)
get_grain_critical_radius () \rightarrow \text{[typing.List[float], typing.List[float]]}
Returns the critical radius of grains in dependency of the time.

**Returns** A tuple of two lists of floats (time [s], critical radius [m])

get_grain_mean_radius () \rightarrow \text{[typing.List[float], typing.List[float]]}
Returns the mean grain size of the matrix phase in dependency of the time.

**Returns** A tuple of two lists of floats (time [s], mean radius [m])

get_grain_number_density () \rightarrow \text{[typing.List[float], typing.List[float]]}
Returns the grain number density in dependency of the time.

**Returns** A tuple of two lists of floats (time [s], grain number density [m^-3])

get_grain_number_density_distribution_for_length (time: float) \rightarrow \text{[typing.List[float], typing.List[float]]}
Returns the number density distribution of grains in dependency of its mean particle length at a certain time.

**Parameters**
- **time** – The time [s]

**Returns** A tuple of two lists of floats (grain length[m], number of grains per unit volume per unit length [m^-4])

get_grain_number_density_distribution_for_radius (time: float) \rightarrow \text{[typing.List[float], typing.List[float]]}
Returns the number density distribution of a grains in dependency of its mean radius at a certain time.

**Parameters**
- **time** – The time [s]

**Returns** A tuple of two lists of floats (radius [m], number of grains per unit volume per unit length [m^-4])

get_grain_size_distribution (time: float) \rightarrow \text{[typing.List[float], typing.List[float]]}
Returns the size distribution of the matrix phase in dependency of its grain radius length at a certain time.

**Parameters**
- **time** – The time [s]

**Returns** A tuple of two lists of floats (grain radius[m], number density of grains[m^-3])

get_matrix_composition_in_mole_fraction_of (element_name: str) \rightarrow \text{[typing.List[float], typing.List[float]]}
Returns the matrix composition (as mole fractions) of a certain element in dependency of the time.

**Parameters**
- **element_name** – The element

**Returns** A tuple of two lists of floats (time [s], mole fraction)

get_matrix_composition_in_weight_fraction_of (element_name: str) \rightarrow \text{[typing.List[float], typing.List[float]]}
Returns the matrix composition (as weight fraction) of a certain element in dependency of the time.

**Parameters**
- **element_name** – The element

**Returns** A tuple of two lists of floats (time [s], weight fraction)

g_mean_aspect_ratio_of (precipitate_id: str) \rightarrow \text{[typing.List[float], typing.List[float]]}
Returns the mean aspect ratio of a precipitate in dependency of the time.

Only available if the morphology is set to PrecipitateMorphology.NEEDLE or PrecipitateMorphology.PLATE.
Parameters **precipitate_id** – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (time [s], mean aspect ratio)

get_mean_cubic_factor_of (**precipitate_id**: str) → [typing.List[float], typing.List[float]]

Returns the mean cubic factor of a precipitate in dependency of the time. Only available if the morphology is set to PrecipitateMorphology.CUBOID.

Parameters **precipitate_id** – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (time [s], mean cubic factor)

get_mean_particle_length_of (**precipitate_id**: str) → [typing.List[float], typing.List[float]]

Returns the mean particle length of a precipitate in dependency of the time.

Only available if the morphology is set to PrecipitateMorphology.NEEDLE or PrecipitateMorphology.PLATE.

Parameters **precipitate_id** – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (time [s], mean particle length [m])

get_mean_radius_of (**precipitate_id**: str) → [typing.List[float], typing.List[float]]

Returns the mean radius of a precipitate in dependency of the time.

Parameters **precipitate_id** – The id of a precipitate can either be phase name or alias

Returns A tuple of two lists of floats (time [s], mean radius [m])

get_normalized_grain_size_distribution (**time**: float) → [typing.List[float], typing.List[float]]

Returns the normalized number density distribution of a grains at a certain time.

Parameters **time** – The time [s]

Returns A tuple of two lists of floats (Normalized size, Frequency)

get_normalized_number_density_distribution_of (**precipitate_id**: str, **time**: float) → [typing.List[float], typing.List[float]]

Returns the normalized number density distribution of a precipitate at a certain time.

Parameters

- **time** – The time [s]
- **precipitate_id** – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (Normalized size, Frequency)

get_nucleation_rate_of (**precipitate_id**: str) → [typing.List[float], typing.List[float]]

Returns the nucleation rate of a precipitate in dependency of the time.

Parameters **precipitate_id** – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (time [s], nucleation rate [m^-3 s^-1])

get_number_density_distribution_for_particle_length_of (**precipitate_id**: str, **time**: float) → [typing.List[float], typing.List[float]]

Returns the number density distribution of a precipitate in dependency of its mean particle length at a
certain time.

Parameters

- `time` – The time [s]
- `precipitate_id` – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (particle length [m], number of particles per unit volume per unit length [m^-4])

get_number_density_distribution_for_radius_of(precipitate_id: str, time: float) -> [typing.List[float], typing.List[float]]

Returns the number density distribution of a precipitate in dependency of its mean radius at a certain time.

Parameters

- `time` – The time [s]
- `precipitate_id` – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (radius [m], number of particles per unit volume per unit length [m^-4])

get_number_density_of(precipitate_id: str) -> [typing.List[float], typing.List[float]]

Returns the particle number density of a precipitate in dependency of the time.

Parameters `precipitate_id` – The id of a precipitate can either be phase name or alias

Returns A tuple of two lists of floats (time [s], particle number density [m^-3])

get_precipitate_composition_in_mole_fraction_of(precipitate_id: str, element_name: str) -> [typing.List[float], typing.List[float]]

Returns the precipitate composition (as mole fractions) of a certain element in dependency of the time.

Parameters

- `precipitate_id` – The id of a precipitate can either be phase name or alias
- `element_name` – The element

Returns A tuple of two lists of floats (time [s], mole fraction)

get_precipitate_composition_in_weight_fraction_of(precipitate_id: str, element_name: str) -> [typing.List[float], typing.List[float]]

Returns the precipitate composition (as weight fraction) of a certain element in dependency of the time.

Parameters

- `precipitate_id` – The id of a precipitate can either be phase name or alias
- `element_name` – The element

Returns A tuple of two lists of floats (time [s], weight fraction)

get_size_distribution_for_particle_length_of(precipitate_id: str, time: float) -> [typing.List[float], typing.List[float]]

Returns the size distribution of a precipitate in dependency of its mean particle length at a certain time.

Parameters

- `time` – The time [s]
- `precipitate_id` – The id of a precipitate can either be the phase name or an alias
Returns A tuple of two lists of floats (particle length [m], number of particles per unit volume per unit length [m^-4])

`get_size_distribution_for_radius_of (precipitate_id: str, time: float) → [typing.List[float], typing.List[float]]`

Returns the size distribution of a precipitate in dependency of its mean radius at a certain time.

Parameters

- **time** – The time [s]
- **precipitate_id** – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (radius [m], number of particles per unit volume per unit length [m^-4])

`get_volume_fraction_of (precipitate_id: str) → [typing.List[float], typing.List[float]]`

Returns the volume fraction of a precipitate in dependency of the time.

Parameters **precipitate_id** – The id of a precipitate can either be the phase name or an alias

Returns A tuple of two lists of floats (time [s], volume fraction)

**5.1. Calculations**

```
class tc_python.precipitation.PrecipitationCalculationTTToCCTResult (result)
    Bases: tc_python.precipitation.PrecipitationCalculationResult
    Result of a TTT or CCT precipitation calculation.

get_result_for_precipitate (precipitate_id: str) → [typing.List[float], typing.List[float]]
    Returns the calculated data of a TTT or CCT diagram for a certain precipitate.

    Parameters **precipitate_id** – The id of a precipitate can either be the phase name or an alias

    Returns A tuple of two lists of floats (time [s], temp [K])
```

```
class tc_python.precipitation.PrecipitationIsoThermalCalculation (calculation)
    Bases: tc_python.abstract_base.AbstractCalculation
    Configuration for an isothermal precipitation calculation.

calculate () → tc_python.precipitation.PrecipitationCalculationSingleResult
    Runs the isothermal precipitation calculation.

    Returns A PrecipitationCalculationSingleResult which later can be used to get specific values from the calculated result

get_system_data () → tc_python.abstract_base.SystemData
    Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using with_system_modifications().
```

Note: Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb-file.

Returns The system data

```
set_composition (element_name: str, value: float)
    Sets the composition of the elements. The unit for the composition can be changed using set_composition_unit(). Default: Mole percent (CompositionUnit.MOLE_PERCENT)

    Parameters

    - **element_name** – The element
```
- **value** – The composition (fraction or percent depending on the composition unit)

  **Returns** This `PrecipitationIsoThermalCalculation` object

  **set_composition_unit** *(unit_enum: tc_python.utils.CompositionUnit = <CompositionUnit.MOLE_PERCENT: 1>)*

  Sets the composition unit. **Default**: Mole percent (CompositionUnit.MOLE_PERCENT).

  **Parameters** unit_enum – The new composition unit

  **Returns** This `PrecipitationIsoThermalCalculation` object

  **set_simulation_time** *(simulation_time: float)*

  Sets the simulation time.

  **Parameters** simulation_time – The simulation time [s]

  **Returns** This `PrecipitationIsoThermalCalculation` object

  **set_temperature** *(temperature: float)*

  Sets the temperature for the isothermal simulation.

  **Parameters** temperature – The temperature [K]

  **Returns** This `PrecipitationIsoThermalCalculation` object

  **with_matrix_phase** *(matrix_phase: tc_python.precipitation.MatrixPhase)*

  Sets the matrix phase.

  **Parameters** matrix_phase – The matrix phase

  **Returns** This `PrecipitationIsoThermalCalculation` object

  **with_numerical_parameters** *(numerical_parameters: tc_python.precipitation.NumericalParameters)*

  Sets the numerical parameters. If not specified, reasonable defaults are be used.

  **Parameters** numerical_parameters – The parameters

  **Returns** This `PrecipitationIsoThermalCalculation` object

  **with_system_modifications** *(system_modifications: tc_python.abstract_base.SystemModifications)*

  Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

  **Note**: This is only possible if the system has been read from unencrypted (i.e. user) databases loaded as a *.tdb*-file.

  **Parameters** system_modifications – The system modification to be performed

  **Returns** This `PrecipitationIsoThermalCalculation` object

**class** `tc_python.precipitation.PrecipitationNonIsoThermalCalculation(calculation)`

**Bases**: `tc_python.abstract_base.AbstractCalculation`

Configuration for a non-isothermal precipitation calculation.

**calculate** () → `tc_python.precipitation.PrecipitationCalculationSingleResult`

Runs the non-isothermal precipitation calculation.

  **Returns** A `PrecipitationCalculationSingleResult` which later can be used to get specific values from the calculated result
**get_system_data() → tc_python.abstract_base.SystemData**

Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using `with_system_modifications()`.

**Note:** Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb*-file.

**Returns** The system data

**set_composition(element_name: str, value: float)**

Sets the composition of the elements. The unit for the composition can be changed using `set_composition_unit()`. Default: Mole percent (CompositionUnit.MOLE_PERCENT)

**Parameters**
- **element_name** – The element
- **value** – The composition (fraction or percent depending on the composition unit)

**Returns** This `PrecipitationIsoThermalCalculation` object

**set_composition_unit(unit_enum: tc_python.utils.CompositionUnit)**

Sets the composition unit. Default: Mole percent (CompositionUnit.MOLE_PERCENT).

**Parameters** unit_enum – The new composition unit

**Returns** This `PrecipitationIsoThermalCalculation` object

**set_simulation_time(simulation_time: float)**

Sets the simulation time.

**Parameters** simulation_time – The simulation time [s]

**Returns** This `PrecipitationNonThermalCalculation` object

**with_matrix_phase(matrix_phase: tc_python.precipitation.MatrixPhase)**

Sets the matrix phase.

**Parameters** matrix_phase – The matrix phase

**Returns** This `PrecipitationIsoThermalCalculation` object

**with_numerical_parameters(numerical_parameters: tc_python.precipitation.NumericalParameters)**

Sets the numerical parameters. If not specified, reasonable defaults are be used.

**Parameters** numerical_parameters – The parameters

**Returns** This `PrecipitationIsoThermalCalculation` object

**with_system_modifications(system_modifications: tc_python.abstract_base.SystemModifications)**

Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

**Note:** This is only possible if the system has been read from unencrypted (i.e. user) databases loaded as a *.tdb*-file.

**Parameters** system_modifications – The system modification to be performed

**Returns** This `PrecipitationNonThermalCalculation` object
with_temperature_profile(temperature_profile: tc_python.utils.TemperatureProfile)
Sets the temperature profile to use with this calculation.

**Parameters**

- **temperature_profile** – the temperature profile object (specifying time / temperature points)

**Returns**

This PrecipitationNonThermalCalculation object

class tc_python.precipitation.PrecipitationTTTCalculation(calculation)
Bases: tc_python.abstract_base.AbstractCalculation

Configuration for a TTT (Time-Temperature-Transformation) precipitation calculation.

calculate() \rightarrow tc_python.precipitation.PrecipitationCalculationTTToCCTResult
Runs the TTT diagram calculation.

**Returns**

A PrecipitationCalculationTTToCCTResult which later can be used to get specific values from the calculated result.

get_system_data() \rightarrow tc_python.abstract_base.SystemData

Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using with_system_modifications().

**Note:** Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb*-file.

**Returns**

The system data

set_composition(element_name: str, value: float)
Sets the composition of the elements. The unit for the composition can be changed using set_composition_unit(). Default: Mole percent (CompositionUnit.MOLE_PERCENT)

**Parameters**

- **element_name** – The element
- **value** – The composition (fraction or percent depending on the composition unit)

**Returns**

This PrecipitationTTTCalculation object

set_composition_unit(unit_enum: tc_python.utils.CompositionUnit)
Sets the composition unit. Default: Mole percent (CompositionUnit.MOLE_PERCENT).

**Parameters**

- **unit_enum** – The new composition unit

**Returns**

This PrecipitationTTTCalculation object

set_max_annealing_time(max_annealing_time: float)
Sets the maximum annealing time, i.e. the maximum time of the simulation if the stopping criterion is not reached.

**Parameters**

- **max_annealing_time** – the maximum annealing time [s]

**Returns**

This PrecipitationTTTCalculation object

set_max_temperature(max_temperature: float)
Sets the maximum temperature for the TTT diagram.

**Parameters**

- **max_temperature** – the maximum temperature [K]

**Returns**

This PrecipitationTTTCalculation object

set_min_temperature(min_temperature: float)
Sets the minimum temperature for the TTT diagram.
Parameters **min_temperature** – the minimum temperature [K]

Returns This *PrecipitationTTTCalculation* object

**set_temperature_step** *(temperature_step: float)*
Sets the temperature step for the TTT diagram. If not set, the default value is 10 K.

Parameters **temperature_step** – the temperature step [K]

Returns This *PrecipitationTTTCalculation* object

**stop_at_percent_of_equilibrium_fraction** *(percentage: float)*
Sets the stop criterion to a percentage of the overall equilibrium phase fraction, alternatively a required volume fraction can be specified (using *stop_at_volume_fraction_of_phase()*).

Parameters **percentage** – the percentage to stop at (value between 0 and 100)

Returns This *PrecipitationTTTCalculation* object

**stop_at_volume_fraction_of_phase** *(volume_fraction: float)*
Sets the stop criterion as a volume fraction of the phase, alternatively a required percentage of the equilibrium phase fraction can be specified (using *stop_at_percent_of_equilibria_fraction()*).

This setting is applied to all phases.

Parameters **volume_fraction** – the volume fraction to stop at (a value between 0 and 1)

Returns This *PrecipitationTTTCalculation* object

**with_matrix_phase** *(matrix_phase: tc_python.precipitation.MatrixPhase)*
Sets the matrix phase.

Parameters **matrix_phase** – The matrix phase

Returns This *PrecipitationTTTCalculation* object

**with_numerical_parameters** *(numerical_parameters: tc_python.precipitation.NumericalParameters)*
Sets the numerical parameters. If not specified, reasonable defaults are be used.

Parameters **numerical_parameters** – The parameters

Returns This *PrecipitationTTTCalculation* object

**with_system_modifications** *(system_modifications: tc_python.abstract_base.SystemModifications)*
Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

Parameters **system_modifications** – The system modification to be performed

Returns This *PrecipitationTTTCalculation* object

---

**Note:** This is only possible if the system has been read from unencrypted (i.e. *user*) databases loaded as a *.tdb*-file.

---

**class** *tc_python.precipitation.TransformationStrainCalculationOption*(value)*

Bases: enum.Enum

Options for calculating the transformation strain.

**CALCULATE_FROM_MOLAR_VOLUME = 2**
Calculates the transformation strain from the molar volume, obtains a purely dilatational strain.
DISREGARD = 1
Ignores the transformation strain, **this is the default setting**.

USER_DEFINED = 3
Transformation strain to be specified by the user.

class tc_python.precipitation.VolumeFractionOfPhaseType(value)
Bases: enum.Enum

Unit of the volume fraction of a phase.

VOLUME_FRACTION = 6
Volume fraction (0 - 1), **this is the default**.

VOLUME_PERCENT = 5
Volume percent (0% - 100%).

### 5.1.4 Module “scheil”

class tc_python.scheil.CalculateSecondaryDendriteArmSpacing
Bases: tc_python.scheil.ScheilBackDiffusion

Configures a secondary dendrite arm spacing calculation used by Scheil with back diffusion. The used equation is $c \times \text{cooling rate}^{-n}$ with $c$ and $n$ being provided either by the user or taken from the defaults.

**set_c** *(c: float = 5e-05)*
Sets the scaling factor $c$ in the governing equation $c \times \text{cooling rate}^{-n}$.

**Default**: 50 µm

**Parameters**
- **c** – The scaling factor [m]

**Returns**
This CalculateSecondaryDendriteArmSpacing object

**set_cooling_rate** *(cooling_rate: float = 1.0)*
Sets the cooling rate.

**Default**: 1.0 K/s

An increased value moves the result from equilibrium toward a Scheil-Gulliver calculation.

**Parameters**
- **cooling_rate** – The cooling rate [K/s]

**Returns**
This CalculateSecondaryDendriteArmSpacing object

**set_fast_diffusing_elements** *(element_names: List[str]*)
Sets elements as fast diffusing. This allows redistribution of these elements in both the solid and liquid parts of the alloy.

**Default**: No fast-diffusing elements.

**Parameters**
- **element_names** – The elements

**Returns**
This CalculateSecondaryDendriteArmSpacing object

**set_n** *(n: float = 0.33)*
Sets the exponent $n$ in the governing equation $c \times \text{cooling rate}^{-n}$.

**Default**: 0.33

**Parameters**
- **n** – The exponent [-]

**Returns**
This CalculateSecondaryDendriteArmSpacing object
```python
set_primary_phasename (primary_phase_name: str = 'AUTOMATIC')

Sets the name of the primary phase.

The primary phase is the phase where the back diffusion takes place. If AUTOMATIC is selected, the program tries to find the phase which will give the most back diffusion. That behavior can be overridden by selecting a specific primary phase.

Default: AUTOMATIC

Parameters primary_phase_name – The phase name (or AUTOMATIC)

Returns This CalculateSecondaryDendriteArmSpacing object
```

```python
class tc_python.scheil.ConstantSecondaryDendriteArmSpacing (secondary_dendrite_arm_spacing: float = 5e-05)

Bases: tc_python.scheil.ScheilBackDiffusion

Configures a constant secondary dendrite arm spacing used by Scheil with back diffusion. The secondary dendrite arm spacing can either be provided by the user or taken from the defaults.

set_cooling_rate (cooling_rate: float = 1.0)

Sets the cooling rate.

Default: 1.0 K/s

An increased value moves the result from equilibrium toward a Scheil-Gulliver calculation.

Parameters cooling_rate – The cooling rate [K/s]

Returns This ConstantSecondaryDendriteArmSpacing object
```

```python
set_fast_diffusing_elements (element_names: List[str])

Sets elements as fast diffusing. This allows redistribution of these elements in both the solid and liquid parts of the alloy.

Default: No fast-diffusing elements.

Parameters element_names – The elements

Returns This ConstantSecondaryDendriteArmSpacing object
```

```python
set_primary_phasename (primary_phase_name: str = 'AUTOMATIC')

Sets the name of the primary phase.

The primary phase is the phase where the back diffusion takes place. If AUTOMATIC is selected, the program tries to find the phase which will give the most back diffusion. That behavior can be overridden by selecting a specific primary phase.

Default: AUTOMATIC

Parameters primary_phase_name – The phase name (or AUTOMATIC)

Returns This ConstantSecondaryDendriteArmSpacing object
```

```python
class tc_python.scheil.ScheilBackDiffusion

Bases: tc_python.scheil.ScheilCalculationType

Configuration for back diffusion in the solid primary phase.
```

**Warning:** This feature has only effect on systems with diffusion data (typically a mobility database). If used for a system without diffusion data, a normal Scheil calculation is done.
**classmethod calculate_secondary_dendrite_arm_spacing()**

Calculate the secondary dendrite arm spacing based on the following equation: \( c \times \text{cooling\_rate}^{-n} \) with \( c \) and \( n \) being provided either by the user or taken from the defaults.

Use the methods provided by \texttt{CalculateSecondaryDendriteArmSpacing} to configure the parameters.

**Returns** A \texttt{CalculateSecondaryDendriteArmSpacing}

**classmethod constant_secondary_dendrite_arm_spacing(** *secondary_dendrite_arm_spacing*: float = 5e-05)**

Assuming constant secondary dendrite arm spacing, provided either by the user or taken from the defaults.

**Default**: 50 \( \mu \text{m} \)

**Parameters** \texttt{secondary\_dendrite\_arm\_spacing} – The dendrite arm spacing [m]

**Returns** A \texttt{ConstantSecondaryDendriteArmSpacing}

**class** \texttt{tc\_python.scheil.ScheilCalculation**(calculator)**}

\texttt{Bases: tc\_python.abstract\_base.AbstractCalculation}

Configuration for a Scheil solidification calculation.

**Note:** Specify the settings, the calculation is performed with \texttt{calculate()}.  

**calculate()** \( \rightarrow \texttt{tc\_python.scheil.ScheilCalculationResult} \)

Runs the Scheil calculation.

**Warning:** Scheil calculations do not support the GAS phase being selected, this means the GAS phase must always be deselected in the system if it is present in the database.

**Returns** A \texttt{ScheilCalculationResult} which later can be used to get specific values from the simulation.

**disable_global_minimization()**

Disables global minimization.

**Default:** Enabled

**Note:** When enabled, a global minimization test is performed when an equilibrium is reached. This costs more computer time but the calculations are more robust.

**Returns** This \texttt{ScheilCalculation} object

**enable_global_minimization()**

Enables global minimization.

**Default:** Enabled

**Note:** When enabled, a global minimization test is performed when an equilibrium is reached. This costs more computer time but the calculations are more robust.

**Returns** This \texttt{ScheilCalculation} object
get_system_data() → tc_python.abstract_base.SystemData
Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using with_system_modifications().

Note: Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb-file.

Returns The system data

set_composition(component_name: str, value: float)
Sets the composition of a component. The unit for the composition can be changed using set_composition_unit().

Default: Mole percent (CompositionUnit.MOLE_PERCENT)

Parameters
• component_name – The component
• value – The composition value [composition unit defined for the calculation]

Returns This ScheilCalculation object

set_composition_unit(unit_enum: tc_python.utils.CompositionUnit = <CompositionUnit.MOLE_PERCENT: 1>)
Sets the composition unit.

Default: Mole percent (CompositionUnit.MOLE_PERCENT).

Parameters unit_enum – The new composition unit

Returns This ScheilCalculation object

set_fast_diffusing_elements(element_names: List[str])
Sets elements as fast diffusing. This allows redistribution of these elements in both the solid and liquid parts of the alloy.

Note: Deprecated in version 2021b: This method has been moved into the class ScheilCalculationType, which is used with the method with_calculation_type(). It will be removed in release 2022b.

Default: No fast-diffusing elements.

Parameters element_names – The elements

Returns This ScheilCalculation object

set_start_temperature(temperature_in_kelvin: float = 2500.0)
Sets the start temperature.

Warning: The start temperature needs to be higher than the liquidus temperature of the alloy.

Default: 2500.0 K

Parameters temperature_in_kelvin – The temperature [K]

Returns This ScheilCalculation object
**with_back_diffusion** *(scheil_back_diffusion: tc_python.scheil.ScheilBackDiffusion)*

Enables back diffusion in the solid primary phase.

**Note:** Deprecated in version 2021b: Use **with_calculation_type()** instead. This method will be removed in release 2022b.

**Warning:** This feature has only effect on systems with diffusion data (typically a mobility database). If used for a system without diffusion data, a normal Scheil calculation is performed.

**Parameters** scheil_back_diffusion – an instance of a ScheilBackDiffusion class, where the options for back diffusion can be specified.

**Returns** This ScheilCalculation object

**with_calculation_type** *(scheil_calculation_type: tc_python.scheil.ScheilCalculationType)*

Chooses a specific Scheil calculation. ClassicScheil for only setting fast diffusers, ScheilBackDiffusion enables back diffusion in the solid primary phase and optionally fast diffusers in all solid phases, and ScheilSoluteTrapping enables solute trapping in the solid primary phase.

**Parameters** scheil_type: Type of Scheil calculation, either ScheilClassic, ScheilBackDiffusion or ScheilSoluteTrapping

**Returns** This ScheilCalculation object

**with_options** *(options: tc_python.scheil.ScheilOptions)*

Sets the Scheil simulation options.

**Parameters** options – The Scheil simulation options

**Returns** This ScheilCalculation object

**with_system_modifications** *(system_modifications: tc_python.abstract_base.SystemModifications)*

Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

**Note:** This is only possible if the system has been read from unencrypted (i.e. user) databases loaded as a *.tdb*-file.

**Parameters** system_modifications – The system modification to be performed

**Returns** This ScheilCalculation object

**class** tc_python.scheil.ScheilCalculationResult *(result)*

Bases: tc_python.abstract_base.AbstractResult

Result of a Scheil calculation.

**get_values_grouped_by_quantity_of** *(x_quantity: Union[tc_python.quantity_factory.ScheilQuantity, str], y_quantity: Union[tc_python.quantity_factory.ScheilQuantity, str], sort_and_merge: bool = True) → Dict[str, tc_python.utils.ResultSetValueGroup]*

Returns x-y-line data grouped by the multiple datasets of the specified quantities (for example in dependency of phases or components). Use get_values_of() instead if you need no separation. The available quantities can be found in the documentation of the factory class ScheilQuantity.
Note: The different datasets might contain NaN-values between different subsections and might not be sorted even if the flag ‘sort_and_merge’ has been set (because they might be unsortable due to their nature).

Parameters

- **x_quantity** – The first Scheil quantity (“x-axis”), Console Mode syntax strings can be used as an alternative (for example “T”)
- **y_quantity** – The second Scheil quantity (“y-axis”), Console Mode syntax strings can be used as an alternative (for example “NV”)
- **sort_and_merge** – If True, the data is sorted and merged into as few subsections as possible (divided by NaN)

Returns Containing the ResultValueGroup dataset objects with their quantity labels as keys

get_values_grouped_by_stable_phases_of(x_quantity: Union[tc_python.quantity_factory.ScheilQuantity, str], y_quantity: Union[tc_python.quantity_factory.ScheilQuantity, str], sort_and_merge: bool = True) → Dict[str, tc_python.utils.ResultValueGroup]

Returns x-y-line data grouped by the sets of “stable phases” (for example “LIQUID” or “LIQUID + FCC_A1”). Use get_values_of() instead if you need no separation. The available quantities can be found in the documentation of the factory class ScheilQuantity.

Note: The different datasets might contain NaN-values between different subsections and might not be sorted even if the flag ‘sort_and_merge’ has been set (because they might be un sortable due to their nature).

Parameters

- **x_quantity** – The first Scheil quantity (“x-axis”), Console Mode syntax strings can be used as an alternative (for example “T”)
- **y_quantity** – The second Scheil quantity (“y-axis”), Console Mode syntax strings can be used as an alternative (for example “NV”)
- **sort_and_merge** – If True, the data will be sorted and merged into as few subsections as possible (divided by NaN)

Returns Containing the ResultValueGroup dataset objects with their “stable phases” labels as keys

get_values_of(x_quantity: Union[tc_python.quantity_factory.ScheilQuantity, str], y_quantity: Union[tc_python.quantity_factory.ScheilQuantity, str]) → [typing.List[float], typing.List[float]]

Returns sorted x-y-line data without any separation. Use get_values_grouped_by_quantity_of() or get_values_grouped_by_stable_phases_of() instead if you need such a separation. The available quantities can be found in the documentation of the factory class ScheilQuantity.

Note: This method will always return sorted data without any NaN-values. In case of ambiguous quantities (for example: CompositionOfPhaseAsWeightFraction(“FCC_A1”, “All”)) that can give data that is...
hard to interpret. In such a case you need to choose the quantity in another way or use one of the other methods.

Parameters

- **x_quantity** – The first Scheil quantity (“x-axis”), Console Mode syntax strings can be used as an alternative (for example “T”)
- **y_quantity** – The second Scheil quantity (“y-axis”), Console Mode syntax strings can be used as an alternative (for example “NV”)

Returns A tuple containing the x- and y-data in lists

```python
save_to_disk(path: str)
```

Saves the result to disc. Note that a result is a folder, containing potentially many files. The result can later be loaded with `load_result_from_disk()`

Parameters

- **path** – the path to the folder you want the result to be saved in.

Returns this `ScheilCalculationResult` object

```python
class tc_python.scheil.ScheilCalculationType
    Bases: object
    Specific configuration for the different Scheil calculation types

    classmethod scheil_back_diffusion()
    Configuration for back diffusion in the solid primary phase.
```

**Warning:** This feature has only effect on systems with diffusion data (typically a mobility database). If used for a system without diffusion data, a normal Scheil calculation is done. :return: A `ScheilBackDiffusion`

```python
class tc_python.scheil.ScheilClassic
    Bases: tc_python.scheil.ScheilCalculationType
    Configuration for Classic Scheil with fast diffusers.
```

```python
classmethod scheil_classic()
    Configuration for Classic Scheil with fast diffusers. :return: A `ScheilClassic`
```

```python
classmethod scheil_solute_trapping()
    Configures the Scheil solute trapping settings. The used solidification speed equation is `Scanning speed * cos(angle)` with `Scanning speed` and `angle` being provided either by the user or taken from the defaults. :return: A `ScheilSoluteTrapping`
```

```python
class tc_python.scheil.ScheilOptions
    Bases: object
    Options for the Scheil simulation.
```

```python
class tc_python.scheil.ScheilOptions
    Bases: object
    Options for the Scheil simulation.
```
calculate_from_liquidus()
Solidification calculation starting from the liquidus temperature. Liquid properties between start tempera-
ture and liquidus are not obtainable.

Default: Calculation starts from liquidus temperature.

Returns This ScheilOptions object

calculate_from_start_temperature()
Calculation of equilibria from start temperature at 50 K intervals until liquidus temperature is reached. This option makes it possible to obtain properties of the liquid phase before the solidification starts.

Default: Calculation starts from liquidus temperature.

Returns This ScheilOptions object

disable_approximate_driving_force_for_metastable_phases()
Disables the approximation of the driving force for metastable phases.

Default: Enabled

Note: When enabled, the metastable phases are included in all iterations. However, these may not have reached their most favorable composition and thus their driving forces may be only approximate.

If it is important that these driving forces are correct, use disable_approximate_driving_force_for_metastable_phases() to force the calculation to converge for the metastable phases.

Returns This ScheilOptions object

disable_control_step_size_during_minimization()
Disables stepsize control during minimization (non-global).

Default: Enabled

Returns This ScheilOptions object

disable_equilibrium_solidification_calculation()
Skips the property (one axis) diagram calculation of solidification under equilibrium conditions, before the Scheil solidification calculation starts.

In general it is not necessary to perform this calculation.

Default: Disabled. The equilibrium solidification calculation is skipped.

Returns This ScheilOptions object

disable_force_positive_definite_phase_hessian()
Disables forcing of positive definite phase Hessian. This determines how the minimum of an equilibrium state in a normal minimization procedure (non-global) is reached. For details, search the Thermo-Calc documentation for “Hessian minimization”.

Default: Enabled

Returns This ScheilOptions object

enable_approximate_driving_force_for_metastable_phases()
Enables the approximation of the driving force for metastable phases.

Default: Enabled
Note: When enabled, the metastable phases are included in all iterations. However, these may not have reached their most favorable composition and thus their driving forces may be only approximate. If it is important that these driving forces are correct, use `disable_approximate_driving_force_for_metastable_phases()` to force the calculation to converge for the metastable phases.

**Returns** This `ScheilOptions` object

`enable_control_step_size_during_minimization()` Enables stepsize control during normal minimization (non-global).

Default: Enabled

**Returns** This `ScheilOptions` object

`enable_equilibrium_solidification_calculation()` Performs a property (one axis) diagram calculation of solidification under equilibrium conditions, before the Scheil solidification calculation starts, in the same way as is typically done in graphical and console mode.

In general it is not necessary to perform this calculation.

Default: Disabled. The equilibrium solidification calculation is skipped.

**Returns** This `ScheilOptions` object

`enable_force_positive_definite_phase_hessian()` Enables forcing of positive definite phase Hessian. This determines how the minimum of an equilibrium state in a normal minimization procedure (non-global) is reached. For details, search the Thermo-Calc documentation for “Hessian minimization”.

Default: Enabled

**Returns** This `ScheilOptions` object

`set_global_minimization_max_grid_points(max_grid_points: int = 2000)` Sets the maximum number of grid points in global minimization. **Only applicable if global minimization is actually used**.

Default: 2000 points

Parameters `max_grid_points` – The maximum number of grid points

**Returns** This `ScheilOptions` object

`set_global_minimization_test_interval(global_test_interval: int = 10)` Sets the interval for the global test.

Default: 10

Parameters `global_test_interval` – The global test interval

**Returns** This `ScheilOptions` object

`set_liquid_phase(phase_name: str = 'LIQUID')` Sets the phase used as the liquid phase.

Default: The phase “LIQUID”.

Parameters `phase_name` – The phase name

**Returns** This `ScheilOptions` object
**set_max_no_of_iterations** *(max_no_of_iterations: int = 500)*

Set the maximum number of iterations.

**Default:** max. 500 iterations

**Note:** As some models give computation times of more than 1 CPU second/iteration, this number is also used to check the CPU time and the calculation stops if 500 CPU seconds/iterations are used.

**Parameters** `max_no_of_iterations` – The max. number of iterations

**Returns** This `ScheilOptions` object

**set_required_accuracy** *(accuracy: float = 1e-06)*

Sets the required relative accuracy.

**Default:** 1.0E-6

**Note:** This is a relative accuracy, and the program requires that the relative difference in each variable must be lower than this value before it has converged. A larger value normally means fewer iterations but less accurate solutions. The value should be at least one order of magnitude larger than the machine precision.

**Parameters** `accuracy` – The required relative accuracy

**Returns** This `ScheilOptions` object

**set_smallest_fraction** *(smallest_fraction: float = 1e-12)*

Sets the smallest fraction for constituents that are unstable.

It is normally only in the gas phase that you can find such low fractions.

The **default value** for the smallest site-fractions is 1E-12 for all phases except for IDEAL phase with one sublattice site (such as the GAS mixture phase in many databases) for which the default value is always as 1E-30.

**Parameters** `smallest_fraction` – The smallest fraction for constituents that are unstable

**Returns** This `ScheilOptions` object

**set_temperature_step** *(temperature_step_in_kelvin: float = 1.0)*

Sets the temperature step. Decreasing the temperature step increases the accuracy, but the default value is usually adequate.

**Default step:** 1.0 K

**Parameters** `temperature_step_in_kelvin` – The temperature step [K]

**Returns** This `ScheilOptions` object

**terminate_on_fraction_of_liquid_phase** *(fraction_to_terminate_at: float = 0.01)*

Sets the termination condition to a specified remaining fraction of liquid phase.

**Default:** Terminates at 0.01 fraction of liquid phase.

**Note:** Either the termination criterion is set to a temperature or fraction of liquid limit, both together are not possible.
Parameters `fraction_to_terminate_at` – the termination fraction of liquid phase (value between 0 and 1)

Returns This `ScheilOptions` object

`terminate_on_temperature`(temperature_in_kelvin: float)
Sets the termination condition to a specified temperature.

Default: Terminates at 0.01 fraction of liquid phase, i.e. not at a specified temperature.

Note: Either the termination criterion is set to a temperature or fraction of liquid limit, both together are not possible.

Parameters `temperature_in_kelvin` – the termination temperature [K]

Returns This `ScheilOptions` object

class `tc_python.scheil.ScheilSoluteTrapping`
Bases: `tc_python.scheil.ScheilCalculationType`

Configures the Scheil solute trapping settings. The used solidification speed equation is `Scanning speed * cos(angle)` with `Scanning speed` and `angle` being provided either by the user or taken from the defaults.

`set_angle`(alpha: float = 45.0)
Sets the transformation angle alpha between the solid/liquid boundary and laser scanning direction.

Default: 45.0

Parameters `alpha` – The transformation angle [degree]

Returns This `ScheilSoluteTrapping` object

`set_primary_phasename`(primary_phase_name: str = ‘AUTOMATIC’)
Sets the name of the primary phase.

The primary phase is the phase where solute trapping takes place. A necessary condition for this phase is that the phase definition contains all of the elements that are chosen in the system. When `AUTOMATIC` is selected, the program tries to find a suitable primary phase that fills this condition.

Default: `AUTOMATIC`

Parameters `primary_phase_name` – The phase name (or `AUTOMATIC`)

Returns This `ScheilSoluteTrapping` object

`set_scanning_speed`(scanning_speed: float = 1.0)
Sets the scanning speed.

Default: 1 m/s

Parameters `scanning_speed` – The scaling factor [m/s]

Returns This `ScheilSoluteTrapping` object
5.1.5 Module “step_or_map_diagrams”

```python
class tc_python.step_or_map_diagrams.AbstractAxisType
    Bases: object
    The abstract base class for all axis types.

class tc_python.step_or_map_diagrams.AbstractPhaseDiagramCalculation(calculator)
    Bases: tc_python.abstract_base.AbstractCalculation
    Abstract configuration required for a property diagram calculation.
```

**Note:** This is an abstract class that cannot be used directly.

```python
add_initial_equilibrium(initial_equilibrium: tc_python.step_or_map_diagrams.InitialEquilibrium)
    Add initial equilibrium start points from which a phase diagram is calculated.
    Scans along the axis variables and generates start points when the scan procedure crosses a phase boundary.
    It may take a little longer to execute than using the minimum number of start points, as some lines may be calculated more than once. But the core remembers all node points and subsequently stops calculations along a line when it finds a known node point.
    It is also possible to create a sequence of start points from one initial equilibria.

    **Parameters**
    initial_equilibrium – The initial equilibrium

    **Returns**
    This `PhaseDiagramCalculation` object

abstract calculate(keep_previous_results: bool = False) → tc_python.step_or_map_diagrams.PhaseDiagramResult
    Disables global minimization.

    **Default**
    Enabled

    **Returns**
    This `PhaseDiagramCalculation` object

dont_keep_default_equilibria()
    Do not keep the initial equilibria added by default.
    This is only relevant in combination with `add_initial_equilibrium()`.
    This is the default behavior.

    **Returns**
    This `PhaseDiagramCalculation` object

enable_global_minimization()
    Enables global minimization.

    **Default**
    Enabled

    **Returns**
    This `PhaseDiagramCalculation` object

get_components() → List[str]
    Returns the names of the components in the system (including all components auto-selected by the database(s)).

    **Returns**
    The component names

get_gibbs_energy_addition_for(phase: str) → float
    Used to get the additional energy term (always being a constant) of a given phase. The value given is added
to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface
tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters**  
**phase** – Specify the name of the (stoichiometric or solution) phase with the addition

**Returns**  
Gibbs energy addition to G per mole formula unit.

### get_system_data() → tc_python.abstract_base.SystemData

Returns the content of the database for the currently loaded system. This can be used to modify the param-
eters and functions and to change the current system by using `with_system_modifications()`.

**Note:** Parameters can only be read from unencrypted (i.e. `user`) databases loaded as `*.tdb-file`.

**Returns**  
The system data

### keep_default_equilibria()

Keep the initial equilibria added by default. This is only relevant in combination with
`add_initial_equilibrium()`.

Default behavior is to not keep default equilibria.

**Returns**  
This `PhaseDiagramCalculation` object

### remove_all_initial_equilibria()

Removes all previously added initial equilibria.

**Returns**  
This `PhaseDiagramCalculation` object

### run_poly_command(command: str)

Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

**Parameters**  
**command** – The Thermo-Calc Console Mode command

**Returns**  
This `PhaseDiagramCalculation` object

**Note:** It should not be necessary for most users to use this method, try to use the corresponding method
implemented in the API instead.

**Warning:** As this method runs raw Thermo-Calc commands directly in the engine, it may hang the
program in case of spelling mistakes (e.g. forgotten equals sign).

### set_gibbs_energy_addition_for(phase: str, gibbs_energy: float)

Used to specify the additional energy term (always being a constant) of a given phase. The value
(`gibbs_energy`) given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent
a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters**

- **phase** – Specify the name of the (stoichiometric or solution) phase with the addition

- **gibbs_energy** – Addition to G per mole formula unit

**Returns**  
This `PhaseDiagramCalculation` object
**set_phase_to_dormant** *(phase: str)*
Sets the phase to the status DORMANT, necessary for calculating the driving force to form the specified phase.

**Parameters**
- **phase** – The phase name or **ALL_PHASES** for all phases

**Returns**
This **PhaseDiagramCalculation** object

**set_phase_to_entered** *(phase: str, amount: float = 1.0)*
Sets the phase to the status ENTERED, that is the default state.

**Parameters**
- **phase** – The phase name or **ALL_PHASES** for all phases
- **amount** – The phase fraction (between 0.0 and 1.0)

**Returns**
This **PhaseDiagramCalculation** object

**set_phase_to_fixed** *(phase: str, amount: float)*
Sets the phase to the status FIXED, i.e. it is guaranteed to have the specified phase fraction after the calculation.

**Parameters**
- **phase** – The phase name
- **amount** – The fixed phase fraction (between 0.0 and 1.0)

**Returns**
This **PhaseDiagramCalculation** object

**set_phase_to_suspended** *(phase: str)*
Sets the phase to the status SUSPENDED, i.e. it is ignored in the calculation.

**Parameters**
- **phase** – The phase name or **ALL_PHASES** for all phases

**Returns**
This **PhaseDiagramCalculation** object

**with_options** *(options: tc_python.step_or_map_diagrams.PhaseDiagramOptions)*
Sets the simulation options.

**Parameters**
- **options** – The simulation options

**Returns**
This **PhaseDiagramCalculation** object

**with_reference_state** *(component: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)*
The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.

All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to set the reference state for a component using **SER**, i.e. the Stable Element Reference (which is usually set as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.
If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

**Parameters**

- **component** – The name of the element must be given.
- **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.
- **temperature** – The Temperature (in K) for the reference state. Or CURRENT_TEMPERATURE which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.
- **pressure** – The Pressure (in Pa) for the reference state.

**Returns** This PhaseDiagramCalculation object

with_system_modifications(system_modifications: tc_python.abstract_base.SystemModifications)

Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

**Note:** This is only possible if the system has been read from unencrypted (i.e. user) databases loaded as a *.tdb*-file.

**Parameters** system_modifications – The system modification to be performed

**Returns** This PhaseDiagramCalculation object

class tc_python.step_or_map_diagrams.AbstractPropertyDiagramCalculation(calculator)

Bases: tc_python.abstract_base.AbstractCalculation

Abstract configuration required for a property diagram calculation.

**Note:** This is an abstract class that cannot be used directly.

**abstract calculate(keep_previous_results: bool = False)**

**disable_global_minimization()**

Disables global minimization.

**Default:** Enabled

**Returns** This PropertyDiagramCalculation object

**disable_step_separate_phases()**

Disables step separate phases. This is the default setting.

**Returns** This PropertyDiagramCalculation object

**enable_global_minimization()**

Enables global minimization.

**Default:** Enabled

**Returns** This PropertyDiagramCalculation object

**enable_step_separate_phases()**

Enables step separate phases.

**Default:** By default separate phase stepping is disabled
Note: This is an advanced option, it is used mostly to calculate how the Gibbs energy for a number of phases varies for different compositions. This is particularly useful to calculate Gibbs energies for complex phases with miscibility gaps and for an ordered phase that is never disordered (e.g. SIGMA-phase, G-phase, MU-phase, etc.).

Returns This PropertyDiagramCalculation object

get_components() → List[str]
Returns the names of the components in the system (including all components auto-selected by the database(s)).

Returns The component names

get_gibbs_energy_addition_for(phase: str) → float
Used to get the additional energy term (always being a constant) of a given phase. The value given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

Parameters phase – Specify the name of the (stoichiometric or solution) phase with the addition

Returns Gibbs energy addition to G per mole formula unit.

get_system_data() → tc_python.abstract_base.SystemData
Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using with_system_modifications().

Note: Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb-file.

Returns The system data

run_poly_command(command: str)
Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

Parameters command – The Thermo-Calc Console Mode command

Returns This PropertyDiagramCalculation object

Note: It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

Warning: As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).

set_gibbs_energy_addition_for(phase: str, gibbs_energy: float)
Used to specify the additional energy term (always being a constant) of a given phase. The value (gibbs_energy) given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.
Parameters

- **phase** – Specify the name of the (stoichiometric or solution) phase with the addition
- **gibbs_energy** – Addition to G per mole formula unit

Returns This `PropertyDiagramCalculation` object

`set_phase_to_dormant(phase: str)`

Sets the phase to the status DORMANT, necessary for calculating the driving force to form the specified phase.

Parameters **phase** – The phase name or `ALL_PHASES` for all phases

Returns This `PropertyDiagramCalculation` object

`set_phase_to_entered(phase: str, amount: float = 1.0)`

Sets the phase to the status ENTERED, that is the default state.

Parameters

- **phase** – The phase name or `ALL_PHASES` for all phases
- **amount** – The phase fraction (between 0.0 and 1.0)

Returns This `PropertyDiagramCalculation` object

`set_phase_to_fixed(phase: str, amount: float)`

Sets the phase to the status FIXED, i.e. it is guaranteed to have the specified phase fraction after the calculation.

Parameters

- **phase** – The phase name
- **amount** – The fixed phase fraction (between 0.0 and 1.0)

Returns This `PropertyDiagramCalculation` object

`set_phase_to_suspended(phase: str)`

Sets the phase to the status SUSPENDED, i.e. it is ignored in the calculation.

Parameters **phase** – The phase name or `ALL_PHASES` for all phases

Returns This `PropertyDiagramCalculation` object

`with_options(options: tc_python.step_or_map_diagrams.PropertyDiagramOptions)`

Sets the simulation options.

Parameters **options** – The simulation options

Returns This `PropertyDiagramCalculation` object

`with_reference_state(component: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)`

The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.

All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set
as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

**Parameters**

- **component** – The name of the element must be given.
- **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.
- **temperature** – The Temperature (in K) for the reference state. Or CURRENT_TEMPERATURE which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.
- **pressure** – The Pressure (in Pa) for the reference state.

**Returns**

This PropertyDiagramCalculation object

**with_system_modifications**(system_modifications: tc_python.abstract_base.SystemModifications)

Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

**Parameters**

*system_modifications* – The system modification to be performed

**Returns**

This PropertyDiagramCalculation object

---

**class** tc_python.step_or_map_diagrams.AxisType

**Bases:** tc_python.step_or_map_diagrams.AbstractAxisType

Factory class providing objects for configuring a logarithmic or linear axis by using `AxisType.linear()` or `AxisType.logarithmic()`.

**classmethod linear()**

Creates an object for configuring a linear calculation axis.

**Default:** A minimum number of 40 steps.

**Note:** The returned object can be configured regarding the maximum step size or the minimum number of steps on the axis.

**Returns**

A new Linear object

**classmethod logarithmic()**

Creates an object for configuring a logarithmic calculation axis.

**Default:** A scale factor of 1.1

---

5.1. Calculations
Note: The returned object can be configured regarding the scale factor.

Returns A new Logarithmic object

class tc_python.step_or_map_diagrams.CalculationAxis (quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity,
str])
Bases: object
A calculation axis used for property and phase diagram calculations.

Note: A calculation axis is defining the varied condition and the range of variation. It is the same concept as in Thermo-Calc Graphical Mode or Console Mode.

Default: A Linear axis with a minimum number of 40 steps

set_max (max: float)
Sets the maximum quantity value of the calculation axis.

There is no default value set, it always needs to be defined.

Parameters max – The maximum quantity value of the axis [unit according to the axis quantity]

Returns This CalculationAxis object

set_min (min: float)
Sets the minimum quantity value of the calculation axis.

There is no default value set, it always needs to be defined.

Parameters min – The minimum quantity value of the axis [unit according to the axis quantity]

Returns This CalculationAxis object

set_start_at (at: float)
Sets the starting point of the calculation on the axis.

Default: The default starting point is the center between the minimum and maximum quantity value

Parameters at – The starting point on the axis [unit according to the axis quantity]

Returns This CalculationAxis object

with_axis_type (axis_type: tc_python.step_or_map_diagrams.AxisType)
Sets the type of the axis.

Default: A Linear axis with a minimum number of 40 steps

Parameters axis_type – The axis type (linear or logarithmic)

Returns This CalculationAxis object

class tc_python.step_or_map_diagrams.Direction (value)
Bases: enum.Enum
An enumeration.

DECREASE_FIRST_AXIS = 3
DECREASE_SECOND_AXIS = 4
INCREASE_FIRST_AXIS = 0
INCREASE_SECOND_AXIS = 1

class tc_python.step_or_map_diagrams.InitialEquilibrium(first_axis: float, second_axis: float):
    Bases: object

    add_equilibria_at_all_phase_changes()
    This generates one start point for each set of phase change in the chosen direction of the specified axis.
    This ensures finding all possible phase boundary lines (not just the first one) along such an axis direction.
    Default behavior is to only generate one start point at the first phase change.
    Returns This InitialEquilibrium object

    add_equilibria_at_first_phase_change()
    This generates one start point at the first phase change.
    This is the default behavior.
    Returns This InitialEquilibrium object

    set_direction(direction_enum: tc_python.step_or_map_diagrams.Direction)
    Specifies along which axes the initial equilibria should be added.
    The default direction is INCREASE_FIRST_AXIS.
    Parameters direction_enum
    Returns This InitialEquilibrium object

class tc_python.step_or_map_diagrams.Linear
    Bases: tc_python.step_or_map_diagrams.AxisType

    Represents a linear axis.

    get_type() → str
    Convenience method for getting axis type.
    Returns The type

    set_max_step_size(max_step_size: float)
    Sets the axis to use the maximum step size configuration.
    Default: This is not the default which is minimum number of steps
    Note: Either maximum step size or minimum number of steps can be used but not both at the same time.
    Parameters max_step_size – The maximum step size [unit according to the axis quantity]
    Returns This Linear object

    set_min_nr_of_steps(min_nr_of_steps: float = 40)
    Sets the axis to use the minimum number of steps configuration.
    Default: This is the default option (with a minimum number of steps of 40)
    Note: Either maximum step size or minimum number of steps can be used but not both at the same time.
    Parameters min_nr_of_steps – The minimum number of steps
    Returns This Linear object
class tc_python.step_or_map_diagrams.Logarithmic(scale_factor: float = 1.1)
    Bases: tc_python.step_or_map_diagrams.AxisType
    Represents a logarithmic axis.

    Note: A logarithmic axis is useful for low fractions like in a gas phase where 1E-7 to 1E-2 might be an interesting range. For the pressure a logarithmic axis is often also useful.

get_type() -> str
    Convenience method for getting axis type.

    Returns The type

set_scale_factor(scale_factor: float = 1.1)
    Sets the scale factor.

    Default: 1.1

    Parameters scale_factor -- The scale factor setting the maximum factor between two calculated values, must be larger than 1.0

    Returns This Logarithmic object

class tc_python.step_or_map_diagrams.PhaseDiagramCalculation(calculator)
    Bases: tc_python.step_or_map_diagrams.AbstractPhaseDiagramCalculation
    Configuration for a phase diagram calculation.

    Note: Specify the conditions, the calculation is performed with calculate().

add_initial_equilibrium(initial_equilibrium: tc_python.step_or_map_diagrams.InitialEquilibrium)
    Add initial equilibrium start points from which a phase diagram is calculated.

    Scans along the axis variables and generates start points when the scan procedure crosses a phase boundary.

    It may take a little longer to execute than using the minimum number of start points, as some lines may be calculated more than once. But the core remembers all node points and subsequently stops calculations along a line when it finds a known node point.

    It is also possible to create a sequence of start points from one initial equilibria.

    Parameters initial_equilibrium -- The initial equilibrium

    Returns This PhaseDiagramCalculation object

calculate(keep_previous_results: bool = False) -> tc_python.step_or_map_diagrams.PhaseDiagramResult
    Performs the phase diagram calculation.

    Warning: If you use keep_previous_results=True, you must not use another calculator or even get results in between the calculations using calculate(). Then the previous results will actually be lost.

    Parameters keep_previous_results -- If True, results from any previous call to this method are appended. This can be used to combine calculations with multiple start points if the mapping fails at a certain condition.

    Returns A new PhaseDiagramResult object which later can be used to get specific values from the calculated result.
disable_global_minimization()
Disables global minimization.

Default: Enabled

Returns This PhaseDiagramCalculation object

dont_keep_default_equilibria()
Do not keep the initial equilibria added by default.
This is only relevant in combination with add_initial_equilibrium().
This is the default behavior.

Returns This PhaseDiagramCalculation object

enable_global_minimization()
Enables global minimization.

Default: Enabled

Returns This PhaseDiagramCalculation object

get_components() \rightarrow \text{List[\text{str}]} 
Returns the names of the components in the system (including all components auto-selected by the database(s)).

Returns The component names

get_gibbs_energy_addition_for(phase: str) \rightarrow \text{float} 
Used to get the additional energy term (always being a constant) of a given phase. The value given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

Parameters phase – Specify the name of the (stoichiometric or solution) phase with the addition

Returns Gibbs energy addition to G per mole formula unit.

get_system_data() \rightarrow \text{tc_python.abstract_base.SystemData} 
Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using with_system_modifications().

Note: Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb-file.

Returns The system data

keep_default_equilibria()
Keep the initial equilibria added by default. This is only relevant in combination with add_initial_equilibrium().
Default behavior is to not keep default equilibria.

Returns This PhaseDiagramCalculation object

remove_all_conditions()
Removes all set conditions.

Returns This PhaseDiagramCalculation object
remove_all_initial_equilibria()
Removes all previously added initial equilibria.

Returns This PhaseDiagramCalculation object

remove_condition(quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str])
Removes the specified condition.

Parameters quantity – The thermodynamic quantity to set as condition; a Console Mode syntax string can be used as an alternative (for example X(Cr))

Returns This ThermodynamicCalculation object

run_poly_command(command: str)
Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

Parameters command – The Thermo-Calc Console Mode command

Returns This PhaseDiagramCalculation object

Note: It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

Warning: As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).

set_condition(quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], value: float)
Sets the specified condition.

Parameters

• quantity – The thermodynamic quantity to set as condition; a Console Mode syntax string can be used as an alternative (for example X(Cr))

• value – The value of the condition

Returns This PhaseDiagramCalculation object

set_gibbs_energy_addition_for(phase: str, gibbs_energy: float)
Used to specify the additional energy term (always being a constant) of a given phase. The value (gibbs_energy) given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

Parameters

• phase – Specify the name of the (stoichiometric or solution) phase with the addition

• gibbs_energy – Addition to G per mole formula unit

Returns This PhaseDiagramCalculation object

set_phase_to_dormant(phase: str)
Sets the phase to the status DORMANT, necessary for calculating the driving force to form the specified phase.

Parameters phase – The phase name or ALL_PHASES for all phases

Returns This PhaseDiagramCalculation object
**set_phase_to_entered** *(phase: str, amount: float = 1.0)*

Sets the phase to the status ENTERED, that is the default state.

**Parameters**

- **phase** – The phase name or `ALL_PHASES` for all phases
- **amount** – The phase fraction (between 0.0 and 1.0)

**Returns** This `PhaseDiagramCalculation` object

**set_phase_to_fixed** *(phase: str, amount: float)*

Sets the phase to the status FIXED, i.e. it is guaranteed to have the specified phase fraction after the calculation.

**Parameters**

- **phase** – The phase name
- **amount** – The fixed phase fraction (between 0.0 and 1.0)

**Returns** This `PhaseDiagramCalculation` object

**set_phase_to_suspended** *(phase: str)*

Sets the phase to the status SUSPENDED, i.e. it is ignored in the calculation.

**Parameters**

- **phase** – The phase name or `ALL_PHASES` for all phases

**Returns** This `PhaseDiagramCalculation` object

**with_first_axis** *(axis: tc_python.step_or_map_diagrams.CalculationAxis)*

Sets the first calculation axis.

**Parameters**

- **axis** – The axis

**Returns** This `PhaseDiagramCalculation` object

**with_options** *(options: tc_python.step_or_map_diagrams.PhaseDiagramOptions)*

Sets the simulation options.

**Parameters**

- **options** – The simulation options

**Returns** This `PhaseDiagramCalculation` object

**with_reference_state** *(component: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)*

The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.

All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

---

**5.1. Calculations**
Parameters

• **component** – The name of the element must be given.

• **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.

• **temperature** – The Temperature (in K) for the reference state. Or CURRENT_TEMPERATURE which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.

• **pressure** – The Pressure (in Pa) for the reference state.

Returns This *PhaseDiagramCalculation* object

with_second_axis *(axis: tc_python.step_or_map_diagrams.CalculationAxis)*
Sets the second calculation axis.

Parameters **axis** – The axis

Returns This *PhaseDiagramCalculation* object

with_system_modifications *(system_modifications: tc_python.abstract_base.SystemModifications)*
Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

Note: This is only possible if the system has been read from unencrypted (i.e. *user*) databases loaded as a *.tdb*-file.

Parameters **system_modifications** – The system modification to be performed

Returns This *PhaseDiagramCalculation* object

---

**class** *tc_python.step_or_map_diagrams.PhaseDiagramOptions*

*Simulation options for phase diagram calculations.*

**disable_approximate_driving_force_for_metastable_phases()**
Disables the approximation of the driving force for metastable phases.

Default: Enabled

Note: When enabled, the metastable phases are included in all iterations. However, these may not have reached their most favorable composition and thus their driving forces may be only approximate.

If it is important that these driving forces are correct, use *disable_approximate_driving_force_for_metastable_phases()* to force the calculation to converge for the metastable phases.

Returns This *PhaseDiagramOptions* object

**disable_control_step_size_during_minimization()**
Disables stepsize control during minimization (non-global).

Default: Enabled

Returns This *PhaseDiagramOptions* object
disable_force_positive_definite_phase_hessian()
Disables forcing of positive definite phase Hessian. This determines how the minimum of an equilibrium state in a normal minimization procedure (non-global) is reached. For details, search the Thermo-Calc documentation for “Hessian minimization”.

Default: Enabled

Returns This PhaseDiagramOptions object

dont_use_auto_start_points()
Switches the usage of automatic starting points for the mapping off.

Default: Switched on

Returns This PhaseDiagramOptions object

dont_use_inside_mesh_points()
Switches the usage of inside meshing points for the mapping off.

Default: Switched off

Returns This PhaseDiagramOptions object

enable_approximate_driving_force_for_metastable_phases()
Enables the approximation of the driving force for metastable phases.

Default: Enabled

Note: When enabled, the metastable phases are included in all iterations. However, these may not have reached their most favorable composition and thus their driving forces may be only approximate.

If it is important that these driving forces are correct, use disable_approximate_driving_force_for_metastable_phases() to force the calculation to converge for the metastable phases.

Returns This PhaseDiagramOptions object

enable_control_step_size_during_minimization()
Enables stepsize control during normal minimization (non-global).

Default: Enabled

Returns This PhaseDiagramOptions object

enable_force_positive_definite_phase_hessian()
Enables forcing of positive definite phase Hessian. This determines how the minimum of an equilibrium state in a normal minimization procedure (non-global) is reached. For details, search the Thermo-Calc documentation for “Hessian minimization”.

Default: Enabled

Returns This PhaseDiagramOptions object

set_global_minimization_max_grid_points(max_grid_points: int = 2000)
Sets the maximum number of grid points in global minimization. ** Only applicable if global minimization is actually used**.

Default: 2000 points

Parameters max_grid_points – The maximum number of grid points

Returns This PhaseDiagramOptions object

5.1. Calculations
**set_global_minimization_test_interval** *(global_test_interval: int = 0)*
Sets the interval for the global test.

Default: 0

Parameters *global_test_interval* – The global test interval

Returns This *PhaseDiagramOptions* object

**set_max_no_of_iterations** *(max_no_of_iterations: int = 500)*
Set the maximum number of iterations.

Default: max. 500 iterations

**Note:** As some models give computation times of more than 1 CPU second/iteration, this number is also used to check the CPU time and the calculation stops if 500 CPU seconds/iterations are used.

Parameters *max_no_of_iterations* – The max. number of iterations

Returns This *PhaseDiagramOptions* object

**set_no_of_mesh_along_axis** *(no_of_mesh_along_axis: int = 3)*
Sets the number of meshes along an axis for the mapping.

Default: 3

Parameters *no_of_mesh_along_axis* – The number of meshes

Returns This *PhaseDiagramOptions* object

**set_required_accuracy** *(accuracy: float = 1e-06)*
Sets the required relative accuracy.

Default: 1.0E-6

**Note:** This is a relative accuracy, and the program requires that the relative difference in each variable must be lower than this value before it has converged. A larger value normally means fewer iterations but less accurate solutions. The value should be at least one order of magnitude larger than the machine precision.

Parameters *accuracy* – The required relative accuracy

Returns This *PhaseDiagramOptions* object

**set_smallest_fraction** *(smallest_fraction: float = 1e-12)*
Sets the smallest fraction for constituents that are unstable.

It is normally only in the gas phase that you can find such low fractions.

The default value for the smallest site-fractions is 1E-12 for all phases except for IDEAL phase with one sublattice site (such as the GAS mixture phase in many databases) for which the default value is always as 1E-30.

Parameters *smallest_fraction* – The smallest fraction for constituents that are unstable

Returns This *PhaseDiagramOptions* object
use_auto_start_points ()
  Switches the usage of automatic starting points for the mapping on.
  Default: Switched on
  Returns This PhaseDiagramOptions object

use_inside_mesh_points ()
  Switches the usage of inside meshing points for the mapping off.
  Default: Switched off
  Returns This PhaseDiagramOptions object

class tc_python.step_or_map_diagrams.PhaseDiagramResult (result)
  Bases: tc_python.abstract_base.AbstractResult
  Result of a phase diagram calculation, it can be evaluated using quantities or Console Mode syntax.

add_coordinate_for_phase_label (x: float, y: float)
  Sets a coordinate in the result plot for which the stable phases will be evaluated and provided in the
  result data object. This can be used to plot the phases of a region into the phase diagram or just to
  programmatically evaluate the phases in certain regions.

  Warning: This method takes coordinates of the plot axes and not of the calculation axis.

Parameters
• x – The coordinate of the first plot axis (“x-axis”) [unit of the plot axis]
• y – The coordinate of the second plot axis (“y-axis”) [unit of the plot axis]

  Returns This PhaseDiagramResult object

get_values_grouped_by_quantity_of (x_quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], y_quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str]) \rightarrow tc_python.step_or_map_diagrams.PhaseDiagramResultValues

Returns x-y-line data grouped by the multiple datasets of the specified quantities (for example in depen-
dency of components). The available quantities can be found in the documentation of the factory class
ThermodynamicQuantity. Usually the result data represents the phase diagram.

Note: The different datasets will contain NaN-values between different subsections and are not sorted
(because they are unsortable due to their nature).

Note: It's possible to use functions as axis variables, either by using ThermodynamicQuan-
tity.user_defined_function, or by using an expression that contains ‘=’.

Example get_values_grouped_by_quantity_of('T', ThermodynamicQuantity.user_defined_function('HM.T'))

Example get_values_grouped_by_quantity_of('T', 'CP=HM.T')

Parameters
• x_quantity – The first quantity (“x-axis”), Console Mode syntax strings can be used
  as an alternative (for example ‘T’), or even a function (for example ‘f=T*1.01’)


- **y_quantity** – The second quantity (“y-axis”), Console Mode syntax strings can be used as an alternative (for example ‘NV’), or even a function (for example ‘CP=HM.T’)

**Returns** The phase diagram data

get_values_grouped_by_stable_phases_of(x_quantity: \(\text{Union}\left[\text{tc\_python\_quantity\_factory.ThermodynamicQuantity}, \text{str}\right]\), y_quantity: \(\text{Union}\left[\text{tc\_python\_quantity\_factory.ThermodynamicQuantity}, \text{str}\right]\) \(\rightarrow\) \(\text{tc\_python\_step\_or\_map\_diagrams.PhaseDiagramResultValues}\)

Returns x-y-line data grouped by the sets of “stable phases” (for example “LIQUID” or “LIQUID + FCC_A1”). The available quantities can be found in the documentation of the factory class ThermodynamicQuantity. Usually the result data represents the phase diagram.

**Note:** The different datasets will contain NaN-values between different subsections and are not sorted (because they are unsortable due to their nature).

**Note:** It’s possible to use functions as axis variables, either by using ThermodynamicQuantity.user_defined_function, or by using an expression that contains ‘=’.

**Example**

```
get_values_grouped_by_quantity_of(‘T’, ThermodynamicQuantity.user_defined_function(‘HM.T’))
```

**Example**

```
get_values_grouped_by_quantity_of(‘T’, ‘CP=HM.T’)
```

**Parameters**

- **x_quantity** – The first quantity (“x-axis”), Console Mode syntax strings can be used as an alternative (for example ‘T’), or even a function (for example ‘f=T*1.01’)

- **y_quantity** – The second quantity (“y-axis”), Console Mode syntax strings can be used as an alternative (for example ‘NV’), or even a function (for example ‘CP=HM.T’)

**Returns** The phase diagram data

remove_phase_labels()

Erases all added coordinates for phase labels.

**Returns** This PhaseDiagramResult object

save_to_disk(path: str)

Saves the result to disc. Note that a result is a folder, containing potentially many files. The result can later be loaded with load_result_from_disk()

**Parameters** path – the path to the folder you want the result to be saved in. It can be relative or absolute.

**Returns** this PhaseDiagramResult object

set_phase_name_style(phase_name_style_enum: \(\text{tc\_python\_step\_or\_map\_diagrams.PhaseNameStyle}\) = \(<\text{PhaseNameStyle.NONE: 0}>\))

Sets the style of the phase name labels that will be used in the result data object (constitution description, ordering description, . . . ).

**Default:** PhaseNameStyle.NONE

**Parameters** phase_name_style_enum – The phase name style

**Returns** This PhaseDiagramResult object
class `tc_python.step_or_map_diagrams.PhaseDiagramResultValues`

Bases: `object`

Represents the data of a phase diagram.

**get_invariants() → `tc_python.utils.ResultValueGroup`**

Returns the x- and y-datasets of all invariants in the phase diagram.

**Note:** The datasets will normally contain different sections separated by NaN-values.

**Returns** The invariants dataset object

**get_lines() → Dict[str, `tc_python.utils.ResultValueGroup`]**

Returns the x- and y-datasets of all phase boundaries in the phase diagram.

**Note:** The datasets will normally contain different sections separated by NaN-values.

**Returns** Containing the phase boundary datasets with the quantities or stable phases as keys (depending on the used method to get the values)

**get_phase_labels() → List[`tc_python.step_or_map_diagrams.PhaseLabel`]**

Returns the phase labels added for certain coordinates using `PhaseDiagramResult.add_coordinate_for_phase_label()`.

**Returns** The list with the phase label data (that contains plot coordinates and stable phases)

**get_tie_lines() → `tc_python.utils.ResultValueGroup`**

Returns the x- and y-datasets of all tie-lines in the phase diagram.

**Note:** The datasets will normally contain different sections separated by NaN-values.

**Returns** The tie-line dataset object

---

**class `tc_python.step_or_map_diagrams.PhaseLabel`**

Bases: `object`

Represents a phase label at a plot coordinate, i.e. the stable phases that are present at that plot coordinate.

**get_text() → str**

Accessor for the phase label :return: the phase label

**get_x() → List[float]**

Accessor for the x-value :return: the x value

**get_y() → List[float]**

Accessor for the y-value :return: the y value

**class `tc_python.step_or_map_diagrams.PhaseNameStyle`**

Bases: `enum.Enum`

The style of the phase names used in the labels.

**ALL = 1**

Adding ordering and constitution description.
CONSTITUTION_DESCRIPTION = 3
Adding only constitution description.

NONE = 0
Only the phase names.

ORDERING_DESCRIPTION = 4
Adding only ordering description.

class tc_python.step_or_map_diagrams.PropertyDiagramCalculation(calculator)
Bases: tc_python.step_or_map_diagrams.AbstractPropertyDiagramCalculation

calculate (keep_previous_results: bool = False) → tc_python.step_or_map_diagrams.PropertyDiagramResult
Performs the property diagram calculation.

**Warning:** If you use `keep_previous_results=True`, you must not use another calculator or even get results in between the calculations using `calculate()`. Then the previous results will actually be lost.

**Parameters** `keep_previous_results` – If `True`, results from any previous call to this method are appended. This can be used to combine calculations with multiple start points if the stepping fails at a certain condition.

**Returns** A new `PropertyDiagramResult` object which later can be used to get specific values from the calculated result

disable_global_minimization ()
Disables global minimization.

**Default:** Enabled

**Returns** This `PropertyDiagramCalculation` object

disable_step_separate_phases ()
Disables step separate phases. This is the default setting.

**Returns** This `PropertyDiagramCalculation` object

enable_global_minimization ()
Enables global minimization.

**Default:** Enabled

**Returns** This `PropertyDiagramCalculation` object

enable_step_separate_phases ()
Enables step separate phases.

**Default:** By default separate phase stepping is disabled

**Note:** This is an advanced option, it is used mostly to calculate how the Gibbs energy for a number of phases varies for different compositions. This is particularly useful to calculate Gibbs energies for complex phases with miscibility gaps and for an ordered phase that is never disordered (e.g. SIGMA-phase, G-phase, MU-phase, etc.).

**Returns** This `PropertyDiagramCalculation` object
**get_components()** → List[str]
Returns the names of the components in the system (including all components auto-selected by the database(s)).

Returns The component names

**get_gibbs_energy_addition_for**(phase: str) → float
Used to get the additional energy term (always being a constant) of a given phase. The value given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

Parameters **phase** – Specify the name of the (stoichiometric or solution) phase with the addition

Returns Gibbs energy addition to G per mole formula unit.

**get_system_data()** → tc_python.abstract_base.SystemData
Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using with_system_modifications().

Note: Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb*-file.

Returns The system data

**remove_all_conditions()**
Removes all set conditions.

Returns This PropertyDiagramCalculation object

**remove_condition**(quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str])
Removes the specified condition.

Parameters **quantity** – The thermodynamic quantity to set as condition; a Console Mode syntax string can be used as an alternative (for example X(Cr))

Returns This PropertyDiagramCalculation object

**run_poly_command**(command: str)
Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

Parameters **command** – The Thermo-Calc Console Mode command

Returns This PropertyDiagramCalculation object

Note: It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

Warning: As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).

**set_condition**(quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], value: float)
Sets the specified condition.

Parameters
• **quantity** – The thermodynamic quantity to set as condition; a Console Mode syntax string can be used as an alternative (for example \( X(\text{Cr}) \))

• **value** – The value of the condition

**Returns** This `PropertyDiagramCalculation` object

**set_gibbs_energy_addition_for** *(phase: str, gibbs_energy: float)*

Used to specify the additional energy term (always being a constant) of a given phase. The value (\( \text{gibbs\_energy} \)) given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters**

• **phase** – Specify the name of the (stoichiometric or solution) phase with the addition

• **gibbs_energy** – Addition to G per mole formula unit

**Returns** This `PropertyDiagramCalculation` object

**set_phase_to_dormant** *(phase: str)*

Sets the phase to the status DORMANT, necessary for calculating the driving force to form the specified phase.

**Parameters** phase – The phase name or `ALL_PHASES` for all phases

**Returns** This `PropertyDiagramCalculation` object

**set_phase_to_entered** *(phase: str, amount: float = 1.0)*

Sets the phase to the status ENTERED, that is the default state.

**Parameters**

• **phase** – The phase name or `ALL_PHASES` for all phases

• **amount** – The phase fraction (between 0.0 and 1.0)

**Returns** This `PropertyDiagramCalculation` object

**set_phase_to_fixed** *(phase: str, amount: float)*

Sets the phase to the status FIXED, i.e. it is guaranteed to have the specified phase fraction after the calculation.

**Parameters**

• **phase** – The phase name

• **amount** – The fixed phase fraction (between 0.0 and 1.0)

**Returns** This `PropertyDiagramCalculation` object

**set_phase_to_suspended** *(phase: str)*

Sets the phase to the status SUSPENDED, i.e. it is ignored in the calculation.

**Parameters** phase – The phase name or `ALL_PHASES` for all phases

**Returns** This `PropertyDiagramCalculation` object

**with_axis** *(axis: tc_python.step_or_map_diagrams.CalculationAxis)*

Sets the calculation axis.

**Parameters** axis – The axis

**Returns** This `PropertyDiagramCalculation` object
**with_options** *(options: tc_python.step_or_map_diagrams.PropertyDiagramOptions)*

Sets the simulation options.

**Parameters**

- **options** – The simulation options

**Returns**

This *PropertyDiagramCalculation* object

**with_reference_state** *(component: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)*

The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.

All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

**Parameters**

- **component** – The name of the element must be given.
- **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.
- **temperature** – The Temperature (in K) for the reference state. Or **CURRENT_TEMPERATURE** which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.
- **pressure** – The Pressure (in Pa) for the reference state.

**Returns**

This *PropertyDiagramCalculation* object

**with_system_modifications** *(system_modifications: tc_python.abstract_base.SystemModifications)*

Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

**Parameters**

- **system_modifications** – The system modification to be performed

**Returns**

This *PropertyDiagramCalculation* object

---

**Note:** This is only possible if the system has been read from unencrypted (i.e. *user*) databases loaded as a *.tdb*-file.

---

**Parameters**

- **system_modifications** – The system modification to be performed

**Returns**

This *PropertyDiagramCalculation* object

---

**class** *tc_python.step_or_map_diagrams.PropertyDiagramOptions*

**Bases:** *object*

Simulation options for the property diagram calculations.
disable_approximate_driving_force_for_metastable_phases()  
Disables the approximation of the driving force for metastable phases.  

**Default:** Enabled

**Note:** When enabled, the metastable phases are included in all iterations. However, these may not have reached their most favorable composition and thus their driving forces may be only approximate. If it is important that these driving forces are correct, use disable_approximate_driving_force_for_metastable_phases() to force the calculation to converge for the metastable phases.

**Returns** This PropertyDiagramOptions object

disable_control_step_size_during_minimization()  
Disables stepsize control during minimization (non-global).  

**Default:** Enabled

**Returns** This PropertyDiagramOptions object

disable_force_positive_definite_phase_hessian()  
Disables forcing of positive definite phase Hessian. This determines how the minimum of an equilibrium state in a normal minimization procedure (non-global) is reached. For details, search the Thermo-Calc documentation for “Hessian minimization”.  

**Default:** Enabled

**Returns** This PropertyDiagramOptions object

enable_approximate_driving_force_for_metastable_phases()  
Enables the approximation of the driving force for metastable phases.  

**Default:** Enabled

**Note:** When enabled, the metastable phases are included in all iterations. However, these may not have reached their most favorable composition and thus their driving forces may be only approximate. If it is important that these driving forces are correct, use disable_approximate_driving_force_for_metastable_phases() to force the calculation to converge for the metastable phases.

**Returns** This PropertyDiagramOptions object

enable_control_step_size_during_minimization()  
Enables stepsize control during normal minimization (non-global).  

**Default:** Enabled

**Returns** This PropertyDiagramOptions object

enable_force_positive_definite_phase_hessian()  
Enables forcing of positive definite phase Hessian. This determines how the minimum of an equilibrium state in a normal minimization procedure (non-global) is reached. For details, search the Thermo-Calc documentation for “Hessian minimization”.  

**Default:** Enabled
Returns This `PropertyDiagramOptions` object

**set_global_minimization_max_grid_points** *(max_grid_points: int = 2000)*
Sets the maximum number of grid points in global minimization. Only applicable if global minimization is actually used.

Default: 2000 points

Parameters **max_grid_points** – The maximum number of grid points

Returns This `PropertyDiagramOptions` object

**set_global_minimization_test_interval** *(global_test_interval: int = 0)*
Sets the interval for the global test.

Default: 0

Parameters **global_test_interval** – The global test interval

Returns This `PropertyDiagramOptions` object

**set_max_no_of_iterations** *(max_no_of_iterations: int = 500)*
Set the maximum number of iterations.

Default: max. 500 iterations

Note: As some models give computation times of more than 1 CPU second/iteration, this number is also used to check the CPU time and the calculation stops if 500 CPU seconds/iterations are used.

Parameters **max_no_of_iterations** – The max. number of iterations

Returns This `PropertyDiagramOptions` object

**set_required_accuracy** *(accuracy: float = 1e-06)*
Sets the required relative accuracy.

Default: 1.0E-6

Note: This is a relative accuracy, and the program requires that the relative difference in each variable must be lower than this value before it has converged. A larger value normally means fewer iterations but less accurate solutions. The value should be at least one order of magnitude larger than the machine precision.

Parameters **accuracy** – The required relative accuracy

Returns This `PropertyDiagramOptions` object

**set_smallest_fraction** *(smallest_fraction: float = 1e-12)*
Sets the smallest fraction for constituents that are unstable.

It is normally only in the gas phase that you can find such low fractions.

The default value for the smallest site-fractions is 1E-12 for all phases except for IDEAL phase with one sublattice site (such as the GAS mixture phase in many databases) for which the default value is always as 1E-30.

Parameters **smallest_fraction** – The smallest fraction for constituents that are unstable

Returns This `PropertyDiagramOptions` object
**PropertyDiagramResult**

Result of a property diagram. This can be used to query for specific values.

### Methods

**get_values_grouped_by_quantity_of**

```python
get_values_grouped_by_quantity_of(x_quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], y_quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], sort_and_merge: bool = True) -> Dict[str, tc_python.utils.ResultValueGroup]
```

Returns x-y-line data grouped by the multiple datasets of the specified quantities (typically the phases). The available quantities can be found in the documentation of the factory class `ThermodynamicQuantity`.

**get_values_grouped_by_stable_phases_of**

```python
get_values_grouped_by_stable_phases_of(x_quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], y_quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], sort_and_merge: bool = True) -> Dict[str, tc_python.utils.ResultValueGroup]
```

Returns x-y-line data grouped by the sets of “stable phases” (for example “LIQUID” or “LIQUID + FCC_A1”). The available quantities can be found in the documentation of the factory class `ThermodynamicQuantity`.

### Parameters

- **x_quantity** – The first quantity (“x-axis”), Console Mode syntax strings can be used as an alternative (for example ‘T’), or even a function (for example ‘f=T*1.01’)
- **y_quantity** – The second quantity (“y-axis”), Console Mode syntax strings can be used as an alternative (for example ‘NV’), or even a function (for example ‘CP=HM.T’)
- **sort_and_merge** – If True, the data is sorted and merged into as few subsections as possible (divided by NaN)

### Returns

Containing the datasets with the quantities as their keys.

### Notes

- The different datasets might contain NaN-values between different subsections and might not be sorted even if the flag `sort_and_merge` has been set (because they might be unsortable due to their nature).
- It’s possible to use functions as axis variables, either by using `ThermodynamicQuantity.user_defined_function`, or by using an expression that contains ‘=’.
Example

\[ \text{get\_values\_grouped\_by\_quantity\_of}(\ 'T', \ \text{ThermodynamicQuantity}().\text{user\_defined\_function}(\ 'HM.T')) \]

Example

\[ \text{get\_values\_grouped\_by\_quantity\_of}(\ 'T', \ 'CP=HM.T') \]

Parameters

- **\textit{x\_quantity}** – The first quantity ("x-axis"), Console Mode syntax strings can be used as an alternative (for example ‘T’), or even a function (for example ‘f=T*1.01’)

- **\textit{y\_quantity}** – The second quantity ("y-axis"), Console Mode syntax strings can be used as an alternative (for example ‘NV’), or even a function (for example ‘CP=HM.T’)

- **\textit{sort\_and\_merge}** – If True, the data will be sorted and merged into as few subsections as possible (divided by NaN)

Returns

Containing the datasets with the quantities as their keys

\[ \text{get\_values\_of}(x\_quantity: \ \text{Union}[\text{tc\_python.quantity\_factory.ThermodynamicQuantity}, \ \text{str}], \ y\_quantity: \ \text{Union}[\text{tc\_python.quantity\_factory.ThermodynamicQuantity}, \ \text{str}]) \rightarrow \ \text{[typing.List[float], typing.List[float]]} \]

Returns sorted x-y-line data without any separation. Use \text{get\_values\_grouped\_by\_quantity\_of()} or \text{get\_values\_grouped\_by\_stable\_phases\_of()} instead if you need such a separation. The available quantities can be found in the documentation of the factory class ThermodynamicQuantity.

Note: This method will always return sorted data without any NaN-values. If it is unsortable that might give data that is hard to interpret. In such a case you need to choose the quantity in another way or use one of the other methods. One example of this is to use quantities with *All*-markers, for example MassFractionOfAComponent("All").

Note: It is possible to use functions as axis variables, either by using \text{ThermodynamicQuantity}.user_defined_function, or by using an expression that contains ‘\textquoteleft;=\textquoteright;”.

Example

\[ \text{get\_values\_grouped\_by\_quantity\_of}(\ 'T', \ \text{ThermodynamicQuantity}().\text{user\_defined\_function}(\ 'HM.T')) \]

Example

\[ \text{get\_values\_grouped\_by\_quantity\_of}(\ 'T', \ 'CP=HM.T') \]

Parameters

- **\textit{x\_quantity}** – The first Thermodynamic quantity (“x-axis”), Console Mode syntax strings can be used as an alternative (for example ‘T’) or even a function (for example ‘f=T*1.01’)

- **\textit{y\_quantity}** – The second Thermodynamic quantity (“y-axis”), Console Mode syntax strings can be used as an alternative (for example ‘NV’), or even a function (for example ‘\text{CP=HM.T}’)

Returns

A tuple containing the x- and y-data in lists

\[ \text{save\_to\_disk}(\ \text{path: str}) \]

Saves the result to disc. Note that a result is a folder, containing potentially many files. The result can later be loaded with \text{load\_result\_from\_disk()}

Parameters

- **\textit{path}** – the path to the folder you want the result to be saved in. It can be relative or absolute.
Returns this `PropertyDiagramResult` object

**set_phase_name_style** (phase_name_style_enum: tc_python.step_or_map_diagrams.PhaseNameStyle = <PhaseNameStyle.NONE: 0>)

Sets the style of the phase name labels that will be used in the result data object (constitution description, ordering description, ...).

Default: PhaseNameStyle.NONE

Parameters **phase_name_style_enum** – The phase name style

Returns This `PropertyDiagramResult` object

### 5.1.6 Module “diffusion”

**class** tc_python.diffusion.AbstractBoundaryCondition

**Bases:** object

The abstract base class for all boundary conditions.

**class** tc_python.diffusion.AbstractCalculatedGrid

**Bases:** tc_python.diffusion.AbstractGrid

The abstract base class for calculated grids.

**class** tc_python.diffusion.AbstractElementProfile

**Bases:** object

The abstract base class for all initial composition profile types.

**class** tc_python.diffusion.AbstractGrid

**Bases:** object

The abstract base class for all grids.

**class** tc_python.diffusion.AbstractSolver

**Bases:** object

Abstract base class for the solvers (Classic, Homogenization and Automatic).

**class** tc_python.diffusion.ActivityFluxFunction

**Bases:** tc_python.diffusion.BoundaryCondition

**get_type** () → str

The type of the boundary condition.

Returns The type

**set_flux_function** (element_name: str, f: str = '0', g: str = '1', n: float = 1.0, to_time: float = 1.7976931348623157e+308)

The flux for the independent components must be given in the format:

\[ J = f(T,P,TIME) \cdot (ACTIVITY^N - g(T,P,TIME)) \]

where \( f \) and \( g \) may be functions of time (TIME), temperature (T), and pressure (P), and \( N \) is an integer.

\( f \) and \( g \) must be expressed in DICTRA Console Mode syntax.

Parameters
- **element_name** – The name of the element
- **f** – the function \( f \) in the formula above
- **g** – the function \( g \) in the formula above
• \( n \) – the constant \( N \) in the formula above

• \texttt{to\_time} – The max-time for which the flux function is used.

```python
class tc_python.diffusion.AutomaticSolver
Bases: tc_python.diffusion.Solver

Solver using the \textit{homogenization model} if any region has more than one phase, otherwise using the \textit{classic model}.
```

**Note:** This is the \texttt{default solver} and recommended for most applications.

```python
def get_type()
    return str

    Returns
    The type of the solver.

def set_flux_balance_equation_accuracy(accuracy: float = 1e-16)

    Only valid if the :class:`ClassicSolver` is actually used (i.e. not more than one phase in each region).

    Sets the required accuracy during the solution of the flux balance equations. \texttt{Default}: 1.0e-16

    Parameters
    \texttt{accuracy} -- The required accuracy

    Returns
    A new \texttt{AutomaticSolver} object

def set_tieline_search_variable_to_activity()

    Only valid if the :class:`ClassicSolver` is actually used (i.e. not more than one phase in each region).

    Configures the solver to use the \textit{activity of a component} to find the correct tie-line at the phase interface. Either activity or chemical potential are applied to reduce the degrees of freedom at the local equilibrium. \texttt{Default}: This is the default setting

    Returns
    A new \texttt{AutomaticSolver} object

def set_tieline_search_variable_to_potential()

    Only valid if the :class:`ClassicSolver` is actually used (i.e. not more than one phase in each region).

    Configures the solver to use the \textit{chemical potential of a component} to find the correct tie-line at the phase interface. Either activity or chemical potential are applied to reduce the degrees of freedom at the local equilibrium. \texttt{Default}: To use the activity

    Returns
    A new \texttt{AutomaticSolver} object
```

```python
class tc_python.diffusion.BoundaryCondition
Bases: tc_python.diffusion.AbstractBoundaryCondition

Contains factory methods for the different boundary conditions available.

```python
def activity_flux_function()

    Factory method that creates a \texttt{new} activity-flux-function boundary condition.

    This type of boundary condition is used to take into account the finite rate of a surface reaction.

    The flux for the independent components must be given in the format:

    \( J = f(T,P,\text{TIME}) \times (\text{ACTIVITY}^N - g(T,P,\text{TIME})) \)

    where \( f \) and \( g \) may be functions of time (\text{TIME}), temperature (\text{T}), and pressure (\text{P}), and \( N \) is an integer. \( f \) and \( g \) must be expressed in DICTRA Console Mode syntax.
```
Note: The activities are those with user-defined reference states. The function mass transfer coefficient is the mass transfer coefficient, activity of the corresponding species in the gas is the activity of the corresponding species in the gas and N is a stoichiometric coefficient.


Returns A new ActivityFluxFunction object

classmethod closed_system() Factory method that creates a new closed-system boundary condition.

Returns A new ClosedSystem object

classmethod fix_flux_value() Factory method that creates a new fix-flux-value boundary condition.

This type of boundary condition makes it possible to enter functions that yield the flux times the molar volume for the independent components. May be a function of time, temperature and pressure: J(T,P,TIME).

Returns A new Fix Flux Value object

classmethod fixed_compositions(unit_enum: tc_python.diffusion.Unit = <Unit.MASS_PERCENT: 3>) Factory method that creates a new fixed-composition boundary condition.

Parameters unit_enum – The composition unit

Returns A new Fixed Compositions object

classmethod mixed_zero_flux_and_activity() Factory method that creates a new mixed zero-flux and activity boundary condition

Returns A new Mixed Zero Flux And Activity object

class tc_python.diffusion.CalculatedGrid Bases: tc_python.diffusion.AbstractCalculatedGrid

Factory class for grids generated by a mathematical series (linear, geometric, ...). Use tc_python.diffusion.PointByPointGrid instead if you want to use an existing grid from experimental data or a previous calculation.

Note: A region must contain a number of grid points. The composition is only known at these grid points and the software assumes that the composition varies linearly between them. The amount and composition of all the phases present at a single grid point in a certain region are those given by thermodynamic equilibrium keeping the over-all composition at the grid point fixed.

classmethod double_geometric(no_of_points: int = 50, lower_geometrical_factor: float = 1.1, upper_geometrical_factor: float = 0.9) Factory method that creates a new double geometric grid.

Note: Double geometric grids have a high number of grid points in the middle or at both ends of a region. One geometrical factor for the lower (left) and upper (right) half of the region need to specified. In both
cases a geometrical factor of larger than one yields a higher density of grid points at the lower end of the half and vice versa for a factor smaller than one.

Parameters

- **no_of_points** – The number of points
- **lower_geometrical_factor** – The geometrical factor for the left half
- **upper_geometrical_factor** – The geometrical factor for the right half

Returns A new DoubleGeometricGrid object

classmethod geometric(**no_of_points**: int = 50, **geometrical_factor**: float = 1.1)

Factory method that creates a new geometric grid.

Note: A grid that yields a varying density of grid points in the region. A geometrical factor larger than one yields a higher density of grid points at the lower end of the region and a factor smaller than one yields a higher density of grid points at the upper end of the region.

Parameters

- **no_of_points** – The number of points
- **geometrical_factor** – The geometrical factor

Returns A new GeometricGrid object

classmethod linear(**no_of_points**: int = 50)

Factory method that creates a new equally spaced grid.

Parameters **no_of_points** – The number of points

Returns A new LinearGrid object

class tc_python.diffusion.ClassicSolver

Bases: tc_python.diffusion.Solver

Solver using the Classic model.

Note: This solver never switches to the homogenization model even if it fails to converge. Use the tc_python.diffusion.AutomaticSolver if necessary instead.

get_type() → str

Convenience method for getting the type of the solver.

Returns The type of the solver

set_flux_balance_equation_accuracy(**accuracy**: float = 1e-16)

Sets the required accuracy during the solution of the flux balance equations. Default: 1.0e-16

Parameters **accuracy** – The required accuracy

Returns A new ClassicSolver object

set_tieline_search_variable_to_activity()

Configures the solver to use the activity of a component to find the correct tie-line at the phase interface.

5.1. Calculations

115
Either activity or chemical potential are applied to reduce the degrees of freedom at the local equilibrium.

**Default:** This is the default setting

```python
set_tieline_search_variable_to_potential()
```

Configures the solver to use the chemical potential of a component to find the correct tie-line at the phase interface. Either activity or chemical potential are applied to reduce the degrees of freedom at the local equilibrium. **Default:** To use the activity

**Returns** A new `ClassicSolver` object

```python
class tc_python.diffusion.ClosedSystem
    Bases: tc_python.diffusion.BoundaryCondition
```

Represents a boundary for a closed system.

```python
get_type() -> str
```

Convenience method for getting the type of the boundary condition.

**Returns** The type of the boundary condition

```python
class tc_python.diffusion.CompositionProfile(unit_enum: tc_python.diffusion.Unit =<Unit.MASS_PERCENT: 3>)
```

Bases: object

Contains initial concentration profiles for the elements.

```python
add(element_name: str, profile: tc_python.diffusion.ElementProfile)
```

Adds a concentration profile for the specified element.

**Parameters**

- `element_name` – The name of the element
- `profile` – The initial concentration profile

**Returns** A `CompositionProfile` object

```python
class tc_python.diffusion.ConstantProfile(value: float)
```

Bases: `tc_python.diffusion.ElementProfile`

Represents a constant initial concentration profile.

```python
get_type() -> str
```

The type of the element profile.

**Returns** The type of the element profile

```python
class tc_python.diffusion.ContinuedDiffusionCalculation(calculation)
```

Bases: `tc_python.abstract_base.AbstractCalculation`

Configuration for a diffusion calculation that is a continuation of a previous isothermal or non-isothermal diffusion calculation. It contains a subset of the settings possible in the original calculation.

Use `set_simulation_time()` to set a simulation time that is higher than the original calculation.

```python
calculate() -> tc_python.diffusion.DiffusionCalculationResult
```

Runs the diffusion calculation.

**Returns** A `DiffusionCalculationResult` which later can be used to get specific values from the calculated result

```python
set_simulation_time(simulation_time: float)
```

Sets the simulation time.

**Parameters** `simulation_time` – The simulation time [s]

**Returns** This `DiffusionIsoThermalCalculation` object
with_left_boundary_condition(boundary_condition: tc_python.diffusion.BoundaryCondition,
    to: float = 1.7976931348623157e+308)

Defines the boundary condition on the left edge of the system.

Default: A closed-system boundary condition.

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

Default: The end of the simulation.

Note: You can specify time-dependent boundary conditions by calling with_left_boundary_condition() many times, with different values of the “to” parameter.

Examples:

• with_left_boundary_condition(BoundaryCondition.closed_system(), to=100)
• with_left_boundary_condition(BoundaryCondition.mixed_zero_flux_and_activity().set_activity_for_element("C", surface_activity), to=500)
• with_left_boundary_condition(BoundaryCondition.closed_system())

This example sets an closed-system-boundary-condition from start up to 100s and a activity-boundary-condition from 100s to 500s and finally a closed-system-boundary-condition from 500s to the end of simulation.

Parameters

• boundary_condition – The boundary condition
• to – The upper time-limit for boundary_condition.

Returns This DiffusionIsoThermalCalculation object

with_options(options: tc_python.diffusion.Options, to: float = 1.7976931348623157e+308)

Sets the general simulation conditions.

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

Default: The end of the simulation.

Parameters

• options – The general simulation conditions
• to – The upper time-limit for options.

Returns This DiffusionIsoThermalCalculation object

with_right_boundary_condition(boundary_condition: tc_python.diffusion.BoundaryCondition,
    to: float = 1.7976931348623157e+308)

Defines the boundary condition on the right edge of the system.

Default: A closed-system boundary condition

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

Default: The end of the simulation.

Note: You can specify time-dependent boundary conditions by calling with_right_boundary_condition() many times, with different values of the “to” parameter.
Examples:

- `with_right_boundary_condition(BoundaryCondition.closed_system(), to=100)`
- `with_right_boundary_condition(BoundaryCondition.mixed_zero_flux_and_activity().set_activity_for_element("C", surface_activity), to=500)`
- `with_right_boundary_condition(BoundaryCondition.closed_system())`

This example sets an closed-system-boundary-condition from start up to 100s and a activity-boundary-condition from 100s to 500s and finally a closed-system-boundary-condition from 500s to the end of simulation.

**Parameters**

- **boundary_condition** – The boundary condition
- **to** – The upper time-limit for boundary_condition.

**Returns** This `DiffusionIsoThermalCalculation` object

**with_solver** (`solver: tc_python.diffusion.Solver, to: float = 1.7976931348623157e+308`)

Sets the solver to use (Classic, Homogenization or Automatic). **Default is Automatic.**

**Default:** The end of the simulation.

**Parameters**

- **solver** – The solver to use
- **to** – The upper time-limit for solver.

**Returns** This `DiffusionIsoThermalCalculation` object

**with_timestep_control** (`timestep_control: tc_python.diffusion.TimestepControl, to: float = 1.7976931348623157e+308`)

Sets the timestep control options.

**Default:** The end of the simulation.

**Parameters**

- **timestep_control** – The new timestep control options
- **to** – The upper time-limit for timestep_control.

**Returns** This `DiffusionIsoThermalCalculation` object

**class** `tc_python.diffusion.DiffusionCalculationResult`(`result`)

**Bases:** `tc_python.abstract_base.AbstractResult`

Result of a diffusion calculation. This can be used to query for specific values. For details of the axis variables, search the Thermo-Calc help.

**get_mass_fraction_at_lower_interface** (`region: str, component: str`) → `[typing.List[float], typing.List[float]]`

Returns the mass fraction of the specified component at the lower boundary of the specified region, in dependency of time.

**Parameters**

- **region** – The name of the region
- **component** – The name of the component
Returns A tuple of two lists of floats (time [s], mass fraction of the specified component)

```python
def get_mass_fraction_at_upper_interface(region: str, component: str) -> [typing.List[float], typing.List[float]]
```

Returns the mass fraction of the specified component at the upper boundary of the specified region, in dependency of time.

Parameters

- `region` – The name of the region
- `component` – The name of the component

Returns A tuple of two lists of floats (time [s], mass fraction of the specified component)

```python
def get_mass_fraction_of_component_at_time(component: str, time: Union[tc_python.diffusion.SimulationTime, float]) -> [typing.List[float], typing.List[float]]
```

Returns the mass fraction of the specified component at the specified time.

Note: Use the enum `tc_python.diffusion.SimulationTime` to choose the first or the last timepoint of the simulation. A timepoint close to the last one should never be specified manually because the actual end of the simulation can slightly deviate.

Parameters

- `component` – The name of the component
- `time` – The time [s]

Returns A tuple of two lists of floats (distance [m], mass fraction of component at the specified time)

```python
def get_mass_fraction_of_phase_at_time(phase: str, time: Union[tc_python.diffusion.SimulationTime, float]) -> [typing.List[float], typing.List[float]]
```

Returns the mass fraction of the specified phase.

Note: Use the enum `tc_python.diffusion.SimulationTime` to choose the first or the last timepoint of the simulation. A timepoint close to the last one should never be specified manually because the actual end of the simulation can slightly deviate.

Parameters

- `phase` – The name of the phase
- `time` – The time [s]

Returns A tuple of two lists of floats (distance [m], mass fraction of the phase at the specified time)

```python
def get_mole_fraction_at_lower_interface(region: str, component: str) -> [typing.List[float], typing.List[float]]
```

Returns the mole fraction of the specified component at the lower boundary of the specified region, in dependency of time.

Parameters

5.1. Calculations
• **region** – The name of the region
• **component** – The name of the component

Returns A tuple of two lists of floats (time [s], mole fraction of the specified component)

```python
def get_mole_fraction_at_upper_interface(region: str, component: str) -> [typing.List[float], typing.List[float]]
```

Returns the mole fraction of the specified component at the upper boundary of the specified region, in dependency of time.

Parameters

• **region** – The name of the region
• **component** – The name of the component

Returns A tuple of two lists of floats (time [s], mole fraction of the specified component)

```python
def get_mole_fraction_of_component_at_time(component: str, time: Union[tc_python.diffusion.SimulationTime, float]) -> [typing.List[float], typing.List[float]]
```

Returns the mole fraction of the specified component at the specified time.

Note: Use the enum `tc_python.diffusion.SimulationTime` to choose the first or the last timepoint of the simulation. A timepoint close to the last one should never be specified manually because the actual end of the simulation can slightly deviate.

Parameters

• **component** – The name of the component

• **time** – The time [s]

Returns A tuple of two lists of floats (distance [m], mole fraction of component at the specified time)

```python
def get_mole_fraction_of_phase_at_time(phase: str, time: Union[tc_python.diffusion.SimulationTime, float]) -> [typing.List[float], typing.List[float]]
```

Returns the mole fraction of the specified phase.

Note: Use the enum `tc_python.diffusion.SimulationTime` to choose the first or the last timepoint of the simulation. A timepoint close to the last one should never be specified manually because the actual end of the simulation can slightly deviate.

Parameters

• **phase** – The name of the phase

• **time** – The time [s]

Returns A tuple of two lists of floats (distance [m], mole fraction of the phase at the specified time)

```python
def get_position_of_lower_boundary_of_region(region: str) -> [typing.List[float], typing.List[float]]
```

Returns the position of the lower boundary of the specified region in dependency of time.
Parameters `region` – The name of the region

Returns A tuple of two lists of floats (time [s], position of lower boundary of region [m])

`get_position_of_upper_boundary_of_region(region: str) → [typing.List[float], typing.List[float]]`

Returns the position of the upper boundary of the specified region in dependency of time.

Parameters `region` – The name of the region

Returns A tuple of two lists of floats (time [s], position of upper boundary of region [m])

`get_regions() → List[str]`

Returns the regions of the diffusion simulation.

Note: Automatically generated regions \( R_{###} \) are included in the list.

Returns The region names

`get_time_steps() → List[float]`

Returns the timesteps of the diffusion simulation.

Returns The timesteps [s]

`get_total_mass_fraction_of_component(component: str) → [typing.List[float], typing.List[float]]`

Returns the total mass fraction of the specified component in dependency of time.

Parameters `component` – The name of the component

Returns A tuple of two lists of floats (time [s], total mass fraction of the component)

`get_total_mass_fraction_of_component_in_phase(component: str, phase: str) → [typing.List[float], typing.List[float]]`

Returns the total mass fraction of the specified component in the specified phase in dependency of time.

Parameters

- `component` – The name of the component
- `phase` – The name of the phase

Returns A tuple of two lists of floats (time [s], total mass fraction of the component in the phase)

`get_total_mass_fraction_of_phase(phase: str) → [typing.List[float], typing.List[float]]`

Returns the total mass fraction of the specified phase in dependency of the time.

Parameters `phase` – The name of the phase

Returns A tuple of two lists of floats (time [s], total mass fraction of the phase)

`get_total_mole_fraction_of_component(component: str) → [typing.List[float], typing.List[float]]`

Returns the total mole fraction of the specified component in dependency of time.

Parameters `component` – The name of the component

Returns A tuple of two lists of floats (time [s], total mole fraction of the component)

`get_total_mole_fraction_of_component_in_phase(component: str, phase: str) → [typing.List[float], typing.List[float]]`

Returns the total mole fraction of the specified component in the specified phase in dependency of time.

Parameters
• \texttt{component} – The name of the component

• \texttt{phase} – The name of the phase

Returns A tuple of two lists of floats (time [s], total mole fraction of the component in the phase)

\texttt{get\_total\_mole\_fraction\_of\_phase (phase: str) \rightarrow [typing.List[float], typing.List[float]]}

Returns the total mole fraction of the specified phase in dependency of time.

Parameters \texttt{phase} – The name of the phase

Returns A tuple of two lists of floats (time [s], total mole fraction of the phase)

\texttt{get\_total\_volume\_fraction\_of\_phase (phase: str) \rightarrow [typing.List[float], typing.List[float]]}

Returns the total volume fraction of the specified phase in dependency of the time.

Parameters \texttt{phase} – The name of the phase

Returns A tuple of two lists of floats (time [s], total volume fraction of the phase)


Returns the specified result from the simulation, allows all possible settings.

Note: As an alternative, DICTRA Console Mode syntax can be used as well for each quantity and condition.

Warning: This is an advanced mode that is equivalent to the possibilities in the DICTRA Console Mode. Not every combination of settings will return a result.

Parameters

• \texttt{x\_axis} – The first result quantity

• \texttt{y\_axis} – The second result quantity

• \texttt{plot\_condition} – The plot conditions

• \texttt{independent\_variable} – The independent variable

Returns A tuple of two lists of floats (the \texttt{x\_axis} quantity result, the \texttt{y\_axis} quantity result) [units according to the quantities]

\texttt{get\_velocity\_of\_lower\_boundary\_of\_region (region: str) \rightarrow [typing.List[float], typing.List[float]]}

Returns the velocity of the lower boundary of the specified region in dependency of time.

Parameters \texttt{region} – The name of the region

Returns A tuple of two lists of floats (time [s], velocity of lower boundary of region [m/s])

\texttt{get\_velocity\_of\_upper\_boundary\_of\_region (region: str) \rightarrow [typing.List[float], typing.List[float]]}

Returns the velocity of the upper boundary of the specified region in dependency of time.

Parameters \texttt{region} – The name of the region

Returns A tuple of two lists of floats (time [s], velocity of upper boundary of region [m/s])
**get_width_of_region** *(region: str) → [typing.List[float], typing.List[float]]*

Returns the width of region, in dependency of time.

**Parameters**
- **region** – The name of the region

**Returns**
- A tuple of two lists of floats (time [s], width of the specified region [m])

**save_to_disk** *(path: str)*

Saves the result to disk. The result can later be loaded using `tc_python.server.SetUp.load_result_from_disk()`.

**Note:** The result data is represented by a whole folder containing multiple files.

**Parameters**
- **path** – The path to the result folder, can be relative or absolute.

**Returns**
- This `DiffusionCalculationResult` object

**with_continued_calculation** ()

Returns a `ContinuedDiffusionCalculation` that is used for continuing a diffusion calculation with altered settings.

**Returns**
- A `ContinuedDiffusionCalculation`

**class** `tc_python.diffusion.DiffusionIsoThermalCalculation` *(calculation)*

**Bases:** `tc_python.abstract_base.AbstractCalculation`

Configuration for an isothermal diffusion calculation.

**add_console_command** *(console_command: str)*

Registers a DICTRA Console Mode command for execution. These commands are executed after all other configuration directly before the calculation starts to run. All commands are stored and used until explicitly deleted using `tc_python.diffusion.DiffusionIsoThermoCalculation.remove_all_console_commands()`.

**Parameters**
- **console_command** – The DICTRA Console Mode command

**Returns**
- This `DiffusionIsoThermalCalculation` object

**Note:** It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

**Warning:** As this method runs raw DICTRA-commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten parenthesis, …).

**add_region** *(region: tc_python.diffusion.Region)*

Adds a region to the calculation. Regions are always added in the simulation domain from left to right.

If you want to replace an already added region, call `remove_all_regions()`, and add the regions that you want to keep.

**Warning:** Regions must have unique names.

**Parameters**
- **region** – The region to be added
Returns This `DiffusionIsoThermalCalculation` object

calculate() \(\rightarrow tc\_python.diffusion.DiffusionCalculationResult\)
Runs the diffusion calculation.

Returns A `DiffusionCalculationResult` which later can be used to get specific values from the calculated result

get_system_data() \(\rightarrow tc\_python.abstract\_base.SystemData\)
Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using `with_system_modifications()`.

Note: Parameters can only be read from unencrypted (i.e. *user*) databases loaded as *.tdb*-file.

Returns The system data

remove_all_console_commands()
Removes all previously added Console Mode commands.

Returns This `DiffusionIsoThermalCalculation` object

remove_all_regions()
Removes all previously added regions.

:returns This `DiffusionIsoThermalCalculation` object

set_simulation_time(simulation_time: float)
Sets the simulation time.

Parameters simulation_time – The simulation time [s]

Returns This `DiffusionIsoThermalCalculation` object

set_temperature(temperature: float)
Sets the temperature for the isothermal simulation.

Parameters temperature – The temperature [K]

Returns This `DiffusionIsoThermalCalculation` object

with_cylindrical_geometry(first_interface_position: float = 0.0)
Sets geometry to `cylindrical`, corresponds to an infinitely long cylinder of a certain radius.

Default: A planar geometry

Note: With a cylindrical or spherical geometry, the system’s zero coordinate (left boundary) is at the centre of the cylinder or sphere by default. By specifying the `first_interface_position`, a different left-most coordinate can be defined. This allows to model a tube or a hollow sphere geometry. The highest coordinate (right boundary) is defined by the cylinder or sphere radius (i.e. by the width of all regions).

Parameters first_interface_position – The position of the left-most coordinate along the axis, only necessary for modeling a tube geometry [m]

Returns This `DiffusionIsoThermalCalculation` object

with_left_boundary_condition(boundary_condition: tc_python.diffusion.BoundaryCondition, to: float = 1.7976931348623157e+308)
Defines the boundary condition on the left edge of the system.
**Default:** A closed-system boundary condition.

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

**Default:** The end of the simulation.

---

**Note:** You can specify time-dependent boundary conditions by calling `with_left_boundary_condition()` many times, with different values of the “to” parameter.

---

Examples:

- `with_left_boundary_condition(BoundaryCondition.closed_system(), to=100)`
- `with_left_boundary_condition(BoundaryCondition.mixed_zero_flux_and_activity().set_activity_for_element("C", surface_activity), to=500)`
- `with_left_boundary_condition(BoundaryCondition.closed_system())`

This example sets an closed-system-boundary-condition from start up to 100s and a activity-boundary-condition from 100s to 500s and finally a closed-system-boundary-condition from 500s to the end of simulation.

**Parameters**

- `boundary_condition` – The boundary condition
- `to` – The upper time-limit for boundary_condition.

**Returns** This `DiffusionIsoThermalCalculation` object

with_options *(options: tc_python.diffusion.Options, to: float = 1.7976931348623157e+308)*

Sets the general simulation conditions.

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

**Default:** The end of the simulation.

**Parameters**

- `options` – The general simulation conditions
- `to` – The upper time-limit for options.

**Returns** This `DiffusionIsoThermalCalculation` object

with_planar_geometry()

Sets geometry to `planar`.

This is default.

**Returns** This `DiffusionIsoThermalCalculation` object

with_reference_state *(element: str, phase: str = 'SER', temperature: float = - 1.0, pressure: float = 100000.0)*

The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.

All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to
set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

**Parameters**

- **element** – The name of the element
- **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.
- **temperature** – The Temperature (in K) for the reference state. Or CURRENT_TEMPERATURE which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.
- **pressure** – The pressure (in Pa) for the reference state

**Returns** This DiffusionIsoThermalCalculation object

```python
with_right_boundary_condition(boundary_condition: tc_python.diffusion.BoundaryCondition, to: float = 1.7976931348623157e+308)
```

Defines the boundary condition on the right edge of the system.

**Default:** A closed-system boundary condition

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

**Default:** The end of the simulation.

**Note:** You can specify time-dependent boundary conditions by calling with_right_boundary_condition() many times, with different values of the “to” parameter.

**Examples:**

- `with_right_boundary_condition(BoundaryCondition.closed_system(), to=100)`
- `with_right_boundary_condition(BoundaryCondition.mixed_zero_flux_and_activity().set_activity_for_element("C", surface_activity), to=500)`
- `with_right_boundary_condition(BoundaryCondition.closed_system())`

This example sets an closed-system-boundary-condition from start up to 100s and a activity-boundary-condition from 100s to 500s and finally a closed-system-boundary-condition from 500s to the end of simulation.

**Parameters**

- **boundary_condition** – The boundary condition
- **to** – The upper time-limit for boundary_condition.

**Returns** This DiffusionIsoThermalCalculation object

```python
with_solver(solver: tc_python.diffusion.Solver, to: float = 1.7976931348623157e+308)
```

Sets the solver to use (Classic, Homogenization or Automatic). **Default is Automatic.**
It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

**Default:** The end of the simulation.

**Parameters**

- `solver` – The solver to use
- `to` – The upper time-limit for solver.

**Returns** This `DiffusionIsoThermalCalculation` object

`with_spherical_geometry(first_interface_position: float = 0.0)`

Sets geometry to *spherical*, corresponds to a sphere with a certain radius.

**Default:** A spherical geometry

**Note:** With a cylindrical or spherical geometry, the system’s zero coordinate (left boundary) is at the centre of the cylinder or sphere by default. By specifying the `first_interface_position`, a different left-most coordinate can be defined. This allows to model a tube or a hollow sphere geometry. The highest coordinate (right boundary) is defined by the cylinder or sphere radius (i.e. by the width of all regions).

**Parameters** `first_interface_position` – The position of the left-most coordinate along the axis, only necessary for modeling a hollow sphere geometry [m]

**Returns** This `DiffusionIsoThermalCalculation` object

`with_system_modifications(system_modifications: tc_python.abstract_base.SystemModifications)`

Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

**Note:** This is only possible if the system has been read from unencrypted (i.e. user) databases loaded as a *.tdb*-file.

**Parameters** `system_modifications` – The system modification to be performed

**Returns** This `DiffusionIsoThermalCalculation` object

`with_timestep_control(timestep_control: tc_python.diffusion.TimestepControl, to: float = 1.7976931348623157e+308)`

Sets the timestep control options.

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

**Default:** The end of the simulation.

**Parameters**

- `timestep_control` – The new timestep control options
- `to` – The upper time-limit for timestep_control.

**Returns** This `DiffusionIsoThermalCalculation` object

```python
class tc_python.diffusion.DiffusionNonIsoThermalCalculation(calculation)
Bases: tc_python.abstract_base.AbstractCalculation
```

Configuration for a non-isothermal diffusion calculation.
**add_console_command** *(console_command: str)*

Registers a DICTRA Console Mode command for execution. These commands are executed after all other configuration directly before the calculation starts to run. All commands are stored and used until explicitly deleted using `tc_python.diffusion.DiffusionNonIsoThermalCalculation.remove_all_console_commands`.

**Parameters**

- **console_command** – The DICTRA Console Mode command

**Returns**

This `DiffusionNonIsoThermalCalculation` object

**Note:** It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

**Warning:** As this method runs raw DICTRA-commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten parenthesis, ...).

**add_region** *(region: tc_python.diffusion.Region)*

Adds a region to the calculation. Regions are always added in the simulation domain from left to right. If you want to replace an already added region, call `remove_all_regions()`, and add the regions that you want to keep.

**Warning:** Regions must have unique names.

**Parameters**

- **region** – The region to be added

**Returns**

This `DiffusionNonIsoThermalCalculation` object

**calculate() → tc_python.diffusion.DiffusionCalculationResult**

Runs the diffusion calculation.

**Returns**

A `DiffusionCalculationResult` which later can be used to get specific values from the calculated result

**get_system_data() → tc_python.abstract_base.SystemData**

Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using `with_system_modifications()`.

**Note:** Parameters can only be read from unencrypted (i.e. `user`) databases loaded as *.tdb*-file.

**Returns**

The system data

**remove_all_console_commands()**

Removes all previously added Console Mode commands.

**Returns**

This `DiffusionNonIsoThermalCalculation` object

**remove_all_regions()**

Removes all previously added regions.

**Returns**

This `DiffusionNonIsoThermalCalculation` object
**set_simulation_time** *(simulation_time: float)*

Sets the simulation time.

**Parameters**  
**simulation_time** – The simulation time [s]

**Returns**  
This `DiffusionNonIsoThermalCalculation` object

**with_cylindrical_geometry** *(first_interface_position: float = 0.0)*

Sets geometry to cylindrical, corresponds to an infinitely long cylinder of a certain radius.

**Default**: A planar geometry

**Note:** With a cylindrical or spherical geometry, the system’s zero coordinate (left boundary) is at the centre of the cylinder or sphere by default. By specifying the `first_interface_position`, a different left-most coordinate can be defined. This allows to model a tube or a hollow sphere geometry. The highest coordinate (right boundary) is defined by the cylinder or sphere radius (i.e. by the width of all regions).

**Parameters**  
**first_interface_position** – The position of the left-most coordinate along the axis, only necessary for modeling a tube geometry [m]

**Returns**  
This `DiffusionNonIsoThermalCalculation` object

**with_left_boundary_condition** *(boundary_condition: tc_python.diffusion.BoundaryCondition, to: float = 1.7976931348623157e+308)*

Defines the boundary condition on the left edge of the system.

**Default**: A closed-system boundary condition.

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

**Default**: The end of the simulation.

**Note:** You can specify time-dependent boundary conditions by calling `with_left_boundary_condition()` many times, with different values of the “to” parameter.

**Examples:**

- `with_left_boundary_condition(BoundaryCondition.closed_system(), to=100)`
- `with_left_boundary_condition(BoundaryCondition.mixed_zero_flux_and_activity().set_activity_for_element("C", surface_activity), to=500)`
- `with_left_boundary_condition(BoundaryCondition.closed_system())`

This example sets an closed-system-boundary-condition from start up to 100s and a activity-boundary-condition from 100s to 500s and finally a closed-system-boundary-condition from 500s to the end of simulation.

**Parameters**

- **boundary_condition** – The boundary condition
- **to** – The upper time-limit for `boundary_condition`.

**Returns**  
This `DiffusionNonIsoThermalCalculation` object

**with_options** *(options: tc_python.diffusion.Options, to: float = 1.7976931348623157e+308)*

Sets the general simulation conditions.

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

**5.1. Calculations**
**Default:** The end of the simulation.

**Parameters**

- **options** – The general simulation conditions
- **to** – The upper time-limit for options.

**Returns** This `DiffusionNonIsoThermalCalculation` object

```python
with_planar_geometry()
```

Sets geometry to `planar`.

This is default.

**Returns** This `DiffusionNonIsoThermalCalculation` object

```python
with_reference_state(element: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)
```

The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.

All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

**Parameters**

- **element** – The name of the element
- **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.
- **temperature** – The Temperature (in K) for the reference state. Or `CURRENT_TEMPERATURE` which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.
- **pressure** – The pressure (in Pa) for the reference state

**Returns** This `DiffusionNonIsoThermalCalculation` object

```python
with_right_boundary_condition(boundary_condition: tc_python.diffusion.BoundaryCondition, to: float = 1.7976931348623157e+308)
```

Defines the boundary condition on the right edge of the system.

**Default:** A closed-system boundary condition

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

**Default:** The end of the simulation.
Note: You can specify time-dependent boundary conditions by calling `with_right_boundary_condition()` many times, with different values of the “to” parameter.

Examples:

- `with_right_boundary_condition(BoundaryCondition.closed_system(), to=100)`
- `with_right_boundary_condition(BoundaryCondition.mixed_zero_flux_and_activity().set_activity_for_element("C", surface_activity), to=500)`
- `with_right_boundary_condition(BoundaryCondition.closed_system())`

This example sets an closed-system-boundary-condition from start up to 100s and a activity-boundary-condition from 100s to 500s and finally a closed-system-boundary-condition from 500s to the end of simulation.

Parameters

- `boundary_condition` – The boundary condition
- `to` – The upper time-limit for `boundary_condition`.

Returns This `DiffusionNonIsoThermalCalculation` object

`with_solver` (solver: `tc_python.diffusion.Solver`, to: float = 1.7976931348623157e+308)

Sets the solver to use (`Classic`, `Homogenization` or `Automatic`). **Default is Automatic.**

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

Default: The end of the simulation.

Parameters

- `solver` – The solver to use
- `to` – The upper time-limit for solver.

Returns This `DiffusionNonIsoThermalCalculation` object

`with_spherical_geometry` (first_interface_position: float = 0.0)

Sets geometry to `spherical`, corresponds to a sphere with a certain radius.

Default: A spherical geometry

Note: With a cylindrical or spherical geometry, the system’s zero coordinate (left boundary) is at the centre of the cylinder or sphere by default. By specifying the `first_interface_position`, a different left-most coordinate can be defined. This allows to model a tube or a hollow sphere geometry. The highest coordinate (right boundary) is defined by the cylinder or sphere radius (i.e. by the width of all regions).

Parameters `first_interface_position` – The position of the left-most coordinate along the axis, only necessary for modeling a hollow sphere geometry [m]

Returns This `DiffusionNonIsoThermalCalculation` object

`with_system_modifications` (system_modifications: `tc_python.abstract_base.SystemModifications`)

Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).
**Note:** This is only possible if the system has been read from unencrypted (i.e. *user*) databases loaded as a `*.tdb`-file.

Parameters **system_modifications** – The system modification to be performed

Returns This `DiffusionNonIsoThermalCalculation` object

**with_temperature_profile** *(temperature_profile: tc_python.utils.TemperatureProfile)*

Sets the temperature profile to use with this calculation.

Parameters **temperature_profile** – The temperature profile object (specifying time / temperature points)

Returns This `DiffusionNonIsoThermalCalculation` object

**with_timestep_control** *(timestep_control: tc_python.diffusion.TimestepControl, to: float = 1.7976931348623157e+308)*

Sets the timestep control options.

It is possible specify the upper time-point for which this setting is valid using the parameter “to”.

Default: The end of the simulation.

Parameters

• **timestep_control** – The new timestep control options

• **to** – The upper time-limit for timestep_control.

Returns This `DiffusionNonIsoThermalCalculation` object

**class tc_python.diffusion.DoubleGeometricGrid** *(no_of_points: int = 50, lower_geometrical_factor: float = 1.1, upper_geometrical_factor: float = 0.9)*

Bases: `tc_python.diffusion.CalculatedGrid`

Represents a double geometric grid.

**get_lower_geometrical_factor** () → float

Returns the lower geometrical factor (for the left half).

Returns The lower geometrical factor

**get_no_of_points** () → int

Returns number of grid points.

Returns The number of grid points

**get_type** () → str

Type of the grid.

Returns The type of the grid

**get_upper_geometrical_factor** ()

Returns the upper geometrical factor (for the right half).

Returns The upper geometrical factor

**set_lower_geometrical_factor** *(geometrical_factor: float = 1.1)*

Sets the lower (left half) geometrical factor.
Note: A geometrical factor of larger than one yields a higher density of grid points at the lower end of the half and vice versa for a factor smaller than one.

**Parameters** `geometrical_factor` – The geometrical factor for the left half

**Returns** This `DoubleGeometricGrid` object

```python
set_no_of_points(no_of_points: int = 50)
```
Sets the number of grid points.

**Parameters** `no_of_points` – The number of points

**Returns** This `DoubleGeometricGrid` object

```python
set_upper_geometrical_factor(geometrical_factor: float = 0.9)
```
Sets the upper (right half) geometrical factor.

Note: A geometrical factor of larger than one yields a higher density of grid points at the lower end of the half and vice versa for a factor smaller than one.

**Parameters** `geometrical_factor` – The geometrical factor for the right half

**Returns** This `DoubleGeometricGrid` object

```python
class tc_python.diffusion.ElementProfile
Bases: tc_python.diffusion.AbstractElementProfile
```
Factory class providing objects for configuring a step, function or linear initial concentration profile.

```python
classmethod constant(value: float)
```
Factory method that creates a new constant initial concentration profile.

**Parameters** `value` – The constant composition in the region. [unit as defined in `CompositionProfile`].

**Returns** A new `ConstantProfile` object

```python
classmethod funct(dictra_console_mode_function: str)
```
Factory method that creates a new initial concentration profile defined by a function in DICTRA Console Mode syntax.

**Parameters** `dictra_console_mode_function` – The function, expressed in DICTRA Console Mode syntax.

**Returns** A new `FunctionProfile` object

Note: This is an advanced feature, preferably a complex concentration profile should be generated using third party libraries and added to the simulation using `tc_python.diffusion.PointByPointGrid`.

```python
classmethod linear(start_value: float, end_value: float)
```
Factory method that creates a new linear initial concentration profile.

**Parameters**
• **start_value** – Composition at the left side of the region [unit as defined in CompositionProfile].

• **end_value** – Composition at the right side of the region [unit as defined in CompositionProfile].

Returns A new *LinearProfile* object

classmethod *step*(lower_boundary: float, upper_boundary: float, step_at: float)
Factory method that creates a new initial concentration profile with a step at the specified distance, otherwise the composition is constant at the specified values.

Parameters

• **lower_boundary** – Composition before the step [unit as defined in CompositionProfile].

• **upper_boundary** – Composition after the step [unit as defined in CompositionProfile].

• **step_at** – The distance where the step should be [m].

Returns A new *StepProfile* object

class tc_python.diffusion.FixFluxValue
Bases: tc_python.diffusion.BoundaryCondition

*get_type* () → str
The type of the boundary condition.

Returns The type

*set_flux*(element_name: str, J: str = '0', to_time: float = 1.7976931348623157e+308)
Enter functions that yield the flux times the molar volume for the specified element. May be a function of time, temperature and pressure: \(J(T, P, TIME)\).

Parameters

• **element_name** – The name of the element

• J – the function \(J(T, P, TIME)\)

• **to_time** – The max-time for which the flux function is used.

class tc_python.diffusion.FixedCompositions
Bases: tc_python.diffusion.BoundaryCondition

Represents a boundary having fixed composition values.

*get_type* () → str
The type of the boundary condition.

Returns The type

*set_composition*(element_name: str, value: float)
Sets the composition for the specified element.

Note: The boundary composition needs to be specified for each element.

Parameters

• **element_name** – The name of the element
• **value** – The composition value [unit according to the constructor parameter]

**class** tc_python.diffusion.FunctionProfile

*dictra_console_mode_function: str*,

Bases: tc_python.diffusion.ElementProfile

Creates an initial concentration profile defined by a function in DICTRA Console Mode syntax.

**Note:** This is an advanced feature, preferably a complex concentration profile should be generated using third party libraries and added to the simulation using tc_python.diffusion.PointByPointGrid.

**get_type() → str**

The type of the element profile.

**Returns** The type

**class** tc_python.diffusion.GeneralLowerHashinShtrikman

Bases: tc_python.diffusion.HomogenizationFunctions

General lower Hashin-Shtrikman bounds: the outermost shell consists of the phase with the most sluggish kinetics.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.

**class** tc_python.diffusion.GeneralLowerHashinShtrikmanExcludedPhase

*excluded_phases: List[str] = []*,

Bases: tc_python.diffusion.HomogenizationFunctions

General lower Hashin-Shtrikman bounds: the outermost shell consists of the phase with the most sluggish kinetics.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.

The excluded phases are not considered when evaluating what phase has the most sluggish kinetics.

**class** tc_python.diffusion.GeneralUpperHashinShtrikman

Bases: tc_python.diffusion.HomogenizationFunctions

General upper Hashin-Shtrikman bounds: the innermost shell consists of the phase with the most sluggish kinetics.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.

**class** tc_python.diffusion.GeneralUpperHashinShtrikmanExcludedPhase

*excluded_phases: List[str] = []*,

Bases: tc_python.diffusion.HomogenizationFunctions

General upper Hashin-Shtrikman bounds: the innermost shell consists of the phase with the most sluggish kinetics.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.

The excluded phases are not considered when evaluating what phase has the most sluggish kinetics.

**class** tc_python.diffusion.GeometricGrid

*no_of_points: int = 50, geometrical_factor: float = 1.1*,

Bases: tc_python.diffusion.CalculatedGrid

---

5.1. Calculations
 Represents a geometric grid.

\texttt{get\_geometrical\_factor()} \rightarrow \texttt{float}

Returns the geometrical factor.

\texttt{Returns}  The geometrical factor

\texttt{get\_no\_of\_points()} \rightarrow \texttt{int}

Returns the number of grid points.

\texttt{Returns}  The number of grid points

\texttt{get\_type()} \rightarrow \texttt{str}

Returns the type of grid.

\texttt{Returns}  The type

\texttt{set\_geometrical\_factor(geometrical\_factor: float = 1.1)}

Sets the geometrical factor.

\textbf{Note:}  A geometrical factor larger than one yields a higher density of grid points at the lower end of the region and a factor smaller than one yields a higher density of grid points at the upper end of the region.

\begin{itemize}
  \item \textbf{Parameters} \texttt{geometrical\_factor} – The geometrical factor
  \item \textbf{Returns} This \texttt{GeometricGrid} object
\end{itemize}

\texttt{set\_no\_of\_points(no\_of\_points: int = 50)}

Sets the number of grid points.

\begin{itemize}
  \item \textbf{Parameters} \texttt{no\_of\_points} – The number of points
  \item \textbf{Returns} This \texttt{GeometricGrid} object
\end{itemize}

\textbf{class} \texttt{tc\_python.diffusion.GridPoint(distance: float)}

\textbf{Bases:} \texttt{object}

Represents a grid point, this is used in combination with grids of the type \texttt{tc\_python.diffusion.PointByPointGrid}.

\texttt{add\_composition(element: str, value: float)}

Adds a composition for the specified element to the grid point.

\begin{itemize}
  \item \textbf{Parameters}
    \begin{itemize}
      \item \texttt{element} – The element
      \item \texttt{value} – The composition value [unit as defined for the grid]
    \end{itemize}
  \item \textbf{Returns} This \texttt{GridPoint} object
\end{itemize}

\textbf{class} \texttt{tc\_python.diffusion.HashinShtrikmanBoundMajority}

\textbf{Bases:} \texttt{tc\_python.diffusion.HomogenizationFunctions}

Hashin-Shtrikman bounds with majority phase as matrix phase: the outermost shell consists of the phase with the highest local volume fraction.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.
class tc_python.diffusion.HashinShtrikmanBoundMajorityExcludedPhase(excluded_phases: List[str] = [])

Bases: tc_python.diffusion.HomogenizationFunctions

Hashin-Shtrikman bounds with majority phase as matrix phase: the outermost shell consists of the phase with the highest local volume fraction.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.

The excluded phases are not considered when evaluating what phase has the most sluggish kinetics.

class tc_python.diffusion.HashinShtrikmanBoundPrescribed(matrix_phase: str)

Bases: tc_python.diffusion.HomogenizationFunctions

Hashin-Shtrikman bounds with prescribed phase as matrix phase: the outermost shell consists of the prescribed phase.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.


Bases: tc_python.diffusion.HomogenizationFunctions

class tc_python.diffusion.HomogenizationFunction(value)

Bases: enum.Enum

*Homogenization function* used for the *homogenization solver*. Many homogenization functions are based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase. Default: *RULE_OF_MIXTURES* (i.e. upper Wiener bounds)

**GENERAL_LOWER_HASHIN_SHTRIKMAN** = 0

General lower Hashin-Shtrikman bounds: the outermost shell consists of the phase with the most sluggish kinetics.

**GENERAL_UPPER_HASHIN_SHTRIKMAN** = 1

General upper Hashin-Shtrikman bounds: the innermost shell consists of the phase with the most sluggish kinetics.

**HASHIN_SHTRIKMAN_BOUND_MAJORITY** = 2

Hashin-Shtrikman bounds with majority phase as matrix phase: the outermost shell consists of the phase with the highest local volume fraction.

**INVERSE_RULE_OF_MIXTURES** = 4

Lower Wiener bounds: the geometrical interpretation are continuous layers of each phase orthogonal to the direction of diffusion

**RULE_OF_MIXTURES** = 3

Upper Wiener bounds: the geometrical interpretation are continuous layers of each phase parallel with the direction of diffusion

class tc_python.diffusion.HomogenizationFunctions

Bases: object

*classmethod general_lower_hashin_shtrikman()*

Factory method that creates a new homogenization function of the type

5.1. Calculations
GeneralLowerHashinShtrikman.

General lower Hashin-Shtrikman bounds: the outermost shell consists of the phase with the most sluggish kinetics.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.

**Returns** A new `GeneralLowerHashinShtrikman` object

`classmethod general_lower_hashin_shtrikman_excluded_phase`

Factory method that creates a new homogenization function of the type `GeneralLowerHashinShtrikmanExcludedPhase`.

General lower Hashin-Shtrikman bounds: the outermost shell consists of the phase with the most sluggish kinetics.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase. The excluded phases are not considered when evaluating what phase has the most sluggish kinetics.

**Parameters** `excluded_phases` – The excluded phases

**Returns** A new `GeneralLowerHashinShtrikmanExcludedPhase` object

`classmethod general_upper_hashin_shtrikman`()

Factory method that creates a new homogenization function of the type `GeneralUpperHashinShtrikman`.

General upper Hashin-Shtrikman bounds: the innermost shell consists of the phase with the most sluggish kinetics.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.

**Returns** A new `GeneralUpperHashinShtrikman` object

`classmethod general_upper_hashin_shtrikman_excluded_phase`  
(excluded_phases: `List[str] = []`)

Factory method that creates a new homogenization function of the type `GeneralUpperHashinShtrikmanExcludedPhase`.

General upper Hashin-Shtrikman bounds: the innermost shell consists of the phase with the most sluggish kinetics.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase. The excluded phases are not considered when evaluating what phase has the most sluggish kinetics.

**Parameters** `excluded_phases` – The excluded phases

**Returns** A new `GeneralUpperHashinShtrikmanExcludedPhase` object

`classmethod hashin_shtrikman_bound_majority`()

Factory method that creates a new homogenization function of the type `HashinShtrikmanBoundMajority`.

Hashin-Shtrikman bounds with majority phase as matrix phase: the outermost shell consists of the phase with the highest local volume fraction.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.

**Returns** A new `HashinShtrikmanBoundMajority` object
**classmethod hashin_shtrikman_bound_majority_excluded_phase** *(excluded_phases: List[str] = [])*

Factory method that creates a **new** homogenization function of the type `HashinShtrikmanBoundMajorityExcludedPhase`.

Hashin-Shtrikman bounds with majority phase as matrix phase: the outermost shell consists of the phase with the highest local volume fraction. Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase. The excluded phases are not considered when evaluating what phase has the most sluggish kinetics.

**Parameters** `excluded_phases` – The excluded phases

**Returns** A new `HashinShtrikmanBoundMajorityExcludedPhase` object

**classmethod hashin_shtrikman_bound_prescribed** *(matrix_phase: str)*

Factory method that creates a **new** homogenization function of the type `HashinShtrikmanBoundPrescribed`.

Hashin-Shtrikman bounds with prescribed phase as matrix phase: the outermost shell consists of the prescribed phase.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase.

**Parameters** `matrix_phase` – The matrix phase

**Returns** A new `HashinShtrikmanBoundPrescribed` object

**classmethod hashin_shtrikman_bound_prescribed_excluded_phase** *(matrix_phase: str, excluded_phases: List[str] = [])*

Factory method that creates a **new** homogenization function of the type `HashinShtrikmanBoundPrescribedExcludedPhase`.

Hashin-Shtrikman bounds with prescribed phase as matrix phase: the outermost shell consists of the prescribed phase.

Based on a variant of the Hashin-Shtrikman bounds, their geometrical interpretation are concentric spherical shells of each phase. The excluded phases are not considered when evaluating what phase has the most sluggish kinetics.

**Parameters**

- `matrix_phase` – The matrix phase
- `excluded_phases` – The excluded phases

**Returns** A new `HashinShtrikmanBoundPrescribedExcludedPhase` object

**classmethod inverse_rule_of_mixtures** ()

Factory method that creates a **new** homogenization function of the type `InverseRuleOfMixtures`.

Lower Wiener bounds: the geometrical interpretation are continuous layers of each phase orthogonal to the direction of diffusion.

**Returns** A new `InverseRuleOfMixtures` object

**classmethod inverse_rule_of_mixtures_excluded_phase** *(excluded_phases: List[str] = [])*

Factory method that creates a **new** homogenization function of the type `InverseRuleOfMixturesExcludedPhase`. 

**5.1. Calculations**
Lower Wiener bounds: the geometrical interpretation are continuous layers of each phase orthogonal to the direction of diffusion. Excluded phases are not considered in the diffusion calculations.

**Parameters**

- `excluded_phases` – The excluded phases

**Returns**

A new `InverseRuleOfMixturesExcludedPhase` object

classmethod `labyrinth_factor_f` *(matrix_phase: str)*

Factory method that creates a new homogenization function of the type `LabyrinthFactorF`.

The labyrinth factor functions implies that all diffusion takes place in a single continuous matrix phase. The impeding effect on diffusion by phases dispersed in the matrix phase is taken into account by multiplying the flux with the volume fraction of the matrix phase.

**Parameters**

- `matrix_phase` – The matrix phase

**Returns**

A new `LabyrinthFactorF` object

classmethod `labyrinth_factor_f2` *(matrix_phase: str)*

Factory method that creates a new homogenization function of the type `LabyrinthFactorF2`.

The labyrinth factor functions implies that all diffusion takes place in a single continuous matrix phase. The impeding effect on diffusion by phases dispersed in the matrix phase is taken into account by multiplying the flux with the volume fraction of the matrix phase squared.

**Parameters**

- `matrix_phase` – The matrix phase

**Returns**

A new `LabyrinthFactorF2` object

classmethod `rule_of_mixtures` ()

Factory method that creates a new homogenization function of the type `RuleOfMixtures`.

Upper Wiener bounds: the geometrical interpretation are continuous layers of each phase parallel with the direction of diffusion.

**Returns**

A new `RuleOfMixtures` object

classmethod `rule_of_mixtures_excluded_phase` *(excluded_phases: List[str] = [])*

Factory method that creates a new homogenization function of the type `RuleOfMixturesExcludedPhase`.

Upper Wiener bounds: the geometrical interpretation are continuous layers of each phase parallel with the direction of diffusion. Excluded phases are not considered in the diffusion calculations.

**Parameters**

- `excluded_phases` – The excluded phases

**Returns**

A new `RuleOfMixturesExcludedPhase` object

class `tc_python.diffusion.HomogenizationSolver`

Bases: `tc_python.diffusion.Solver`

Solver using the Homogenization model.

**Note:** This solver always uses the homogenization model, even if all regions have only one phase. The solver is significantly slower than the Classic model. Use the `tc_python.diffusion.AutomaticSolver` instead if you do not need that behavior.

**disable_global_minimization** ()

Disables global minimization to be used in equilibrium calculations. **Default:** Disabled
**Note:** In general, using global minimization **significantly increases the simulation time**, but there is also a significantly reduced risk for non-converged equilibrium calculations.

**Returns** A new `HomogenizationSolver` object

### disable_interpolation_scheme()
Configures the simulation not use *any* interpolation scheme. Default: To use the logarithmic interpolation scheme with 10000 discretization steps

**Note:** The homogenization scheme can be switched on by using `with_linear_interpolation_scheme` or `with_logarithmic_interpolation_scheme`.

### enable_global_minimization()
Enables global minimization to be used in equilibrium calculations. Default: Disabled

**Note:** In general, using global minimization **significantly increases the simulation time**, but there is also a significantly reduced risk for non-converged equilibrium calculations.

**Returns** A new `HomogenizationSolver` object

### get_type() → str
The type of solver.

**Returns** The type

### set_fraction_of_free_memory_to_use(fraction: float)
Sets the maximum fraction of free physical memory to be used by the interpolation scheme. Default: 1 / 10 of the free physical memory

**Parameters** fraction – The maximum free physical memory fraction to be used

**Returns** A new `HomogenizationSolver` object

### set_homogenization_function(homogenization_function_enum:
    tc_python.diffusion.HomogenizationFunction = <HomogenizationFunction.RULE_OF_MIXTURES: 3>)
Sets the homogenization function used by the homogenization model.

**Note:** Deprecated in version 2021b: Use `with_function()` instead. This method will be removed in release 2022b.

Default is RULE_OF_MIXTURES.

**Parameters** homogenization_function_enum – The homogenization function used by the homogenization model

**Returns** A new `HomogenizationSolver` object

### set_memory_to_use(memory_in_megabytes: float)
Sets the maximum physical memory in megabytes to be used by the interpolation scheme. Default: 1000 MBytes of the free physical memory

**Parameters** memory_in_megabytes – The maximum physical memory to be used

5.1. Calculations
Returns A new `HomogenizationSolver` object

with_function (homogenization_function: tc_python.diffusion.HomogenizationFunctions)

Sets the homogenization function used by the homogenization model.

Parameters homogenization_function – The homogenization function used by the homogenization model

Returns A new `HomogenizationSolver` object

with_linear_interpolation_scheme (steps: int = 10000)

Configures the simulation to use the linear interpolation scheme. Default: To use the logarithmic interpolation scheme with 10000 discretization steps.

Parameters steps – The number of discretization steps in each dimension

Returns A new `HomogenizationSolver` object

with_logarithmic_interpolation_scheme (steps: int = 10000)

Configures the simulation to use the linear interpolation scheme. Default: To use the logarithmic interpolation scheme with 10000 discretization steps.

Parameters steps – The number of discretization steps in each dimension

Returns A new `HomogenizationSolver` object

class tc_python.diffusion.InverseRuleOfMixtures

Bases: tc_python.diffusion.HomogenizationFunctions

Lower Wiener bounds: the geometrical interpretation are continuous layers of each phase orthogonal to the direction of diffusion.

class tc_python.diffusion.InverseRuleOfMixturesExcludedPhase (excluded_phases: List[str] = [])

Bases: tc_python.diffusion.HomogenizationFunctions

Lower Wiener bounds: the geometrical interpretation are continuous layers of each phase orthogonal to the direction of diffusion.

Excluded phases are not considered in the diffusion calculations.

class tc_python.diffusion.LabyrinthFactorF (matrix_phase: str)

Bases: tc_python.diffusion.HomogenizationFunctions

The labyrinth factor functions implies that all diffusion takes place in a single continuous matrix phase. The impeding effect on diffusion by phases dispersed in the matrix phase is taken into account by multiplying the flux with the volume fraction of the matrix phase.

class tc_python.diffusion.LabyrinthFactorF2 (matrix_phase: str)

Bases: tc_python.diffusion.HomogenizationFunctions

The labyrinth factor functions implies that all diffusion takes place in a single continuous matrix phase. The impeding effect on diffusion by phases dispersed in the matrix phase is taken into account by multiplying the flux with the volume fraction squared of the matrix phase.

class tc_python.diffusion.LinearGrid (no_of_points: int = 50)

Bases: tc_python.diffusion.CalculatedGrid

Represents an equally spaced grid.

get_no_of_points () -> int

Returns the number of grid points.

Returns The number of grid points
**get_type()** → **str**  
Type of the grid.  

**Returns** The type

**set_no_of_points** *(no_of_points: int = 50)*  
Sets the number of grid points.  

**Parameters** no_of_points – The number of points  

**Returns** This LinearGrid object

---

**class** tc_python.diffusion.LinearProfile *(start_value: float, end_value: float)*  
**Bases:** tc_python.diffusion.ElementProfile

Represents a linear initial concentration profile.

**get_type()** → **str**  
The type of the element profile.

**Returns** The type

---

**class** tc_python.diffusion.MixedZeroFluxAndActivity  
**Bases:** tc_python.diffusion.BoundaryCondition

Represents a boundary having zero-flux as well as fixed-activity conditions.

**Default** On that boundary for every element without an explicitly defined condition, a zero-flux boundary condition is used.

**get_type()** → **str**  
The type of the boundary condition.

**Returns** The type

---

**set_activity_for_element** *(element_name: str, activity: str, to_time: float = 1.7976931348623157e+308)*  
Sets an activity expression for an element at the boundary. Enter a formula that the software evaluates during the calculation.

The formula can be:

- a function of the variable \( \text{TIME} \)
- a constant

The formula must be written with these rules:

- a number must begin with a number (not a .)
- a number must have a dot or an exponent (E)

The operators +, -, *, /, ** (exponentiation) can be used and with any level of parenthesis. As shown, the following operators must be followed by open and closed parentheses ()

- \( \text{SQRT}(X) \) is the square root
- \( \text{EXP}(X) \) is the exponential
- \( \text{LOG}(X) \) is the natural logarithm
- \( \text{LOG10}(X) \) is the base 10 logarithm
- \( \text{SIN}(X), \text{COS}(X), \text{TAN}(X), \text{ASIN}(X), \text{ACOS}(X), \text{ATAN}(X) \)
- \( \text{SINH}(X), \text{COSH}(X), \text{TANH}(X), \text{ASINH}(X), \text{ACOSH}(X), \text{ATANH}(X) \)
- \( \text{SIGN}(X) \)
• \( ERF(X) \) is the error function

**Default:** the expression entered is used for the entire simulation.

**Parameters**
- `element_name` – The name of the element
- `activity` – The activity
- `to_time` – The max-time for which the activity is used.

**set_zero_flux_for_element**(element_name: str)
Sets a zero-flux condition for an element at the boundary. **Default for all elements at the boundary without an explicitly defined condition**

**Parameters**
- `element_name` – The name of the element

```python
class tc_python.diffusion.Options
    Bases: object

    General simulation conditions for the diffusion calculations.

    disable_forced_starting_values_in_equilibrium_calculations()
    Disables forced starting values for the equilibrium calculations. **The default is ‘enable_automatic_forced_starting_values_in_equilibrium_calculations’**.

    Returns This Options object

disable_save_results_to_file()
Disables the saving of results to file during the simulation. **Default:** Saving of the results at every timestep

    Returns This Options object

enable_automatic_forced_starting_values_in_eq_calculations()
Lets calculation engine decide if forced start values for the equilibrium calculations should be used. **This is the default setting.**

    Returns This Options object

enable_forced_starting_values_in_equilibrium_calculations()
Enables forced start values for the equilibrium calculations. **The default is ‘enable_automatic_forced_starting_values_in_equilibrium_calculations’**.

    Returns This Options object

enable_save_results_to_file(every_nth_step: int = -1)
Enables and configures saving of results to file during the simulation. They can be saved for every n-th or optionally for every timestep (-1). **Default:** Saving of the results at every timestep

    Parameters
    - `every_nth_step` – -1 or a value ranging from 0 to 99

    Returns This Options object

enable_time_integration_method_automatic()
Enables automatic selection of integration method. **This is the default method.**

    Returns This Options object

enable_time_integration_method_euler_backwards()
Enables Euler backwards integration. **The default method is enable_time_integration_method_automatic.**
Note: This method is more stable but less accurate and may be necessary if large fluctuations occur in the profiles.

Returns This Options object

def enable_time_integration_method_trapezoidal():
    Enables trapezoidal integration.

Note: If large fluctuations occur in the profiles, it may be necessary to use the more stable but less accurate Euler backwards method.

Returns This Options object

def set_default_driving_force_for_phases_allowed_to_form_at_interf(driving_force: float = 1e-05):
    Sets the default required driving force for phases allowed to form at the interfaces. Default: 1.0e-5

Note: The required driving force (evaluated as $DGM(ph)$) is used for determining whether an inactive phase is stable, i.e. actually formed. $DGM$ represents the driving force normalized by $RT$ and is dimensionless.

Parameters driving_force – The driving force ($DGM(ph)$) [-]

Returns This Options object

class tc_python.diffusion.PointByPointGrid(
    unit_enum: tc_python.diffusion.Unit = <Unit.MASS_PERCENT: 3>)
Bases: tc_python.diffusion.AbstractGrid
Represents a point-by-point grid. This is setting the grid and the compositions at once, it is typically used to enter a measured composition profile or the result from a previous calculation.

Note: If a point-by-point grid is used, it is not necessary to specify the grid and composition profile separately.

def add_point(grid_point: tc_python.diffusion.GridPoint):
    Adds a grid point to the grid.

    Parameters grid_point – The grid point

    Returns This PointByPointGrid object

get_type() → str
    Type of the grid.

    Returns The type

class tc_python.diffusion.Region(name: str)
Bases: object
    Represents a region of the simulation domain that can contain more that one phase.

5.1. Calculations
**Note:** The first added phase represents the matrix phase, while all later added phases are *spheriod phases*, i.e. precipitate phases.

```python
add_phase(phase_name: str, is_matrix_phase: bool = False)
```

Adds a phase to the region, each region must contain at least one phase.

**Note:** Normally the *matrix phase* and the *precipitate phases* are automatically chosen based on the presence of all profile elements in the phase and if it has diffusion data. If multiple phases have equal properties, the phase that was added first is chosen. The matrix phase can be explicitly set by using `is_matrix_phase=True`.

**Note:** If multiple phases are added to a region, the *homogenization model* is applied. That means that average properties of the local phase mixture are used.

**Parameters**
- `phase_name` – The phase name
- `is_matrix_phase` – If set to `True` this phase is explicitly set as matrix phase for the region, if no phase is set to `True`, the matrix phase is chosen automatically

**Returns** This `Region` object

### add_phase_allowed_to_form_at_left_interface(phase_name: str, driving_force: float = 1e-05)

Adds a phase allowed to form at the left boundary of the region (an *inactive phase*). The phase will only appear at the interface as a new automatic region if the driving force to form it is sufficiently high.

**Parameters**
- `phase_name` – The phase name
- `driving_force` – The driving force for the phase to form ($DGM(ph)$)

**Returns** This `Region` object

### add_phase_allowed_to_form_at_right_interface(phase_name: str, driving_force: float = 1e-05)

Adds a phase allowed to form at the right boundary of the region (an *inactive phase*). The phase will only appear at the interface as a new automatic region if the driving force to form it is sufficiently high.

**Parameters**
- `phase_name` – The phase name
- `driving_force` – The driving force for the phase to form ($DGM(ph)$)

**Returns** This `Region` object

### remove_all_phases()

Removes all previously added phases from the region.

**Returns** This `Region` object

### set_width(width: float)

Defined the width of the region.
**Note:** This method needs only to be used if a calculated grid has been defined (using `with_grid()`).

**Parameters** width – The width [m]

**Returns** This `Region` object

### with_composition_profile

```
(with_composition_profile: initial_compositions: tc_python.diffusion.CompositionProfile)
```

Defines the initial composition profiles for all elements in the region.

**Note:** This method needs only to be used if a calculated grid has been defined (using `with_grid()`).

**Parameters** initial_compositions – The initial composition profiles for all elements

**Returns** This `Region` object

### with_grid

```
(with_grid: grid: tc_python.diffusion.CalculatedGrid)
```

Defines a calculated grid in the region. If measured composition profiles or the result from a previous calculation should be used, instead `with_point_by_point_grid_containing_compositions()` needs to be applied.

**Note:** The composition profiles need to be defined separately using `with_composition_profile()`, additionally the region width needs to be specified using `set_width()`.

**Parameters** grid – The grid

**Returns** This `Region` object

### with_point_by_point_grid_containing_compositions

```
(with_point_by_point_grid_containing_compositions: grid: tc_python.diffusion.PointByPointGrid)
```

Defines a point-by-point grid in the region. This is setting the grid and the compositions at once, it is typically used to enter a measured composition profile or the result from a previous calculation. If the composition profile should be calculated (linear, geometric, ...) `with_grid()` should be used instead.

**Note:** If a point-by-point grid is used, `with_grid()`, `with_composition_profile()` and `set_width()` are unnecessary and must not be used.

**Parameters** grid – The point-by-point grid

**Returns** This `Region` object

---

5.1. Calculations
Upper Wiener bounds: the geometrical interpretation are continuous layers of each phase parallel with the direction of diffusion.

Excluded phases are not considered in the diffusion calculations.

```python
class tc_python.diffusion.SimulationTime(value)
Bases: enum.Enum

Specifying special time steps for the evaluation of diffusion results.

**Note:** These placeholders should be used because especially the actual last timestep will slightly differ from the specified end time of the simulation.

**FIRST = 0**
Represents the first timestep of the simulation

**LAST = 1**
Represents the last timestep of the simulation
```

```python
class tc_python.diffusion.Solver
Bases: tc_python.diffusion.AbstractSolver
Factory class providing objects representing a solver.

classmethod automatic()
Factory method that creates a **new automatic solver**. **This is the default solver and recommended for most applications.**

**Note:** This solver uses the homogenization model if any region has more than one phase, otherwise it uses the classic model.

**Returns** A new `AutomaticSolver` object

classmethod classic()
Factory method that creates a **new classic solver.**

**Note:** This solver never switches to the homogenization model even if the solver fails to converge. Use the `tc_python.diffusion.AutomaticSolver` if necessary instead.

**Returns** A new `ClassicSolver` object

classmethod homogenization()
Factory method that creates a **new homogenization solver.**

**Note:** This solver always uses the homogenization model, even if all regions have only one phase. The solver is **significantly slower than the Classic model.** Use the `tc_python.diffusion.AutomaticSolver` instead if you do not need that behavior.

**Returns** A new `HomogenizationSolver` object
class tc_python.diffusion.StepProfile(lower_boundary: float, upper_boundary: float, step_at: float)

Bases: tc_python.diffusion.ElementProfile

Represents an initial constant concentration profile with a step at the specified position.

get_type() → str

The type of the element profile.

Returns: The type

class tc_python.diffusion.TimestepControl

Bases: object

Settings that control the time steps in the simulation.

disable_check_interface_position()

Disables checking of the interface position, i.e. the timesteps are not controlled by the phase interface displacement during the simulation. The default setting is :func:`enable_automatic_check_interface_position`.

Returns: This TimestepControl object

enable_automatic_check_interface_position()

Lets calculation engine decide if checking of the interface position should be used. This is the default setting.

Returns: This TimestepControl object

enable_check_interface_position()

Enables checking of the interface position, i.e. the timesteps are controlled by the phase interface displacement during the simulation. The default setting is :func:`enable_automatic_check_interface_position`.

Returns: This TimestepControl object

set_initial_time_step(initial_time_step: float = 1e-07)

Sets the initial timestep. Default: 1.0e-7 s

Parameters: initial_time_step – The initial timestep [s]

Returns: This TimestepControl object

set_max_absolute_error(absolute_error: float = 1e-05)

Sets the maximum absolute error. Default: 1.0e-5

Parameters: absolute_error – The maximum absolute error

Returns: This TimestepControl object

set_max_relative_error(relative_error: float = 0.05)

Sets the maximum relative error. Default: 0.05

Parameters: relative_error – The maximum relative error

Returns: This TimestepControl object

set_max_timestep_allowed_as_percent_of_simulation_time(max_timestep_allowed_as_percent_of_simulation_time: float = 10.0)

The maximum timestep allowed during the simulation, specified in percent of the simulation time. Default: 10.0%

Parameters: max_timestep_allowed_as_percent_of_simulation_time – The maximum timestep allowed [%]

Returns: This TimestepControl object
**set_max_timestep_increase_factor** *(max_timestep_increase_factor: float = 2.0)*

Sets the maximum timestep increase factor. **Default**: 2

**Note:** For example, if 2 is entered the maximum time step is twice as long as the previous time step taken.

**Parameters**
- **max_timestep_increase_factor** – The maximum timestep increase factor

**Returns**
- This `TimestepControl` object

**set_smallest_time_step_allowed** *(smallest_time_step_allowed: float = 1e-07)*

Sets the smallest time step allowed during the simulation. This is required when using the automatic procedure to determine the time step. **Default**: 1.0e-7s

**Parameters**
- **smallest_time_step_allowed** – The smallest timestep allowed [s]

**Returns**
- This `TimestepControl` object

```python
class tc_python.diffusion.Unit(value):
    Bases: enum.Enum

    Represents a composition unit.

    MASS_FRACTION = 2
    Mass fraction.

    MASS_PERCENT = 3
    Mass percent.

    MOLE_FRACTION = 0
    Mole fraction.

    MOLE_PERCENT = 1
    Mole percent.

    U_FRACTION = 4
    U fraction
```

### 5.1.7 Module “propertymodel”

```python
class tc_python.propertymodel.PropertyModelCalculation(calculator):
    Bases: tc_python.abstract_base.AbstractCalculation

    Configuration for a Property Model calculation.

    **Note:** Specify the settings, the calculation is performed with `calculate()`.
```

**add_poly_command** *(poly_command: str)*

Registers a POLY Console Mode command for execution. These commands are executed after all other configuration directly before the calculation starts to run. All commands are stored and used until explicitly deleted using `remove_all_poly_commands`.

**Parameters**
- **poly_command** – The POLY Console Mode command

**Returns**
- This `PropertyModelCalculation` object
**Note:** It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

**Warning:** As this method runs raw POLY-commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten parenthesis, ...).

**calculate()** → `tc_python.propertymodel.PropertyModelResult`
Runs the Property Model calculation.

**Returns** A `PropertyModelResult` which later can be used to get specific values from the simulation.

**get_argument_default**(`argument_id: str`) → `object`
Returns the default value for the specified argument. The argument id can be obtained with `get_arguments()`.

**Parameters** `argument_id` - The argument id

**Returns** The default value (the type depends on the argument)

**get_argument_description**(`argument_id: str`) → `str`
Returns the detailed description of the argument. The id can be obtained with `get_arguments()`.

**Parameters** `argument_id` - The argument id

**Returns** The detailed description

**get_arguments()** → `Set[str]`
Returns a list of the arguments of the Property Model.

**Note:** The arguments are the ‘UI-panel components’ defined in the Property Model interface method `provide_ui_panel_components()`. They have the same id as specified in the Property Model. The naming is different because there is no UI present.

**Returns** The ids of the available arguments

**get_dynamic_arguments()** → `Set[str]`
Returns a list of the dynamic arguments of the Property Model.

**Note:** Dynamic arguments are “extra” arguments created by pressing the “plus” button that can occur next to the UI-panel for some models, when running the Property Model from within Thermo-Calc. You can use them also from the API by `invoke_dynamic_argument()`.

**Returns** The ids of the available dynamic arguments

**get_model_description()** → `str`
Returns the description text of the current model.

**Returns** The description

**get_model_parameter_value**(`model_parameter_id: str`) → `float`
Returns the current value of an optimizable model parameter. The id can be obtained with `get_model_parameters()`.

---

5.1. Calculations
Parameters `model_parameter_id` – The model parameter id

Returns The current value [unit according to the parameter meaning]

`get_model_parameters()` → `Set[str]`

Returns a list of the optimizable model parameters.

**Note:** The model parameters are an optional set of variables that can be used within the Property Model. Typically they are used to provide the possibility to inject parameter values during an optimization into the model. This allows the dynamic development of Property Models that need to be fitted to experimental data. The model parameters are controlled with the Property Model interface methods `provide_model_parameters` and `set_model_parameter`.

Returns The ids of the optimizable model parameters

`get_system_data()` → `tc_python.abstract_base.SystemData`

Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using `with_system_modifications()`.

**Note:** Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb*-file.

Returns The system data

`invoke_dynamic_argument` (*argument_id*: `str`)

Increases the number of instances of this dynamic argument by one, the argument will have an id such as `argument_1`, `argument_2`, … if the dynamic argument is called `argument`.

**Note:** You can obtain all available dynamic arguments by using `get_dynamic_arguments()`.

Parameters `argument_id` – `argument_id`: The argument id

Returns This `PropertyModelCalculation` object

`remove_all_conditions()`

Removes all set classic POLY conditions.

**Note:** This does not affect the compositions set by `set_composition()`.

Returns This `PropertyModelCalculation` object

`remove_all_poly_commands()`

Removes all previously added POLY Console Mode commands.

Returns This `PropertyModelCalculation` object

`remove_dependent_element()`

Removes a manually set dependent element. This method does not affect the automatic choice of the dependent element if `set_composition()` is used.

Returns This `PropertyModelCalculation` object
**set_argument** *(argument: str, value: str)*
Sets the specified model argument to the specified value. The id can be obtained with `get_arguments()`.

**Parameters**
- **argument** – The argument id
- **value** – The value [unit according to the argument meaning]

**Returns** This `PropertyModelCalculation` object

**set_composition** *(element_name: str, value: float)*
Sets the composition of an element. The unit for the composition can be changed using `set_composition_unit()`.

**Default:** Mole percent (`CompositionUnit.MOLE_PERCENT`)

**Parameters**
- **element_name** – The element
- **value** – The composition value [composition unit defined for the calculation]

**Returns** This `PropertyModelCalculation` object

**set_composition_unit** *(unit_enum: tc_python.utils.CompositionUnit = <CompositionUnit.MOLE_PERCENT: 1>)*
Sets the composition unit.

**Default:** Mole percent (`CompositionUnit.MOLE_PERCENT`).

**Parameters**
- **unit_enum** – The new composition unit

**Returns** This `PropertyModelCalculation` object

**set_condition** *(classic_condition: str, value: float)*
Adds a classic POLY condition. If that method is used, all conditions need to be specified in such a way. If this method is used, it is necessary to set the dependent element manually using `set_dependent_element()`.

**Default if not specified:** pressure P = 1e5 Pa, system size N = 1, Temperature T = 1000 K

**Warning:** There is not possible to mix POLY-commands and compositions using `set_composition()`.

**Note:** It should not be necessary for most users to use this method, try to use `set_composition()` instead.

**Warning:** As this method runs raw POLY-commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten parenthesis, ...).
**set_dependent_element** *(dependent_element_name: str)*
Sets the dependent element manually.

**Parameters**
- **dependent_element_name** – The name of the dependent element

**Returns**
This `PropertyModelCalculation` object

**Note:** It should not be necessary for most users to use this method. Setting the dependent element manually is only necessary and allowed if `set_condition()` is used.

**set_model_parameter** *(model_parameter_id: str, value)*
Resets an optimizable model parameter. The id can be obtained with `get_model_parameters()`.

**Parameters**
- **model_parameter_id** – The model parameter id
- **value** – The new value of the parameter

**Returns**
This `PropertyModelCalculation` object

**set_temperature** *(temperature: float = 1000)*
Sets the temperature.

**Default:** 1000 K

**Parameters**
- **temperature** – The temperature [K]

**Returns**
This `PropertyModelCalculation` object

**with_system_modifications** *(system_modifications: tc_python.abstract_base.SystemModifications)*
Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

**Note:** This is only possible if the system has been read from unencrypted (i.e. *user*) databases loaded as a *.tdb*-file.

**Parameters**
- **system_modifications** – The system modification to be performed

**Returns**
This `PropertyModelCalculation` object

---

**class** `tc_python.propertymodel.PropertyModelResult` *(result)*

**Bases:** `tc_python.abstract_base.AbstractResult`

The result of a Property Model calculation.

**get_result_quantities** () → Set[str]
Returns a list of the available result quantities defined in the Property Model.

**Returns**
The ids of the defined result quantities

**get_result_quantity_description** *(result_quantity_id) → str*
Returns the detailed description of the result quantity. The id can be obtained by `get_result_quantities()`.

**Parameters**
- **result_quantity_id** – The result quantity id

**Returns**
The detailed description
get_single_equilibrium_result(result_quantity_id: str = '') → tc_python.single_equilibrium.SingleEquilibriumResult

Returns a result quantity value. The available result quantities can be obtained by get_result_quantities().

Parameters result_quantity_id – The id of the result quantity.

Returns The requested value [unit depending on the quantity], if the result is a SingleEquilibriumResult, is returned.

get_value_of(result_quantity_id: str) → Union[Float, Dict[str, float]]

Returns a result quantity value. The available result quantities can be obtained by get_result_quantities().

Parameters result_quantity_id – The id of the result quantity

Returns The requested value [unit depending on the quantity]. If the result is parameterized, parameter-value pairs are returned.

save_to_disk(path: str)

Saves the result to disk. The result can later be loaded using tc_python.server.SetUp.load_result_from_disk().

Note: The result data is represented by a whole folder possibly containing multiple files.

Parameters path – The path to the result folder, can be relative or absolute.

Returns This PropertyModelResult object

5.1.8 Module “material_to_material”

class tc_python.material_to_material.AbstractConstantCondition
    Bases: object

The abstract base class for all constant conditions.

class tc_python.material_to_material.AbstractMaterialToMaterialCalculationAxis
    Bases: object

The abstract base class of all calculation axis.

class tc_python.material_to_material.ConstantCondition
    Bases: tc_python.material_to_material.AbstractConstantCondition

A constant condition.

classmethod fraction_of_material_b(fraction_of_material_b: float = 0.5)

Creates a constant fraction of material B condition object.

Note: The unit depends on the composition unit setting in the calculator object.

Parameters fraction_of_material_b – The fraction of material B [weight-fraction or mole-fraction]

Returns The condition object
**classmethod temperature** *(temperature: float = 1000)*

Creates a constant temperature condition object.

**Parameters**
- **temperature** – The temperature [K]

**Returns**
- The condition object

**class** tc_python.material_to_material.FractionOfMaterialBAxis *(from_fraction: float = 0.0, to_fraction: float = 1.0, start_fraction: float = 0.5)*

Bases: tc_python.material_to_material.MaterialToMaterialCalculationAxis

A fraction of material B axis.

**class** tc_python.material_to_material.FractionOfMaterialBCondition *(fraction_of_material_b: float = 0.5)*

Bases: tc_python.material_to_material.ConstantCondition

A constant fraction of material B condition.

**class** tc_python.material_to_material.MaterialToMaterialCalculationAxis

Bases: tc_python.material_to_material.AbstractMaterialToMaterialCalculationAxis

A calculation axis.

**classmethod fraction_of_material_b** *(from_fraction: float = 0.0, to_fraction: float = 1.0, start_fraction: float = 0.5)*

Creates a fraction of material B axis object.

**Note:** The unit depends on the composition unit setting in the calculator.

**classmethod temperature** *(from_temperature: float = 1000, to_temperature: float = 3000, start_temperature: float = 2000)*

Creates a temperature calculation axis object.

**Parameters**
- **from_temperature** – The left axis limit [K]
- **to_temperature** – The right axis limit [K]
- **start_temperature** – The start temperature of the calculation [K]

**Returns**
- A new TemperatureAxis condition object

**class** tc_python.material_to_material.MaterialToMaterialCalculationContainer *(instance)*

Bases: object

Provides access to the calculation objects for all Material to Material calculations.
These are specialised calculations for mixtures of two materials A and B. Otherwise they behave identical to the corresponding regular single equilibrium, property diagram and phase diagram calculations.

```python
with_phase_diagram_calculation (default_conditions: bool = True, components: List[str] = [])
```

Creates a Material to Material phase diagram (map) calculation.

**Parameters**

- **default_conditions** – If True, automatically sets the conditions \( N=1 \) and \( P=100000 \)
- **components** – Specify here the components of the system (for example: \([\text{AL}_2\text{O}_3, \ldots]\)), only necessary if they differ from the elements. If this option is used, all elements of the system need to be replaced by a component.

**Returns** A new `MaterialToMaterialPhaseDiagramCalculation` object

```python
with_property_diagram_calculation (default_conditions: bool = True, components: List[str] = [])
```

Creates a Material to Material property diagram (step) calculation.

**Parameters**

- **default_conditions** – If True, automatically sets the conditions \( N=1 \) and \( P=100000 \)
- **components** – Specify here the components of the system (for example: \([\text{AL}_2\text{O}_3, \ldots]\)), only necessary if they differ from the elements. If this option is used, all elements of the system need to be replaced by a component.

**Returns** A new `MaterialToMaterialPropertyDiagramCalculation` object

```python
with_single_equilibrium_calculation (default_conditions: bool = True, components: List[str] = [])
```

Creates a Material to Material single equilibrium calculation.

**Parameters**

- **default_conditions** – If True, automatically sets the conditions \( N=1 \) and \( P=100000 \)
- **components** – Specify here the components of the system (for example: \([\text{AL}_2\text{O}_3, \ldots]\)), only necessary if they differ from the elements. If this option is used, all elements of the system need to be replaced by a component.

**Returns** A new `MaterialToMaterialSingleEquilibriumCalculation` object

```python
class tc_python.material_to_material.MaterialToMaterialPhaseDiagramCalculation (calculator)
```

Configuration for a Material to Material phase diagram calculation.

**Note:** Specify the conditions, the calculation is performed with `calculate()`.

```python
add_initial_equilibrium (initial_equilibrium: tc_python.step_or_map_diagrams.InitialEquilibrium)
```

Add initial equilibrium start points from which a phase diagram is calculated.

Scans along the axis variables and generates start points when the scan procedure crosses a phase boundary.
It may take a little longer to execute than using the minimum number of start points, as some lines may be calculated more than once. But the core remembers all node points and subsequently stops calculations along a line when it finds a known node point.

It is also possible to create a sequence of start points from one initial equilibria.

**Parameters** initial_equilibrium – The initial equilibrium

**Returns** This `MaterialToMaterialPhaseDiagramCalculation` object

**calculate** (keep_previous_results: bool = False) → `tc_python.material_to_material.MaterialToMaterialPhaseDiagramResult`

Performs the phase diagram calculation.

**Warning:** If you use `keep_previous_results=True`, you must not use another calculator or even get results in between the calculations using `calculate()`. Then the previous results will actually be lost.

**Parameters** keep_previous_results – If `True`, results from any previous call to this method are appended. This can be used to combine calculations with multiple start points if the mapping fails at a certain condition.

**Returns** A new `MaterialToMaterialPhaseDiagramResult` object which later can be used to get specific values from the calculated result.

**disable_global_minimization** ()

Disables global minimization.

**Default:** Enabled

**Returns** This `MaterialToMaterialPhaseDiagramCalculation` object

**dont_keep_default_equilibria** ()

Do not keep the initial equilibria added by default.

This is only relevant in combination with `add_initial_equilibrium()`.

This is the default behavior.

**Returns** This `MaterialToMaterialPhaseDiagramCalculation` object

**enable_global_minimization** ()

Enables global minimization.

**Default:** Enabled

**Returns** This `MaterialToMaterialPhaseDiagramCalculation` object

**get_components** () → List[str]

Returns the names of the components in the system (including all components auto-selected by the database(s)).

**Returns** The component names

**get_gibbs_energy_addition_for** (phase: str) → float

Used to get the additional energy term (always being a constant) of a given phase. The value given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters** phase – Specify the name of the (stoichiometric or solution) phase with the addition
Returns Gibbs energy addition to G per mole formula unit.

**get_system_data() → tc_python.abstract_base.SystemData**

Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using `with_system_modifications()`.

**Note:** Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb*-file.

Returns The system data

**keep_default_equilibria()**

Keep the initial equilibria added by default. This is only relevant in combination with `add_initial_equilibrium()`.

Default behavior is to not keep default equilibria.

**Returns** This `MaterialToMaterialPhaseDiagramCalculation` object

**remove_all_initial_equilibria()**

Removes all previously added initial equilibria.

**Returns** This `MaterialToMaterialPhaseDiagramCalculation` object

**run_poly_command(command: str)**

Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

**Parameters** `command` – The Thermo-Calc Console Mode command

**Returns** This `MaterialToMaterialPhaseDiagramCalculation` object

**Note:** It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

**Warning:** As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).

**set_activities(activities: Dict[str, float])**

Sets the constant activity conditions.

**Note:** The activity conditions are identical for both materials.

**Parameters** `activities` – The constant activities

**Returns** This `MaterialToMaterialPhaseDiagramCalculation` object

**set_composition_unit(unit: tc_python.utils.CompositionUnit = <CompositionUnit.MASS_PERCENT: 3>)**

Sets the composition unit of both materials A and B.

**Default:** Weight percent

**Parameters** `unit` – The composition unit of both materials A and B

**Returns** This `MaterialToMaterialPhaseDiagramCalculation` object

5.1. Calculations
**set_gibbs_energy_addition_for** *(phase: str, gibbs_energy: float)*

Used to specify the additional energy term (always being a constant) of a given phase. The value *(gibbs_energy)* given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters**

- **phase** – Specify the name of the (stoichiometric or solution) phase with the addition
- **gibbs_energy** – Addition to G per mole formula unit

**Returns** This *MaterialToMaterialPhaseDiagramCalculation* object

**set_material_a** *(composition: Dict[str, float], dependent_component: str = None)*

Sets the composition of the material A.

The unit is set with *set_composition_unit()*.

**Tip:** The material can also have constant activity conditions, they are set in *set_activities()*.

**Parameters**

- **composition** – The composition of the material A
- **dependent_component** – The dependent component of the material A

**Returns** This *MaterialToMaterialPhaseDiagramCalculation* object

**set_material_b** *(composition: Dict[str, float], dependent_component: str = None)*

Sets the composition of the material B.

The unit is set with *set_composition_unit()*.

**Tip:** The material can also have constant activity conditions, they are set in *set_activities()*.

**Parameters**

- **composition** – The composition of the material B
- **dependent_component** – The dependent component of the material B

**Returns** This *MaterialToMaterialPhaseDiagramCalculation* object

**set_phase_to_dormant** *(phase: str)*

Sets the phase to the status DORMANT, necessary for calculating the driving force to form the specified phase.

**Parameters** **phase** – The phase name or **ALL_PHASES** for all phases

**Returns** This *MaterialToMaterialPhaseDiagramCalculation* object

**set_phase_to_entered** *(phase: str, amount: float = 1.0)*

Sets the phase to the status ENTERED, that is the default state.

**Parameters**

- **phase** – The phase name or **ALL_PHASES** for all phases
- **amount** – The phase fraction (between 0.0 and 1.0)
Returns This `MaterialToMaterialPhaseDiagramCalculation` object

`set_phase_to_fixed(phase: str, amount: float)`
Sets the phase to the status FIXED, i.e. it is guaranteed to have the specified phase fraction after the calculation.

Parameters
- `phase` – The phase name
- `amount` – The fixed phase fraction (between 0.0 and 1.0)

Returns This `MaterialToMaterialPhaseDiagramCalculation` object

`set_phase_to_suspended(phase: str)`
Sets the phase to the status SUSPENDED, i.e. it is ignored in the calculation.

Parameters `phase` – The phase name or `ALL_PHASES` for all phases

Returns This `MaterialToMaterialPhaseDiagramCalculation` object

`set_pressure(pressure: float)`
Sets the pressure (i.e. the condition \( P \)).

Note: If the flag `default_conditions=True` has been set during the creation of the calculator, the pressure is set to 1000 hPa by default.

Parameters `pressure` – The pressure [Pa]

Returns This `MaterialToMaterialPhaseDiagramCalculation` object

`set_system_size(system_size: float)`
Sets the system size (i.e. the condition ‘N’, the number of moles).

Note: If the flag `default_conditions=True` has been set during the creation of the calculator, the system size is set to 1.0 moles by default.

Parameters `system_size` – The system size [mole]

Returns This `MaterialToMaterialPhaseDiagramCalculation` object

`with_first_axis(axis: tc_python.material_to_material.MaterialToMaterialCalculationAxis)`
Sets the first axis (either temperature or fraction of material B). This calculation type requires that both temperature and fraction of material B axis are set.

Parameters `axis` – The axis

Returns This `MaterialToMaterialPhaseDiagramCalculation` object

`with_options(options: tc_python.step_or_map_diagrams.PhaseDiagramOptions)`
Sets the simulation options.

Parameters `options` – The simulation options

Returns This `PhaseDiagramCalculation` object

`with_reference_state(component: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)`
The reference state for a component is important when calculating activities, chemical potentials and...
enthallpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.

All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

**Parameters**

- **component** – The name of the element must be given.
- **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.
- **temperature** – The Temperature (in K) for the reference state. Or CURRENT_TEMPERATURE which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.
- **pressure** – The Pressure (in Pa) for the reference state.

**Returns**

This `MaterialToMaterialPhaseDiagramCalculation` object

**with_second_axis** *(axis: tc_python.material_to_material.MaterialToMaterialCalculationAxis)*

Sets the second axis (either temperature or fraction of material B). This calculation type requires that both temperature and fraction of material B axis are set.

**Parameters**

- **axis** – The axis

**Returns**

This `MaterialToMaterialPhaseDiagramCalculation` object

**with_system_modifications** *(system_modifications: tc_python.abstract_base.SystemModifications)*

Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

**Note:** This is only possible if the system has been read from unencrypted (i.e. *user*) databases loaded as a *.tdb*-file.

**Parameters**

- **system_modifications** – The system modification to be performed

**Returns**

This `MaterialToMaterialPhaseDiagramCalculation` object

**class** `tc_python.material_to_material.MaterialToMaterialPhaseDiagramResult` *(result)*

**Bases:** `tc_python.step_or_map_diagrams.PhaseDiagramResult`

Result of a Material to Material phase diagram calculation, it can be evaluated using quantities or Console Mode syntax.
add_coordinate_for_phase_label \((x: \text{float}, y: \text{float})\)
Sets a coordinate in the result plot for which the stable phases will be evaluated and provided in the result data object. This can be used to plot the phases of a region into the phase diagram or just to programmatically evaluate the phases in certain regions.

**Warning:** This method takes coordinates of the \textit{plot} axes and not of the calculation axis.

**Parameters**

- \textbf{\textit{x}} – The coordinate of the first \textit{plot} axis (“x-axis”) [unit of the \textit{plot} axis]
- \textbf{\textit{y}} – The coordinate of the second \textit{plot} axis (“y-axis”) [unit of the \textit{plot} axis]

**Returns** This \texttt{MaterialToMaterialPhaseDiagramResult} object

get_values_grouped_by_quantity_of \((x\_quantity: \text{Union}[\text{tc\_python.quantity\_factory.ThermodynamicQuantity, str}], y\_quantity: \text{Union}[\text{tc\_python.quantity\_factory.ThermodynamicQuantity, str}]) \rightarrow \text{tc\_python.step\_or\_map\_diagrams.PhaseDiagramResultValues}\)
Returns x-y-line data grouped by the multiple datasets of the specified quantities (for example in dependency of components). The available quantities can be found in the documentation of the factory class \texttt{ThermodynamicQuantity}. Usually the result data represents the phase diagram.

**Note:** The different datasets will contain NaN-values between different subsections and are not sorted (because they are unsortable due to their nature).

**Note:** It is possible to use functions as axis variables, either by using \texttt{ThermodynamicQuantity}. \texttt{user\_defined\_function}, or by using an expression that contains ‘=’.

**Example**
get_values_grouped_by_quantity_of('T', \texttt{ThermodynamicQuantity.user\_defined\_function('HM.T'))}

**Example** get_values_grouped_by_quantity_of('T', 'CP=HM.T')

**Parameters**

- \textbf{\textit{x\_quantity}} – The first quantity (“x-axis”), Console Mode syntax strings can be used as an alternative (for example ‘T’), \texttt{MATERIAL\_B\_FRACTION}, or even a function (for example ‘f=T*1.01’)
- \textbf{\textit{y\_quantity}} – The second quantity (“y-axis”), Console Mode syntax strings can be used as an alternative (for example ‘NV’), \texttt{MATERIAL\_B\_FRACTION}, or even a function (for example ‘CP=HM.T’)

**Returns** The phase diagram data

get_values_grouped_by_stable_phases_of \((x\_quantity: \text{Union}[\text{tc\_python.quantity\_factory.ThermodynamicQuantity, str}], y\_quantity: \text{Union}[\text{tc\_python.quantity\_factory.ThermodynamicQuantity, str}]) \rightarrow \text{tc\_python.step\_or\_map\_diagrams.PhaseDiagramResultValues}\)
Returns x-y-line data grouped by the sets of “stable phases” (for example “LIQUID” or “LIQUID + FCC\_A1”). The available quantities can be found in the documentation of the factory class \texttt{ThermodynamicQuantity}. Usually the result data represents the phase diagram.
Note: The different datasets will contain NaN-values between different subsections and are not sorted (because they are unsortable due to their nature).

Note: It's possible to use functions as axis variables, either by using ThermodynamicQuantity.
user_defined_function, or by using an expression that contains '='.

Example:
```python
get_values_grouped_by_quantity_of('T', ThermodynamicQuantity.user_defined_function('HM.T'))
```

Example:
```python
get_values_grouped_by_quantity_of('T', 'CP=HM.T')
```

Parameters

- **x_quantity** – The first quantity (“x-axis”), Console Mode syntax strings can be used as an alternative (for example ‘T’), MATERIAL_B_FRACTION, or even a function (for example ‘f=T*1.01’)

- **y_quantity** – The second quantity (“y-axis”), Console Mode syntax strings can be used as an alternative (for example ‘NV’), MATERIAL_B_FRACTION, or even a function (for example ‘CP=HM.T’)

Returns: The phase diagram data

```python
remove_phase_labels()
```
Erases all added coordinates for phase labels.

Returns: This `MaterialToMaterialPhaseDiagramResult` object

```python
save_to_disk(path: str)
```
Saves the result to disc. Note that a result is a folder, containing potentially many files. The result can later be loaded with `load_result_from_disk()`

Parameters: path – the path to the folder you want the result to be saved in. It can be relative or absolute.

Returns: this `MaterialToMaterialPhaseDiagramResult` object

```python
set_phase_name_style(phase_name_style_enum: tc_python.step_or_map_diagrams.PhaseNameStyle = <PhaseNameStyle.NONE: 0>)
```
Sets the style of the phase name labels that will be used in the result data object (constitution description, ordering description, …).

Default: PhaseNameStyle.NONE

Parameters: phase_name_style_enum – The phase name style

Returns: This `MaterialToMaterialPhaseDiagramResult` object

```python
class tc_python.material_to_material.MaterialToMaterialPropertyDiagramCalculation(calculator)
```
Bases: `tc_python.step_or_map_diagrams.AbstractPropertyDiagramCalculation`

Configuration for a Material to Material property diagram calculation.

Note: Specify the conditions and possibly other settings, the calculation is performed with `calculate()`.

```python
calculate(keep_previous_results: bool = False) → tc_python.material_to_material.MaterialToMaterialPropertyDiagramResult
```
Performs the Material to Material property diagram calculation.
**Warning:** If you use `keep_previous_results=True`, you must not use another calculator or even get results in between the calculations using `calculate()`. Then the previous results will actually be lost.

**Parameters** `keep_previous_results` – If `True`, results from any previous call to this method are appended. This can be used to combine calculations with multiple start points if the stepping fails at a certain condition.

**Returns** A new `MaterialToMaterialPropertyDiagramResult` object which later can be used to get specific values from the calculated result

**disable_global_minimization()**
Disables global minimization.

**Default:** Enabled

**Returns** This `MaterialToMaterialPropertyDiagramCalculation` object

**disable_step_separate_phases()**
Disables step separate phases. This is the default setting.

**Returns** This `MaterialToMaterialPropertyDiagramCalculation` object

**enable_global_minimization()**
Enables global minimization.

**Default:** Enabled

**Returns** This `MaterialToMaterialPropertyDiagramCalculation` object

**enable_step_separate_phases()**
Enables step separate phases.

**Default:** By default separate phase stepping is disabled

**Note:** This is an advanced option, it is used mostly to calculate how the Gibbs energy for a number of phases varies for different compositions. This is particularly useful to calculate Gibbs energies for complex phases with miscibility gaps and for an ordered phase that is never disordered (e.g. SIGMA-phase, G-phase, MU-phase, etc.).

**Returns** This `MaterialToMaterialPropertyDiagramCalculation` object

**get_components() → List[str]**
Returns the names of the components in the system (including all components auto-selected by the database(s)).

**Returns** The component names

**get_gibbs_energy_addition_for**(phase: str) → float
Used to get the additional energy term (always being a constant) of a given phase. The value given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters** `phase` – Specify the name of the (stoichiometric or solution) phase with the addition

**Returns** Gibbs energy addition to G per mole formula unit.

5.1. Calculations
**get_system_data** () → tc_python.abstract_base.SystemData

Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using `with_system_modifications()`.

**Note:** Parameters can only be read from unencrypted (i.e. user) databases loaded as *.*tdb-file.

**Returns** The system data

**run_poly_command**(command: str)

Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

**Parameters** command – The Thermo-Calc Console Mode command

**Returns** This MaterialToMaterialPropertyDiagramCalculation object

**Note:** It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

**Warning:** As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).

**set_activities**(activities: Dict[str, float])

Sets the constant activity conditions.

**Note:** The activity conditions are identical for both materials.

**Parameters** activities – The constant activities

**Returns** This MaterialToMaterialPropertyDiagramCalculation object

**set_composition_unit**(unit: tc_python.utils.CompositionUnit = <CompositionUnit.MASS_PERCENT: 3>)

Sets the composition unit of both materials A and B.

**Default:** Weight percent

**Parameters** unit – The composition unit of both materials A and B

**Returns** This MaterialToMaterialPropertyDiagramCalculation object

**set_gibbs_energy_addition_for**(phase: str, gibbs_energy: float)

Used to specify the additional energy term (always being a constant) of a given phase. The value (gibbs_energy) given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters**

- **phase** – Specify the name of the (stoichiometric or solution) phase with the addition
- **gibbs_energy** – Addition to G per mole formula unit

**Returns** This MaterialToMaterialPropertyDiagramCalculation object
set_material_a (composition: Dict[str, float], dependent_component: str = None)
Sets the composition of the material A.

The unit is set with `set_composition_unit()`.

**Tip:** The material can also have constant activity conditions, they are set in `set_activities()`.

**Parameters**
- **composition** – The composition of the material A
- **dependent_component** – The dependent component of the material A

**Returns** This `MaterialToMaterialPropertyDiagramCalculation` object

set_material_b (composition: Dict[str, float], dependent_component: str = None)
Sets the composition of the material B.

The unit is set with `set_composition_unit()`.

**Tip:** The material can also have constant activity conditions, they are set in `set_activities()`.

**Parameters**
- **composition** – The composition of the material B
- **dependent_component** – The dependent component of the material B

**Returns** This `MaterialToMaterialPropertyDiagramCalculation` object

set_phase_to_dormant (phase: str)
Sets the phase to the status DORMANT, necessary for calculating the driving force to form the specified phase.

**Parameters** phase – The phase name or `ALL_PHASES` for all phases

**Returns** This `MaterialToMaterialPropertyDiagramCalculation` object

set_phase_to_entered (phase: str, amount: float = 1.0)
Sets the phase to the status ENTERED, that is the default state.

**Parameters**
- **phase** – The phase name or `ALL_PHASES` for all phases
- **amount** – The phase fraction (between 0.0 and 1.0)

**Returns** This `MaterialToMaterialPropertyDiagramCalculation` object

set_phase_to_fixed (phase: str, amount: float)
Sets the phase to the status FIXED, i.e. it is guaranteed to have the specified phase fraction after the calculation.

**Parameters**
- **phase** – The phase name
- **amount** – The fixed phase fraction (between 0.0 and 1.0)

**Returns** This `MaterialToMaterialPropertyDiagramCalculation` object
**set_phase_to_suspended** *(phase: str)*
Sets the phase to the status SUSPENDED, i.e. it is ignored in the calculation.

**Parameters**
- **phase** – The phase name or *ALL_PHASES* for all phases

**Returns**
This *MaterialToMaterialPropertyDiagramCalculation* object

**set_pressure** *(pressure: float)*
Sets the pressure (i.e. the condition \( P \)).

**Note:** If the flag *default_conditions=True* has been set during the creation of the calculator, the pressure is set to 1000 hPa by default.

**Parameters**
- **pressure** – The pressure [Pa]

**Returns**
This *MaterialToMaterialPropertyDiagramCalculation* object

**set_system_size** *(system_size: float)*
Sets the system size (i.e. the condition ‘N’, the number of moles).

**Note:** If the flag *default_conditions=True* has been set during the creation of the calculator, the system size is set to 1.0 moles by default.

**Parameters**
- **system_size** – The system size [mole]

**Returns**
This *MaterialToMaterialPropertyDiagramCalculation* object

**with_axis** *(axis: tc_python.material_to_material.MaterialToMaterialCalculationAxis)*
Sets the axis (either temperature or fraction of material B). This calculation type requires that either temperature or fraction of material B is set as a constant condition - the other one is set as an axis.

**Parameters**
- **axis** – The axis

**Returns**
This *MaterialToMaterialPropertyDiagramCalculation* object

**with_constant_condition** *(condition: tc_python.material_to_material.ConstantCondition)*
Sets the constant condition (either temperature of fraction of material B). This calculation type requires that either temperature or fraction of material B is set as a constant condition - the other one is set as an axis.

**Parameters**
- **condition** – The condition

**Returns**
This *MaterialToMaterialPropertyDiagramCalculation* object

**with_options** *(options: tc_python.step_or_map_diagrams.PropertyDiagramOptions)*
Sets the simulation options.

**Parameters**
- **options** – The simulation options

**Returns**
This *MaterialToMaterialPropertyDiagramCalculation* object

**with_reference_state** *(component: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)*
The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state.
All data in all phases where this component dissolves must use the same reference state. However, different
databases can use different reference states for the same element/component. It is important to be careful
when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by
the database. If the reference state in the database is not suitable for your purposes, use this command to
set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set
as default for a major component in alloys dominated by the component). In such cases, the temperature
and pressure for the reference state is not needed.

For a phase to be usable as a reference for a component, the component needs to have the same composition
as an end member of the phase. The reference state is an end member of a phase. The selection of the end
member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the
end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

**Parameters**

- **component** – The name of the element must be given.
- **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element
  Reference.
- **temperature** – The Temperature (in K) for the reference state. Or CURRENT_TEMPERATURE which means
  that the current temperature is used at the time of evaluation of the reference energy for the calculation.
- **pressure** – The Pressure (in Pa) for the reference state.

**Returns**

This `MaterialToMaterialPropertyDiagramCalculation` object

**with_system_modifications**

Updates the system of this calculator with the supplied system modification (containing new phase param-
eters and system functions).

---

**Note:** This is only possible if the system has been read from unencrypted (i.e. `user`) databases loaded as
a *.tdb*-file.

---

**Parameters**

- **system_modifications** – The system modification to be performed

**Returns**

This `MaterialToMaterialPropertyDiagramCalculation` object

class `tc_python.material_to_material.MaterialToMaterialPropertyDiagramResult`

Result of a Material to Material property diagram. It can be used to query for specific values.

**get_values_grouped_by_quantity_of**

Returns x-y-line data grouped by the multiple datasets of the specified quantities (typically the phases). The
available quantities can be found in the documentation of the factory class `ThermodynamicQuantity`.

**Note:** The different datasets might contain `NaN`-values between different subsections and might not be
sorted even if the flag `sort_and_merge` has been set (because they might be unsortable due to their
nature).

5.1. Calculations 169
Note: It's possible to use functions as axis variables, either by using `ThermodynamicQuantity.user_defined_function`, or by using an expression that contains '=', i.e., `ThermodynamicQuantity.user_defined_function('HM.T')`.

Example

```python
get_values_grouped_by_quantity_of('T', ThermodynamicQuantity.user_defined_function('HM.T'))
```

Examples

```python
get_values_grouped_by_quantity_of('T', 'CP=HM.T')
```

Parameters

- **x_quantity** – The first quantity ("x-axis"), Console Mode syntax strings can be used as an alternative (for example ‘T’), `MATERIAL_B_FRACTION`, or even a function (for example ‘f=T*1.01’)

- **y_quantity** – The second quantity ("y-axis"), Console Mode syntax strings can be used as an alternative (for example ‘NV’), `MATERIAL_B_FRACTION`, or even a function (for example ‘CP=HM.T’)

- **sort_and_merge** – If True, the data is sorted and merged into as few subsections as possible (divided by `NaN`)

Returns

Containing the datasets with the quantities as their keys

```python
def get_values_grouped_by_stable_phases_of(x_quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], y_quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str], sort_and_merge: bool = True) -> Dict[str, tc_python.utils.ResultValueGroup]
```

Returns x-y-line data grouped by the sets of “stable phases” (for example “LIQUID” or “LIQUID + FCC_A1”). The available quantities can be found in the documentation of the factory class `ThermodynamicQuantity`.

Note: The different datasets might contain `NaN`-values between different subsections and different lines of an ambiguous dataset. They might not be sorted even if the flag `sort_and_merge` has been set (because they might be unsortable due to their nature).

Note: It's possible to use functions as axis variables, either by using `ThermodynamicQuantity.user_defined_function`, or by using an expression that contains '=', i.e., `ThermodynamicQuantity.user_defined_function('HM.T')`.

Example

```python
get_values_grouped_by_quantity_of('T', ThermodynamicQuantity.user_defined_function('HM.T'))
```

Examples

```python
get_values_grouped_by_quantity_of('T', 'CP=HM.T')
```

Parameters

- **x_quantity** – The first quantity ("x-axis"), Console Mode syntax strings can be used as an alternative (for example ‘T’), `MATERIAL_B_FRACTION`, or even a function (for example ‘f=T*1.01’)

- **y_quantity** – The second quantity ("y-axis"), Console Mode syntax strings can be used as an alternative (for example ‘NV’), `MATERIAL_B_FRACTION`, or even a function (for example ‘CP=HM.T’)

Chapter 5. API Reference
• **sort_and_merge** – If True, the data will be sorted and merged into as few subsections as possible (divided by NaN)

**Returns** Containing the datasets with the quantities as their keys

```python
get_values_of(x_quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str],
y_quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str]) \rightarrow
typing.List[float], typing.List[float]
```

Returns sorted x-y-line data without any separation. Use `get_values_grouped_by_quantity_of()` or `get_values_grouped_by_stable_phases_of()` instead if you need such a separation. The available quantities can be found in the documentation of the factory class `ThermodynamicQuantity`.

**Note:** This method will always return sorted data without any NaN-values. If it is unsortable that might give data that is hard to interpret. In such a case you need to choose the quantity in another way or use one of the other methods. One example of this is to use quantities with All-markers, for example `MassFractionOfAComponent("All")`.

**Note:** It's possible to use functions as axis variables, either by using `ThermodynamicQuantity.user_defined_function()` or by using an expression that contains ‘=’.

**Example**

```python
get_values_grouped_by_quantity_of('T', ThermodynamicQuantity.user_defined_function('HM.T'))
```

**Example**

```python
get_values_grouped_by_quantity_of('T', 'CP=HM.T')
```

**Parameters**

- **x_quantity** – The first thermodynamic quantity (“x-axis”), Console Mode syntax strings can be used as an alternative (for example ‘T’, `MATERIAL_B_FRACTION`, or even a function (for example ‘f=T*1.01’).
- **y_quantity** – The second thermodynamic quantity (“y-axis”), Console Mode syntax strings can be used as an alternative (for example ‘NV’, `MATERIAL_B_FRACTION`, or even a function (for example ‘CP=HM.T’)

**Returns** A tuple containing the x- and y-data in lists

```python
save_to_disk(path: str)
```

Saves the result to disc. Note that a result is a folder, containing potentially many files. The result can later be loaded with `load_result_from_disk()`

**Parameters** **path** – the path to the folder you want the result to be saved in. It can be relative or absolute.

**Returns** This `MaterialToMaterialPropertyDiagramResult` object

```python
set_phase_name_style(phase_name_style_enum: tc_python.step_or_map_diagrams.PhaseNameStyle = <PhaseNameStyle.NONE: 0>)
```

Sets the style of the phase name labels that will be used in the result data object (constitution description, ordering description, ...).

**Default:** PhaseNameStyle.NONE

**Parameters** **phase_name_style_enum** – The phase name style

**Returns** This `MaterialToMaterialPropertyDiagramResult` object

---

5.1. Calculations
class tc_python.material_to_material.MaterialToMaterialSingleEquilibriumCalculation(calculator)
Bases: tc_python.single_equilibrium.AbstractSingleEquilibriumCalculation

Configuration for a Material to Material single fraction of B calculation.

**Note:** Specify the conditions and possibly other settings, the calculation is performed with `calculate()`.

`calculate()` \rightarrow tc_python.material_to_material.MaterialToMaterialSingleEquilibriumResult
Performs the material to material calculation.

**Note:** The calculation result is **no** temporary result object.

**Returns** A new `MaterialToMaterialSingleEquilibriumResult` object which can be used to get specific values from the calculated result. It is undefined behavior to use that object after the state of the calculation has been changed.

disable_global_minimization()
Turns the global minimization completely off.

**Returns** This `MaterialToMaterialSingleEquilibriumCalculation` object

enable_global_minimization()
Turns the global minimization on (using the default settings).

**Returns** This `MaterialToMaterialSingleEquilibriumCalculation` object

get_components() \rightarrow List[str]
Returns a list of components in the system (including all components auto-selected by the database(s)).

**Returns** The components

get_gibbs_energy_addition_for(phase: str) \rightarrow float
Used to get the additional energy term (always being a constant) of a given phase. The value given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

**Parameters** phase – Specify the name of the (stoichiometric or solution) phase with the addition

**Returns** Gibbs energy addition to G per mole formula unit.

get_system_data() \rightarrow tc_python.abstract_base.SystemData
Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using `with_system_modifications()`.

**Note:** Parameters can only be read from unencrypted (i.e. **user**) databases loaded as *.tdb-file.

**Returns** The system data

run_poly_command(command: str)
Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine.

**Parameters** command – The Thermo-Calc Console Mode command
set_activities(activities: Dict[str, float])
Sets the constant activity conditions.

Note: The activity conditions are identical for both materials.

Parameters activities – The constant activities

Returns This MaterialToMaterialSingleEquilibriumCalculation object

set_component_to_entered(component: str)
Sets the specified component to the status ENTERED, that is the default state.

Parameters component – The component name or ALL_COMPONENTS

Returns This MaterialToMaterialSingleEquilibriumCalculation object

set_component_to_suspended(component: str, reset_conditions: bool = False)
Sets the specified component to the status SUSPENDED, i.e. it is ignored in the calculation.

Parameters

• reset_conditions – if ‘True’ also remove composition conditions for the component if they are defined
• component – The component name or ALL_COMPONENTS

Returns This MaterialToMaterialSingleEquilibriumCalculation object

set_composition_unit(unit: tc_python.utils.CompositionUnit = <CompositionUnit.MASS_PERCENT: 3>)
Sets the composition unit of both materials A and B.

Default: Weight percent

Parameters unit – The composition unit of both materials A and B

Returns This MaterialToMaterialSingleEquilibriumCalculation object

set_gibbs_energy_addition_for(phase: str, gibbs_energy: float)
Used to specify the additional energy term (always being a constant) of a given phase. The value (gibbs_energy) given is added to the Gibbs energy of the (stoichiometric or solution) phase. It can represent a nucleation barrier, surface tension, elastic energy, etc.

It is not composition-, temperature- or pressure-dependent.

Parameters

• phase – Specify the name of the (stoichiometric or solution) phase with the addition
• gibbs_energy – Addition to G per mole formula unit
set_material_a (composition: Dict[str, float], dependent_component: str = None)
Sets the composition of the material A.
The unit is set with set_composition_unit().

Parameters
- composition – The composition of the material A
- dependent_component – The dependent component of the material A

Returns This MaterialToMaterialSingleEquilibriumCalculation object

Tip: The material can also have constant activity conditions, they are set in set_activities().

set_material_b (composition: Dict[str, float], dependent_component: str = None)
Sets the composition of the material B.
The unit is set with set_composition_unit().

Parameters
- composition – The composition of the material B
- dependent_component – The dependent component of the material B

Returns This MaterialToMaterialSingleEquilibriumCalculation object

Tip: The material can also have constant activity conditions, they are set in set_activities().

set_phase_to_dormant (phase: str)
Sets the phase to the status DORMANT, necessary for calculating the driving force to form the specified phase.

Parameters phase – The phase name or ALL_PHASES for all phases

Returns This MaterialToMaterialSingleEquilibriumCalculation object

set_phase_to_entered (phase: str, amount: float = 1.0)
Sets the phase to the status ENTERED, that is the default state.

Parameters
- phase – The phase name or ALL_PHASES for all phases
- amount – The phase fraction (between 0.0 and 1.0)

Returns This MaterialToMaterialSingleEquilibriumCalculation object

set_phase_to_fixed (phase: str, amount: float)
Sets the phase to the status FIXED, i.e. it is guaranteed to have the specified phase fraction after the calculation.

Parameters
- phase – The phase name
- amount – The fixed phase fraction (between 0.0 and 1.0)
set_phase_to_suspended(phase: str)
Sets the phase to the status SUSPENDED, i.e. it is ignored in the calculation.

Parameters phase – The phase name or ALL_PHASES for all phases

Returns This MaterialToMaterialSingleEquilibriumCalculation object

set_pressure(pressure: float)
Sets the pressure (i.e. the condition $P$).

Note: If the flag default_conditions=True has been set during the creation of the calculator, the pressure is set to 1000 hPa by default.

Parameters pressure – The pressure [Pa]

Returns This MaterialToMaterialSingleEquilibriumCalculation object

set_system_size(system_size: float)
Sets the system size (i.e. the condition ‘N’, the number of moles).

Note: If the flag default_conditions=True has been set during the creation of the calculator, the system size is set to 1.0 moles by default.

Parameters system_size – The system size [mole]

Returns This MaterialToMaterialSingleEquilibriumCalculation object

with_first_constant_condition(condition: tc_python.material_to_material.ConstantCondition)
Sets the first constant condition (either temperature or fraction of material B).

Parameters condition – The condition

Returns This MaterialToMaterialSingleEquilibriumCalculation object

with_options(options: tc_python.single_equilibrium.SingleEquilibriumOptions)
Sets the simulation options.

Parameters options – The simulation options

Returns This SingleEquilibriumCalculation object

with_reference_state(component: str, phase: str = 'SER', temperature: float = -1.0, pressure: float = 100000.0)
The reference state for a component is important when calculating activities, chemical potentials and enthalpies and is determined by the database being used. For each component the data must be referred to a selected phase, temperature and pressure, i.e. the reference state. All data in all phases where this component dissolves must use the same reference state. However, different databases can use different reference states for the same element/component. It is important to be careful when combining data obtained from different databases.

By default, activities, chemical potentials and so forth are computed relative to the reference state used by the database. If the reference state in the database is not suitable for your purposes, use this command to set the reference state for a component using SER, i.e. the Stable Element Reference (which is usually set as default for a major component in alloys dominated by the component). In such cases, the temperature and pressure for the reference state is not needed.

5.1. Calculations
For a phase to be usable as a reference for a component, the component needs to have the same composition as an end member of the phase. The reference state is an end member of a phase. The selection of the end member associated with the reference state is only performed once this command is executed.

If a component has the same composition as several end members of the chosen reference phase, then the end member that is selected at the specified temperature and pressure will have the lowest Gibbs energy.

**Parameters**

- **component** – The name of the element must be given.
- **phase** – Name of a phase used as the new reference state. Or SER for the Stable Element Reference.
- **temperature** – The Temperature (in K) for the reference state. Or CURRENT_TEMPERATURE which means that the current temperature is used at the time of evaluation of the reference energy for the calculation.
- **pressure** – The Pressure (in Pa) for the reference state.

**Returns**

This MaterialToMaterialSingleEquilibriumCalculation object

```python
def with_second_constant_condition(condition: tc_python.material_to_material.ConstantCondition):
    """Sets the second constant condition (either temperature of fraction of material B)."
    Parameters condition -- The condition
    Returns This MaterialToMaterialSingleEquilibriumCalculation object
```

**with_system_modifications**

```python
def with_system_modifications(system_modifications: tc_python.abstract_base.SystemModifications):
    """Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions)."
    Parameters system_modifications -- The system modification to be performed
    Returns This MaterialToMaterialSingleEquilibriumCalculation object
```

**Note:**

This is only possible if the system has been read from unencrypted (i.e. user) databases loaded as a *.tdb*-file.

**class**

```python
class tc_python.material_to_material.MaterialToMaterialSingleEquilibriumResult(
r
    result
```

**Parameters**

- **pressure** – The pressure [Pa]

**Returns**

This MaterialToMaterialSingleEquilibriumResult object

```python
def change_pressure(pressure: float):
    """Change the pressure and re-evaluate the results from the equilibrium without minimizing Gibbs energy, i.e. with high performance. The properties are calculated at the new pressure using the phase amount, temperature and composition of phases from the initial equilibrium. Use get_value_of() to obtain them."
    Parameters pressure -- The pressure [Pa]
    Returns This MaterialToMaterialSingleEquilibriumResult object
```

```python
def change_temperature(temperature: float):
    """Change the temperature and re-evaluate the results from the equilibrium without minimizing Gibbs energy, i.e. with high performance. The properties are calculated at the new temperature using the phase amount, pressure and composition of phases from the initial equilibrium. Use get_value_of() to obtain them."
```
Note: This is typically used when calculating room temperature properties (e.g. density) for a material when it is assumed that the equilibrium phase amount and composition freeze-in at a higher temperature during cooling.

**Parameters**

- **temperature** – The temperature [K]

**Returns**

This `MaterialToMaterialSingleEquilibriumResult` object

---

**get_components** () → List[str]

Returns the names of the components selected in the system (including any components auto-selected by the database(s)).

**Returns**

The names of the selected components

**get_conditions** () → List[str]

Returns the conditions.

**Returns**

The selected conditions

**get_phases** () → List[str]

Returns the phases present in the system due to its configuration. It also contains all phases that have been automatically added during the calculation, this is the difference to the method `System.get_phases_in_system()`.

**Returns**

The names of the phases in the system including automatically added phases

**get_stable_phases** () → List[str]

Returns the stable phases (i.e. the phases present in the current equilibrium).

**Returns**

The names of the stable phases

**get_value_of** (`quantity: Union[tc_python.quantity_factory.ThermodynamicQuantity, str]`) → float

Returns a value from a single equilibrium calculation.

**Parameters**

- **quantity** – The thermodynamic quantity to get the value of; a Console Mode syntax strings can be used as an alternative (for example “NPM(FCC_A1)"

**Returns**

The requested value

**run_poly_command** (`command: str`)  
Runs a Thermo-Calc command from the Console Mode POLY module immediately in the engine. This affects only the state of the result object.

**Parameters**

- **command** – The Thermo-Calc Console Mode command

**Returns**

This `MaterialToMaterialSingleEquilibriumResult` object

---

**Note:** It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

---

**Warning:** As this method runs raw Thermo-Calc commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten equals sign).

**save_to_disk** (`path: str`)  
Saves the result to disk. Note that the result is a folder, containing potentially many files. The result can later be loaded with `load_result_from_disk()`
Parameters `path` – the path to the folder you want the result to be saved in. It can be relative or absolute.

Returns this `MaterialToMaterialSingleEquilibriumResult` object

class tc_python.material_to_material.TemperatureAxis (from_temperature: float = 1000, to_temperature: float = 3000, start_temperature: float = 2000)

Bases: `tc_python.material_to_material.MaterialToMaterialCalculationAxis` A temperature calculation axis.

class tc_python.material_to_material.TemperatureCondition (temperature: float = 1000.0)

Bases: `tc_python.material_to_material.ConstantCondition` A constant temperature condition.

5.2 Module “system”

class tc_python.system.MultiDatabaseSystemBuilder (multi_database_system_builder)

Bases: object

Used to select databases, elements, phases etc. and create a System object. The difference to the class SystemBuilder is that the operations are performed on all the previously selected databases. The system is then used to create calculations.

create_and_select_species (stoichiometry: str)

Specify a species from the already entered elements. The stoichiometry of the species is the chemical formula of the species. The created species will also be automatically selected.

Note: The elements in the chemical formula are normally separated by stoichiometric numbers. Neither parenthesis “()” nor an underscore “_” is allowed in the chemical formula, while the special combination “/-” or “/+” can be used. Consult the Thermo-Calc database documentation for details about the syntax.

Parameters `stoichiometry` – The stoichiometry of the species

Returns This `MultiDatabaseSystemBuilder` object

deselect_constituent_on_sublattice (phase_name: str, sublattice_no: int, constituent_name_to_deselect: str)

Rejects a constituent on a sublattice in a phase in both the thermodynamic and the kinetic database.

Parameters

• `phase_name` – The name of the phase

• `sublattice_no` – The number of the sublattice (starting with 1)

• `constituent_name_to_deselect` – The name of the constituent to deselect

Returns This `MultiDatabaseSystemBuilder` object

deselect_phase (phase_name_to_deselect: str)

Rejects a phase for both the thermodynamic and the kinetic database.

Parameters `phase_name_to_deselect` – The phase name

Returns This `MultiDatabaseSystemBuilder` object
deselect_species

Removes the species from the system.

Parameters

species_name – The species

Returns

This :class:`MultiDatabaseSystemBuilder` object

get_system

Creates a new System object that is the basis for all calculation types. Several calculation types can be defined later from the object; these are independent.

Returns

A new :class:`System` object

select_constituent_on_sublattice

Selects a constituent on a sublattice in a phase in both the thermodynamic and the kinetic database.

Note: Previously the third parameter constituent_name_to_select had a wrong name, it has been corrected in version 2021b.

Parameters

- phase_name – The name of the phase
- sublattice_no – The number of the sublattice (starting with 1)
- constituent_name_to_select – The name of the constituent to select

Returns

This :class:`MultiDatabaseSystemBuilder` object

select_phase

Selects a phase for both the thermodynamic and the kinetic database.

Parameters

phase_name_to_select – The phase name

Returns

This :class:`MultiDatabaseSystemBuilder` object

select_species

Adds the species to the system. Up to 1000 species can be defined in a single system.

Parameters

species_name – The species

Returns

This :class:`MultiDatabaseSystemBuilder` object

with_new_composition_set

Used to enter two or more composition sets for a phase. If a phase has a miscibility gap it is necessary to have two composition sets, one for each possible composition that can be stable simultaneously.

The databases often create the typical composition sets for phases automatically when data are retrieved. The equilibrium calculations (using the default settings with global minimization) will usually add new composition sets if needed.

Parameters

composition_set – the composition set

Returns

This :class:`MultiDatabaseSystemBuilder` object
without_default_phases()
Rejects all the default phases from both the thermodynamic and the kinetic database, any phase now needs
to be selected manually for the databases.

Returns This MultiDatabaseSystemBuilder object

class tc_python.system.System(system_instance)
Bases: object
A system containing selections for databases, elements, phases etc.

Note: For the defined system, different calculations can be configured and run. Instances of this class should always be created from a SystemBuilder.

Note: The system object is immutable, i.e. it cannot be changed after is has been created. If you want to change the system, you must instead create a new one.

convert_composition(input_composition: Dict[str, float], input_unit: tc_python.utils.ConversionUnit, output_unit: tc_python.utils.ConversionUnit, dependent_component: str = '') → Dict[str, float]
Provides conversion between composition units for any combination of chemical compounds. It is fast because no thermodynamic equilibrium calculation is involved.


Note: It is not required that the chemical compounds are components of the database. The only requirement is that all elements are present in the database.

Parameters
- input_composition – Composition (for example: {“Al2O3”: 25.0, “FeO”: 75.0})
- input_unit – Unit of the input composition
- output_unit – Requested output unit
- dependent_component – The dependent component (optional), for example: “Fe”. If no dependent component is specified the sum of the input composition needs to match 100% / 1

Returns The composition in the requested output unit

get_all_elements_in_databases() → List[str]
Returns the names of all elements present in the selected databases, regardless of the actual selection of elements.

Returns A list of element names

get_all_phases_in_databases() → List[str]
Returns all phase names present in the selected databases, regardless of selected elements, phases etc.

Returns A list of phase names

get_all_species_in_databases() → List[str]
Returns all species names present in the selected databases, regardless of the actual selection of elements, phases, . . .
Returns A list of species names

get_element_object (element_name: str) → tc_python.entities.Element

Returns the Element object of an element. This can be used to obtain detailed information about the element.

Parameters element_name – The element name

Returns A Element: object

get_elements_in_system () → List[str]

Returns the names of all elements present in the selected system.

Note: The list does not contain any elements or components that have been auto-selected by the database(s) in a calculator. Use the get_components() of the calculator object instead to get the complete information.

Returns A list of element names

get_phase_object (phase_name: str) → tc_python.entities.Phase

Returns the Phase object of a phase. This can be used to obtain detailed information about the phase.

Parameters phase_name – The phase name

Returns A Phase: object

get_phases_in_system () → List[str]

Returns all phase names present in the system due to its configuration (selected elements, phases, etc.).

Returns A list of phase names

get_references () → Dict[str, List[str]]

Provides a dictionary with database references per database in the selected system.

Returns The database references

get_species_in_system () → List[str]

Returns the names of all species present in the selected system.

Note: The list does not contain any species or components that have been auto-selected by the database(s) in a calculator. Use the get_components() of the calculator object instead to get the complete information.

Returns The list of species names

get_species_object (species_name: str) → tc_python.entities.Species

Returns the Species object of an species. This can be used to obtain detailed information about the species.

Parameters species_name – The species name

Returns A Species: object

get_system_data () → tc_python.abstract_base.SystemData

Returns the content of the database. This can be used to modify the parameters and functions and to change the current system by using with_system_modifications().
Note: Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb-file.

Returns The system data

```
with_batch_equilibrium_calculation (default_conditions: bool = True,
                                 components: List[str] = []) →
```

Creates a batch-equilibrium calculation (a vectorized equilibrium calculation).

Note: Use this instead of looping if you want to calculate equilibria for a larger number of compositions
and know the conditions in advance. This calculation type has improved performance when calculating a
large number of equilibria when each individual calculations is quick. E.g. when evaluating single phase
properties for thousands of compositions.

Parameters

- **default_conditions** – If True, automatically sets the conditions N=1 and
  P=100000

- **components** – Specify here the components of the system (for example: [AL2O3, ...]),
  only necessary if they differ from the elements. If this option is used, all elements
  of the system need to be replaced by a component.

Returns A new BatchEquilibriumCalculation object

```
with_cct_precipitation_calculation () → tc_python.precipitation.PrecipitationCCTCalculation
```

Creates a CCT diagram calculation.

Returns A new PrecipitationCCTCalculation object

```
with_isothermal_diffusion_calculation () → tc_python.diffusion.DiffusionIsoThermalCalculation
```

Creates an isothermal diffusion calculation.

Returns A new DiffusionIsoThermalCalculation object

```
with_isothermal_precipitation_calculation () → tc_python.precipitation.PrecipitationIsoThermalCalculation
```

Creates an isothermal precipitation calculation.

Returns A new PrecipitationIsoThermalCalculation object

```
with_material_to_material () → tc_python.material_to_material.MaterialToMaterialCalculationContainer
```

Provides access to all Material to Material calculations. The actual calculation needs to be chosen in the
returned object.

Returns A new MaterialToMaterialCalculationContainer object

```
with_non_isothermal_diffusion_calculation () → tc_python.diffusion.DiffusionNonIsoThermalCalculation
```

Creates a non-isothermal precipitation calculation.

Returns A new PrecipitationNonIsoThermalCalculation object

```
with_non_isothermal_precipitation_calculation () → tc_python.precipitation.PrecipitationNonIsoThermalCalculation
```

Creates a non-isothermal precipitation calculation.

Returns A new PrecipitationNonIsoThermalCalculation object
with_phase_diagram_calculation (default_conditions: bool = True, components: List[str] = []) → tc_python.step_or_map_diagrams.PhaseDiagramCalculation

Creates a phase diagram (map) calculation.

Parameters

• **default_conditions** – If True, automatically sets the conditions N=1 and P=100000

• **components** – Specify here the components of the system (for example: `[AL2O3, ...]`), only necessary if they differ from the elements. If this option is used, all elements of the system need to be replaced by a component.

Returns A new PhaseDiagramCalculation object

with_property_diagram_calculation (default_conditions: bool = True, components: List[str] = []) → tc_python.step_or_map_diagrams.PropertyDiagramCalculation

Creates a property diagram (step) calculation.

Parameters

• **default_conditions** – If True, automatically sets the conditions N=1 and P=100000

• **components** – Specify here the components of the system (for example: `[AL2O3, ...]`), only necessary if they differ from the elements. If this option is used, all elements of the system need to be replaced by a component.

Returns A new PropertyDiagramCalculation object

with_property_model_calculation (model: str, path_to_models: str = '', debug_model: bool = False) → tc_python.propertymodel.PropertyModelCalculation

Creates a Property Model calculation.

The parameter **debug_model** is only used when debugging self-developed models.

Parameters

• **model** – The Property Model to be calculated.

• **path_to_models** – The path where the Property Models are installed. If no value is entered, the Property Models folder used by the normal Thermo-Calc application is used.

• **debug_model** – Used when debugging self-developed models.

Returns A new PropertyModelCalculation object

with_scheil_calculation () → tc_python.scheil.ScheilCalculation

Creates a Scheil solidification calculation.

**Warning:** Scheil calculations do not support the `GAS` phase being selected, this means the `GAS` phase must always be deselected in the system if it is present in the database.

Returns A new ScheilCalculation object

with_single_equilibrium_calculation (default_conditions: bool = True, components: List[str] = []) → tc_python.single_equilibrium.SingleEquilibriumCalculation

Creates a single equilibrium calculation.
Parameters

- **default_conditions** – If True, automatically sets the conditions $N=1$ and $P=100000$

- **components** – Specify here the components of the system (for example: \([AL2O3, \ldots]\)), only necessary if they differ from the elements. If this option is used, all elements of the system need to be replaced by a component.

Returns A new `SingleEquilibriumCalculation` object

```python
with_ttt_precipitation_calculation() \rightarrow tc_python.precipitation.PrecipitationTTTCalculation
```

Creates a TTT diagram calculation.

Returns A new `PrecipitationTTTCalculation` object

```python
class tc_python.system.SystemBuilder(system_builder)
Bases: object
```

Used to select databases, elements, phases etc. and create a System object. The system is then used to create calculations.

```python
create_and_select_species(stoichiometry: str)
```

Specify a species from the already entered elements. The stoichiometry of the species is the chemical formula of the species. The created species will also be automatically selected.

**Note:** The elements in the chemical formula are normally separated by stoichiometric numbers. Neither parenthesis “()” nor an underscore “_” is allowed in the chemical formula, while the special combination “/-” or “/+” can be used. Consult the Thermo-Calc database documentation for details about the syntax.

**Parameters** stoichiometry – The stoichiometry of the species

**Returns** This `SystemBuilder` object

```python
deselect_constituent_on_sublattice(phase_name: str, sublattice_no: int, constituent_name_to_deselect: str)
```

Rejects a constituent on a sublattice in a phase in the last specified database only.

**Parameters**

- **phase_name** – The name of the phase
- **sublattice_no** – The number of the sublattice (starting with 1)
- **constituent_name_to_deselect** – The name of the constituent to deselect

**Returns** This `SystemBuilder` object

```python
deselect_phase(phase_name_to_deselect: str)
```

Rejects a phase in the last specified database only.

**Parameters** phase_name_to_deselect – The name of the phase

**Returns** This `SystemBuilder` object

```python
deselect_species(stoichiometry: str)
```

Removes the species from the system.

**Parameters** stoichiometry – The species

**Returns** This `SystemBuilder` object
get_system() \rightarrow tc\_python.system.System

Creates a new System object that is the basis for all calculation types. Several calculation types can be defined later from the object; these are independent.

Returns A new System object

get_system_for_scheil_calculations() \rightarrow tc\_python.system.System

Creates a new System object without gas phases being selected, that is the basis for all calculation types, but its particularly useful for Scheil solidification calculations, where the model does not allow that a gas phase is selected in the system.

Several calculation types can be defined later from the object; these are independent.

Returns A new System object

select_constituent_on_sublattice(phase_name: str, sublattice_no: int, constituent_name_to_select: str)

Selects a constituent on a sublattice in a phase in the last specified database only.

Note: Previously the third parameter constituent_name_to_select had a wrong name, it has been corrected in version 2021b.

Parameters

- phase_name – The name of the phase
- sublattice_no – The number of the sublattice (starting with 1)
- constituent_name_to_select – The name of the constituent to select

Returns This SystemBuilder object

select_database_and_elements(database_name: str, list_of_element_strings: List[str])

Selects a thermodynamic or kinetic database and its selected elements (that will be appended). After that, phases can be selected or unselected.

Parameters

- database_name – The database name, for example “FEDEMO”
- list_of_element_strings – A list of one or more elements as strings, for example [“Fe”, “C”]

Returns This SystemBuilder object

select_phase(phase_name_to_select: str)

Selects a phase in the last specified database only.

Parameters phase_name_to_select – The name of the phase

Returns This SystemBuilder object

select_species(stoichiometry: str)

Adds the species to the system. Up to 1000 species can be defined in a single system.

Parameters stoichiometry – The species

Returns This SystemBuilder object

select_user_database_and_elements(path_to_user_database: str, list_of_element_strings: List[str])

Selects a thermodynamic database which is a user-defined database and select its elements (that will be appended).
Note: By using a r-literal, it is possible to use slashes on all platforms, also on Windows: `select_user_database_and_elements(r"my path/user_db.tdb", ["Fe", "Cr"])`

Otherwise it is required to use **double** backslashes on Windows as separator.

---

Note: On Linux and Mac the path is case-sensitive, also the file ending.

---

**Parameters**

- **path_to_user_database** – The path to the database file (“database”.TDB), defaults to the current working directory. Only the filename is required if the database is located in the same folder as the script.

- **list_of_element_strings** – A list of one or more elements as strings, for example [“Fe”, “C”]

**Returns** This `SystemBuilder` object

```python
with_new_composition_set (composition_set: tc_python.entities.CompositionSet)
```

Used to enter composition sets for a phase. If a phase has a miscibility gap it is necessary to have two composition sets, one for each possible composition that can be stable simultaneously.

**Parameters** composition_set – The composition set

**Returns** This `SystemBuilder` object

```python
without_default_phases ()
```

Rejects all default phases in the last specified database only, any phase needs now to be selected manually for that database.

**Returns** This `SystemBuilder` object

---

### 5.3 Module “entities”

```python
class tc_python.entities.CompositionSet (phase_name: str)
```

Used by the method `tc_python.system.SystemBuilder.with_new_composition_set()` to enter two or more composition sets for a phase.

**Parameters** phase_name – The name of the phase for which a new composition set is required

```python
set_major_constituents_for_sublattice (sublattice_index: int, major_constituents: List[str])
```

Specify the new major constituent(s) for the sublattice.

**Default:** If not specified, a default is automatically chosen based on the specified composition set.

**Note:** This is useful in order to make calculations converge faster and more easily (because it may simplify giving start values when calculating the equilibrium as those phases with miscibility gaps should have different major constituents for each composition set). The databases often set major constituents for several phases automatically when the data is retrieved.
Parameters

- **sublattice_index** – Index of the sublattice to set the major constituents for (starting with 1)
- **major_constituents** – Optional list of the major constituents, which must be selected from the phase constitution of the current system.

Returns This `CompositionSet` object

```python
class tc_python.entities.Element(element)
    Bases: object

    Represents an element, making detailed information about the element accessible.

    get_enthalpy() → float
        Returns the enthalpy of the element at 298 K, part of the stable element reference state (SER).
        Returns The enthalpy [J]

    get_entropy_diff_0_to_298k() → float
        Returns the entropy difference 0 - 298 K of the element, part of the stable element reference state (SER).
        Returns The entropy difference 0 - 298 K [J/K]

    get_molar_mass() → float
        Returns the molar mass of the element.
        Returns The molar mass [g/mol]

    get_name() → str
        Returns the name of the element.
        Returns The element name

    get_stable_element_reference() → str
        Returns the stable element reference (i.e. the stable phase at 298.15 K and 1 bar, reference for all element thermodynamic data).
        Returns The name of the stable element reference

    is_interstitial() → bool
        Returns if the element is interstitial.

    is_special() → bool
        Returns if the element is special (i.e. vacancies (VA) and electrons (denoted either as /- in gaseous, liquid or solid phases, or ZE in an aqueous solution phase)).

    is_valid() → bool
        Returns if the element is valid. Non-valid elements are represented by an empty name.
```

**Note:** In the diffusion simulations (DICTRA), the assumption that the volume is carried by the substitutional elements only is applied. The interstitial elements are assumed to have zero molar volumes.
class tc_python.entities.Phase(\textit{phase})

Bases: \texttt{object}

Represents a phase, making detailed information about the phase accessible.

\textbf{get\_name}() \rightarrow \texttt{str}

Returns the name of the phase.

\textbf{Returns} The phase name

\textbf{get\_species}() \rightarrow \texttt{Set[tc_python.entities.Species]}

Returns the species of the phase.

\textbf{Returns} A set containing the species

\textbf{get\_species\_for\_composition\_profile}() \rightarrow \texttt{Set[tc_python.entities.Species]}

Returns all species that need to be defined in a composition profile of the phase for diffusion simulations - except for one species that needs to be the dependent species.

\textbf{Note:} In a composition profile of a phase for diffusion simulations it is necessary to specify all non-stoichiometric and non-special species. In case of a DILUTE diffusion model, the database enforces the choice of a certain dependent species.

\textbf{Returns} Set with the species

\textbf{get\_sublattices}() \rightarrow \texttt{List[tc_python.entities.Sublattice]}

Returns the sublattices of the phase in a well-defined contiguous order.

\textbf{Returns} A list containing the \texttt{Sublattice} objects

\textbf{get\_type}() \rightarrow \texttt{tc_python.entities.PhaseType}

Returns the type of the phase (liquid, ionic liquid, solid, gas).

\textbf{Returns} The type of a phase

\textbf{has\_diffusion\_data}() \rightarrow \texttt{bool}

Returns if diffusion data exists for the phase.

\textbf{Returns} If diffusion data exists for the phase

\textbf{has\_molar\_volume\_data}() \rightarrow \texttt{bool}

Returns if molar volume data exists for the phase.

\textbf{Returns} If molar volume data exists for the phase

\textbf{is\_dilute\_diffusion\_model}() \rightarrow \texttt{bool}

Returns if diffusion is described using the DILUTE model for the phase. This will always return \textit{False} if no diffusion data is available.

\textbf{Returns} If the DILUTE model is used

\textbf{is\_gas}() \rightarrow \texttt{bool}

Returns if the phase is a gas phase.

\textbf{Returns} If the phase is a gas phase

\textbf{is\_ionic\_liquid}() \rightarrow \texttt{bool}

Returns if the phase is an ionic liquid phase.

\textbf{Returns} If the phase is an ionic liquid phase
is_liquid() → bool
    Returns if the phase is a liquid or ionic liquid phase.
    **Returns** If the phase is a liquid phase

is_solid() → bool
    Returns if the phase is a solid phase.
    **Returns** If the phase is a solid phase

class tc_python.entities.PhaseType(value)
    Bases: enum.Enum
    The type of a phase.
    GAS = 0
        Gas phase.
    IONIC_LIQUID = 2
        Ionic liquid phase.
    LIQUID = 1
        Liquid phase.
    SOLID = 3
        Solid phase.

class tc_python.entities.Species(species)
    Bases: object
    Represents a species, making detailed information about the species accessible.
    get_all_elements() → List[Tuple[tc_python.entities.Element, float]]
        Returns all the elements that the species is composed of.
        **Returns** List of all elements of the species and their stoichiometry
    get_charge() → int
        Returns the charge of the species.
        **Returns** The charge of the species
    get_name() → str
        Returns the name of the species.
        **Returns** The species name
    is_element() → bool
        Returns if the species actually represents an element.
        **Returns** If the species represents an element
    is_interstitial() → bool
        Returns if the species is interstitial.
        **Returns** If the species is interstitial

Note: In the diffusion simulations (DICTRA), the assumption that the volume is carried by the substitutional elements only is applied. The interstitial elements are assumed to have zero molar volumes.
is_special() → bool
Returns if the species is special (i.e. vacancies (VA) and electrons (denoted either as /- in gaseous, liquid or solid phases, or ZE in an aqueous solution phase)).

Returns If the species is special

is_valid() → bool
Returns if the species is valid. Non-valid species are represented by an empty name.

Returns If the species is valid

to_element() → tc_python.entities.Element
Returns the Element representation of the species - if the species actually represents an element.

Returns The Element object

class tc_python.entities.Sublattice(sublattice)
Bases: object
Represents a sublattice of a phase.

get_constituents() → Set[tc_python.entities.Species]
Returns the constituents of the sublattice.

Returns A set containing the constituents

get_nr_of_sites() → float
Returns the number of sites in the sublattice.

Returns A float number

5.4 Module “server”

class tc_python.server.LoggingPolicy(value)
Bases: enum.Enum
Logging policy that determines how the TC-Python logs are presented to the user.

FILE = 1
Logging to a file.

NONE = 2
No logging at all.

SCREEN = 0
Logging to the screen.

class tc_python.server.ResultLoader(result_loader)
Bases: object
Contains methods for loading results from previously done calculations.

diffusion(path: str) → tc_python.diffusion.DiffusionCalculationResult
Loads a DiffusionCalculationResult from disc.

Parameters path – path to the folder where result was previously saved.

Returns A new DiffusionCalculationResult object which later can be used to get specific values from the calculated result

phase_diagram(path: str) → tc_python.step_or_map_diagrams.PhaseDiagramResult
Loads a PhaseDiagramResult from disc.
Parameters path – path to the folder where result was previously saved.

Returns A new PhaseDiagramResult object which later can be used to get specific values from the calculated result

`precipitation_TTT_or_CCT` *(path: str) → tc_python.precipitation.PrecipitationCalculationTTTorCCTResult*

Loads a PrecipitationCalculationTTTorCCTResult from disc.

Parameters path – path to the folder where result was previously saved.

Returns A new PrecipitationCalculationTTTorCCTResult object which later can be used to get specific values from the calculated result

`precipitation_single` *(path: str) → tc_python.precipitation.PrecipitationCalculationSingleResult*

Loads a PrecipitationCalculationSingleResult from disc.

Parameters path – path to the folder where result was previously saved.

Returns A new PrecipitationCalculationSingleResult object which later can be used to get specific values from the calculated result

`property_diagram` *(path: str) → tc_python.step_or_map_diagrams.PropertyDiagramResult*

Loads a PropertyDiagramResult from disc.

Parameters path – path to the folder where result was previously saved.

Returns A new PropertyDiagramResult object which later can be used to get specific values from the calculated result

`property_model` *(path: str) → tc_python.propertymodel.PropertyModelResult*

Loads a PropertyModelResult from disc.

Parameters path – path to the folder where result was previously saved.

Returns A new PropertyModelResult object which later can be used to get specific values from the calculated result

`scheil` *(path: str) → tc_python.scheil.ScheilCalculationResult*

Loads a ScheilCalculationResult from disc.

Parameters path – path to the folder where result was previously saved.

Returns A new ScheilCalculationResult object which later can be used to get specific values from the calculated result

`single_equilibrium` *(path: str) → tc_python.single_equilibrium.SingleEquilibriumResult*

Loads a SingleEquilibriumResult from disc.

Parameters path – path to the folder where result was previously saved.

Returns A new SingleEquilibriumResult object which later can be used to get specific values from the calculated result

`class tc_python.server.SetUp(debug_logging=False)`

Bases: object

Starting point for all calculations.

---

**Note:** This class exposes methods that have no precondition, it is used for choosing databases and elements.

`disable_caching()`

A previously set cache folder is no longer used.
Note: Within the session, caching is activated and used through the default temporary directory.

**Returns** This SetUp object

**get_database_info** *(database_short_name: str) → str*
Obtains the short information available for the specified database.

**Parameters** database_short_name – The name of the database (i.e. “FEDEMO”, …)

**Returns** The short information about the database

**get_database_path_on_disk** *(database_short_name: str) → str*
Obtains the path to the database file on disk. TCPATH is a placeholder for the root path of the used Thermo-Calc installation.

Note: Encrypted databases (*.TDC) cannot be edited.

**Parameters** database_short_name – The name of the database (i.e. “FEDEMO”, …)

**Returns** The path to the database on disk

**get_databases** () → List[str]
Obtains the short names of all databases available in the used Thermo-Calc installation.

Note: Only databases with a valid license are listed.

**Returns** List of the available databases

**get_property_models** *(path_to_models: str = '') → Set[str]*
Lists the names of all Property Models in the specified directory.

If the directory is not specified, the Property Model folder used by the normal Thermo-Calc application is used.

**Parameters** path_to_models – The path where the Property Models are installed. If no value is entered, the Property Model folder used by the normal Thermo-Calc application is used.

**Returns** Set containing all Property Model names

**load_result_from_disk** () → tc_python.server.ResultLoader
Loads a previously calculated result from disk.

Note: This only works for results created by calling one of the save_result() methods on a Result class created from a calculation.

**Returns** A new ResultLoader object

**select_database_and_elements** *(database_name: str, list_of_elements: List[str]) → tc_python.system.SystemBuilder*
Selects a first thermodynamic or kinetic database and selects the elements in it.
Parameters

- **database_name** – The name of the database, for example “FEDEMO”

- **list_of_elements** – The list of the selected elements in that database, for example [“Fe”, “C”]

Returns A new `SystemBuilder` object

```python
def select_thermodynamic_and_kinetic_databases_with_elements(thermodynamic_db_name: str, kinetic_db_name: str, list_of_elements: List[str]) -> tc_python.system.MultiDatabaseSystemBuilder:
    Selects the thermodynamic and kinetic database at once, guarantees that the databases are added in the correct order. Further rejection or selection of phases applies to both databases.
```

Parameters

- **thermodynamic_db_name** – The thermodynamic database name, for example “FEDEMO”

- **kinetic_db_name** – The kinetic database name, for example “MFEDEMO”

- **list_of_elements** – The list of the selected elements in that database, for example [“Fe”, “C”]

Returns A new `MultiDatabaseSystemBuilder` object

```python
def select_user_database_and_elements(path_to_user_database: str, list_of_elements: List[str]) -> tc_python.system.SystemBuilder:
    Selects a user-defined database and selects the elements in it.
```

Note: By using a `r`-literal, it is possible to use slashes on all platforms, also on Windows: `select_user_database_and_elements(r"my path/user_db.tdb", ["Fe", "Cr"])`

Otherwise it is required to use `double` backslashes on Windows as separator.

Note: On Linux and Mac the path is case-sensitive, also the file ending.

Parameters

- **path_to_user_database** – The path to the database file (“database”.TDB), defaults to the current working directory. Only filename is required if the database is located in the same folder as the script.

- **list_of_elements** – The list of the selected elements in that database, for example [“Fe”, “C”]

Returns A new `SystemBuilder` object

```python
def set_cache_folder(path: str = '', precision_for_floats: int = 12):
    Sets a folder where results from calculations and state of systems are saved. If at any time a calculation is run which has the exact same setting as a previous, the calculation is not re-run. The result is instead loaded from this folder.
```
Note: The same folder can be used in several scripts, and it can even be shared between different users. It can be a network folder.

Parameters

- **path** – path to the folder where results should be stored. It can be relative or absolute.
- **precision_for_floats** – The number of significant figures used when comparing if the calculation has the same setting as a previous.

Returns This SetUp object

```python
def set_ges_version(version: int = 6)
    """Setting the version of the Gibbs Energy System (GES)."""
    Parameters version – The GES-version (currently version 5 or 6)
    Returns This SetUp object
```

```python
def set_log_level_to_debug()
    """Sets log level to DEBUG"""
    Returns This SetUp object
```

```python
def set_log_level_to_info()
    """Sets log level to INFO"""'
    Returns This SetUp object
```

`class tc_python.server.TCython(logging_policy=<LoggingPolicy.SCREEN: 0>, log_file=None, debug_mode=False, debug_logging=False, do_throw_on_backend_hard_crash=True, port_number=0)`

Bases: object

Starting point of the API. Typical syntax:

```python
with TCython() as session:
    session.select_database_and_elements(...)  
```

Note: Each usage of `with TCython()` causes significant overhead (starting a new process, stopping the old one, cleaning up the temporary disk space). Usually it is recommendable to call `with TCython()` only once for each process, even if working in a loop. Instead you should pass the session or calculator object into the loop and use them there.

If necessary, beginning from version 2019a it is however possible to call `with TCython` safely multiple times.

```python
def start_api_server(logging_policy=<LoggingPolicy.SCREEN: 0>, log_file=None, debug_mode=False, is_unittest=False, do_throw_on_backend_hard_crash=True, port_number=0)
    """Starts a process of the API server and sets up the socket communication with it."""
    Parameters
        - **logging_policy** – Determines if the TC-Python log output is sent to the screen (LoggingPolicy.SCREEN), to file (LoggingPolicy.FILE) or nothing is logged at all (LoggingPolicy.NONE) Default: LoggingPolicy.SCREEN. Note that the log-handlers can also be adapted through the `tc_python.LOGGER` object at any time.
```

194 Chapter 5. API Reference
• **log_file** – The log-file relative to the current path or absolute, only relevant if `logging_policy=LoggingPolicy.FILE`. Log-output will be appended.

• **debug_mode** – If `True` it is tried to open a connection to an already running API-server. This is only used for debugging the API itself.

• **is_unittest** – Should be `True` if called by a unit test, only to be used internally for development.

• **do_throw_on_backend_hard_crash** – If `True` an `UnrecoverableCalculationException` will be thrown if the Java-backend crashes hard, if `False` the application will simply crash with a FORTRAN-stacktrace. If `True` the exception can be caught outside of the ‘with’-clause and the application can continue, if `False` more information about the error is shown by the stacktrace.

• **port_number** – The port number for the communication with the Java-backend server. This is not required to be changed by normal users.

**Warning:** Most users should use **TCPython** using a `with`-statement for automatic management of the resources (network sockets and temporary files). If you anyway need to use that method, make sure to call `stop_api_server()` in any case using the try-finally-pattern.

```python
tc_python.server.start_matlab_server(logging_policy=<LoggingPolicy.SCREEN: 0>, log_file=None, debug_mode=False, is_unittest=False, do_throw_on_backend_hard_crash=True, port_number=0)
```

```python
tc_python.server.stop_api_server(gateway_id: str = None)
```

Clears all resources used by the session (i.e. shuts down the API server and deletes all temporary files). The disk usage of temporary files might be significant.

**Warning:** Call this method only if you used `start_api_server()` initially. It should never be called when the API has been initialized in a `with`-statement using **TCPython**.

### 5.5 Module “quantity_factory”

**class** `tc_python.quantity_factory.DiffusionQuantity`

**Bases:** `tc_python.quantity.AbstractQuantity`

Factory class providing quantities used for defining diffusion simulations and their results.

**Note:** In this factory class only the most common quantities are defined, you can always use the **Console Mode** syntax strings in the respective methods as an alternative (for example: “NPM(*)”).

```python
classmethod activity_of_component(component: str, use_ser: bool = False) →
tc_python.quantity.ActivityOfComponent
```

Creates a quantity representing the activity of a component.

**Parameters**

• **component** – The name of the component, use `ALL_COMPONENTS` to choose all components
• **use_ser** – Use Stable-Element-Reference(SER). The user-defined reference state is be used if this setting is set to `False`.

**Returns** A new `ActivityOfComponent` object.

```python
classmethod chemical_diffusion_coefficient(  
    phase: str,  
    diffusing_element: str,  
    gradient_element: str,  
    reference_element: str) → tc_python.quantity.ChemicalDiffusionCoefficient
```

Creates a quantity representing the chemical diffusion coefficient of a phase \( \text{[m}^2/\text{s}] \).

**Parameters**

• **phase** – The name of the phase

• **diffusing_element** – The diffusing element

• **gradient_element** – The gradient element

• **reference_element** – The reference element (for example “Fe” in a steel)

**Returns** A new `ChemicalDiffusionCoefficient` object.

```python
classmethod chemical_potential_of_component(  
    component: str,  
    use_ser: bool = False) → tc_python.quantity.ChemicalPotentialOfComponent
```

Creates a quantity representing the chemical potential of a component \( \text{[J]} \).

**Parameters**

• **component** – The name of the component, use `ALL_COMPONENTS` to choose all components

• **use_ser** – Use Stable-Element-Reference(SER). The user-defined reference state is used if this setting is set to `False`.

**Returns** A new `ChemicalPotentialOfComponent` object.

```python
classmethod distance(region: str = 'All') → tc_python.quantity.Distance
```

Creates a quantity representing the distance \( \text{[m]} \).

**Parameters** **region** – The name of the region or `All` to choose global.

```python
classmethod intrinsic_diffusion_coefficient(  
    phase: str,  
    diffusing_element: str,  
    gradient_element: str,  
    reference_element: str) → tc_python.quantity.IntrinsicDiffusionCoefficient
```

Creates a quantity representing the intrinsic diffusion coefficient of a phase \( \text{[m}^2/\text{s}] \).

**Parameters**

• **phase** – The name of the phase

• **diffusing_element** – The diffusing element

• **gradient_element** – The gradient element

• **reference_element** – The reference element (for example “Fe” in a steel)

**Returns** A new `IntrinsicDiffusionCoefficient` object.

```python
```

Creates a quantity representing \( L^2 \) of a phase \( \text{[m}^2/\text{s}] \).

**Parameters**
• **phase** – The name of the phase

• **diffusing_element** – The diffusing element

• **gradient_element** – The gradient element

• **reference_element** – The reference element (for example “Fe” in a steel)

Returns A new `Lbis` object.

```python
classmethod mass_fraction_of_a_component (component: str) →
tc_python.quantity.MassFractionOfAComponent
```

Creates a quantity representing the mass fraction of a component.

Parameters **component** – The name of the component or `ALL_COMPONENTS` to choose all components

Returns A new `MassFractionOfAComponent` object.

```python
classmethod mass_fraction_of_a_phase (phase: str) →
tc_python.quantity.MassFractionOfAPhase
```

Creates a quantity representing the mass fraction of a phase.

Parameters **phase** – The name of the phase or `ALL_PHASES` to choose all phases.

Returns A new `MassFractionOfAPhase` object.

```python
classmethod mobility_of_component_in_phase (phase: str, component: str) →
tc_python.quantity.MobilityOfComponentInPhase
```

Creates a quantity representing the mobility of a component in a phase \([\text{m}^2/\text{Js}]\).

Parameters

• **phase** – The name of the phase

• **component** – The name of the component

Returns A new `MobilityOfComponentInPhase` object.

```python
classmethod mole_fraction_of_a_component (component: str) →
tc_python.quantity.MoleFractionOfAComponent
```

Creates a quantity representing the mole fraction of a component.

Parameters **component** – The name of the component or `ALL_COMPONENTS` to choose all components

Returns A new `MoleFractionOfAComponent` object.

```python
classmethod mole_fraction_of_a_phase (phase: str) →
tc_python.quantity.MoleFractionOfAPhase
```

Creates a quantity representing the mole fraction of a phase.

Parameters **phase** – The name of the phase or `ALL_PHASES` to choose all phases.

Returns A new `MoleFractionOfAPhase` object.

```python
classmethod position_of_lower_boundary_of_region (region: str) →
tc_python.quantity.PositionOfLowerBoundaryOfRegion
```

Creates a quantity representing the position of lower boundary of a region \([\text{m}]\).

Parameters **region** – The name of the region

Returns A new `PositionOfLowerBoundaryOfRegion` object.

```python
classmethod position_of_upper_boundary_of_region (region: str) →
tc_python.quantity.PositionOfUpperBoundaryOfRegion
```

Creates a quantity representing the position of upper boundary of a region \([\text{m}]\).

5.5. Module “quantity_factory”
Parameters `region` – The name of the region

Returns A new `PositionOfUpperBoundaryOfRegion` object.

classmethod `temperature()` → `tc_python.quantity.Temperature`

Creates a quantity representing the temperature [K].

Returns A new `Temperature` object.


Creates a quantity representing thermodynamic factor of a phase.

Parameters

- `phase` – The name of the phase
- `diffusing_element` – The diffusing element
- `gradient_element` – The gradient element
- `reference_element` – The reference element (for example “Fe” in a steel)

Returns A new `ThermoDynamicFactor` object.

classmethod `time()` → `tc_python.quantity.Time`

Creates a quantity representing the time [s].

classmethod `total_mass_fraction_of_component(component: str)` → `tc_python.quantity.TotalMassFractionOfComponent`

Creates a quantity representing the total mass fraction of a component.

Parameters `component` – The name of the component

Returns A new `TotalMassFractionOfComponent` object.

classmethod `total_mass_fraction_of_component_in_phase(phase: str, component: str)` → `tc_python.quantity.TotalMassFractionOfComponentInPhase`

Creates a quantity representing the total mass fraction of a component in a phase.

Parameters

- `phase` – The name of the phase
- `component` – The name of the component

Returns A new `TotalMassFractionOfComponentInPhase` object.

classmethod `total_mass_fraction_of_phase(phase: str)` → `tc_python.quantity.TotalMassFractionOfPhase`

Creates a quantity representing the total mass fraction of a phase.

Parameters `phase` – The name of the phase.

Returns A new `TotalMassFractionOfPhase` object.

classmethod `total_mole_fraction_of_component(component: str)` → `tc_python.quantity.TotalMoleFractionOfComponent`

Creates a quantity representing the total mole fraction of a component.

Parameters `component` – The name of the component

Returns A new `TotalMoleFractionOfComponent` object.
classmethod total_mole_fraction_of_component_in_phase(phase: str, component: str) → tc_python.quantity.TotalMoleFractionOfComponentInPhase

Creates a quantity representing the total mole fraction of a component in a phase.

Parameters

- **phase** – The name of the phase
- **component** – The name of the component

Returns A new TotalMoleFractionOfComponentInPhase object.

classmethod total_volume_fraction_of_phase(phase: str) → tc_python.quantity.TotalVolumeFractionOfPhase

Creates a quantity representing the total volume fraction of a phase.

Parameters **phase** – The name of the phase.

Returns A new TotalVolumeFractionOfPhase object.

classmethod tracer_diffusion_coefficient(phase: str, diffusing_element: str) → tc_python.quantity.TracerDiffusionCoefficient

Creates a quantity representing tracer diffusion coefficient of a phase [m^2/s].

Parameters

- **phase** – The name of the phase
- **diffusing_element** – The diffusing element

Returns A new TracerDiffusionCoefficient object.

classmethod u_fraction_of_a_component(component: str) → tc_python.quantity.UFractionOfAComponent

Creates a quantity representing the u-fraction of a component.

Parameters **component** – The name of the component

Returns A new UFractionOfAComponent object.

classmethod user_defined_function(expression: str) → tc_python.quantity.Function

Creates a quantity representing a user-defined function.

Parameters **expression** – The function expression

Returns A new Function object

classmethod velocity_of_lower_boundary_of_region(region: str) → tc_python.quantity.VelocityOfLowerBoundaryOfRegion

Creates a quantity representing the velocity of lower boundary of a region [m/s].

Parameters **region** – The name of the region

Returns A new VelocityOfLowerBoundaryOfRegion object.

classmethod velocity_of_upper_boundary_of_region(region: str) → tc_python.quantity.VelocityOfUpperBoundaryOfRegion

Creates a quantity representing the velocity of upper boundary of a region [m/s].

Parameters **region** – The name of the region

Returns A new VelocityOfUpperBoundaryOfRegion object.

classmethod width_of_region(region: str) → tc_python.quantity.Function

Creates a quantity representing the width of a region [m].

Parameters **region** – The name of the region

5.5. Module “quantity_factory”
Returns A new WidthOfRegion object.

```python
class tc_python.quantity_factory.IndependentVariable
    Bases: tc_python.quantity.AbstractQuantity

Factory class providing quantities used for defining the independent variable in general diffusion result querying.

    classmethod distance(region: str = 'All') → tc_python.quantity.Distance
    Creates an independent variable representing the distance [m].
    Returns A new Distance object

    classmethod time() → tc_python.quantity.Time
    Creates an independent variable representing the time [s].
    Returns A new Time object
```

```python
class tc_python.quantity_factory.PlotCondition
    Bases: tc_python.quantity.AbstractQuantity

Factory class providing quantities used for defining the plot condition in general diffusion result querying.

    Note: In this factory class only the most common quantities are defined, you can always use the Console Mode syntax strings in the respective methods as an alternative (for example: “time last”).

    classmethod distance(distancepoint: float, region: str = 'All') → tc_python.quantity.DistanceCondition
    Creates a plot condition representing the distance [m].
    Change in version 2019b: Mandatory parameter distancepoint added
    Parameters
        • distancepoint – The distance from the lower interface of the region
        • region – The name of the region or All to choose global.
    Returns A new DistanceCondition object

    classmethod integral() → tc_python.quantity.IntegralCondition
    Creates an integral plot condition.
    Returns A new IntegralCondition object

    classmethod interface(region: str, interface_position: tc_python.utils.InterfacePosition) → tc_python.quantity.InterfaceCondition
    Creates a plot condition representing an interface between two regions.
    Parameters
        • region – The name of the region used for defining the interface
        • interface_position – The position of the interface relative to that region (lower or upper)
    Returns A new InterfaceCondition object

    classmethod time(timepoint: Union[float, str] = 'Last') → tc_python.quantity.TimeCondition
    Creates a plot condition representing the time [s].
    Change in version 2019b: Lists of timepoints are no longer supported
    Parameters timepoint – The timepoint. Optionally “Last” can be used for the end of the simulation
Returns A new TimeCondition object

class tc_python.quantity_factory.ScheilQuantity
    Bases: tc_python.quantity.AbstractQuantity

Factory class providing quantities used for defining a Scheil calculation result (tc_python.scheil.ScheilCalculationResult).

classmethod apparent_heat_capacity_per_gram() → tc_python.quantity.ApparentHeatCapacityPerGram
    Creates a quantity representing the apparent heat capacity [J/g/K].
    Returns A new ApparentHeatCapacityPerGram object.

classmethod apparent_heat_capacity_per_mole() → tc_python.quantity.ApparentHeatCapacityPerMole
    Creates a quantity representing the apparent heat capacity [J/mol/K].
    Returns A new ApparentHeatCapacityPerMole object.

classmethod apparent_volumetric_thermal_expansion_coefficient() → tc_python.quantity.ApparentVolumetricThermalExpansionCoefficient
    Creates a quantity representing the apparent volumetric thermal expansion coefficient of the system [1/K].
    Returns A new ApparentVolumetricThermalExpansionCoefficient object.

classmethod composition_of_phase_as_mole_fraction(phase: str, component: str) → tc_python.quantity.CompositionOfPhaseAsMoleFraction
    Creates a quantity representing the composition of a phase [mole-fraction].
    Parameters
        • phase – The name of the phase, use ALL_PHASES to choose all stable phases
        • component – The name of the component, use ALL_COMPONENTS to choose all components
    Returns A new CompositionOfPhaseAsMoleFraction object.

classmethod composition_of_phase_as_weight_fraction(phase: str, component: str) → tc_python.quantity.CompositionOfPhaseAsWeightFraction
    Creates a quantity representing the composition of a phase [weight-fraction].
    Parameters
        • phase – The name of the phase, use ALL_PHASES to choose all stable phases
        • component – The name of the component, use ALL_COMPONENTS to choose all components
    Returns A new CompositionOfPhaseAsWeightFraction object.

classmethod density_of_phase(phase: str) → tc_python.quantity.DensityOfPhase
    Creates a quantity representing the average density of a phase [g/cm^3].
    Parameters phase – The name of the phase or ALL_PHASES to choose all phases
    Returns A new DensityOfPhase object.

classmethod density_of_solid_phase(phase: str) → tc_python.quantity.DensityOfSolidPhase
    Creates a quantity representing the average density of a solid phase [g/cm^3].

Note: Deprecated in version 2022a: This quantity has been renamed to density_of_phase(). It will be removed in release 2023a.
Parameters **phase** – The name of the phase or **ALL_PHASES** to choose all solid phases

**Returns** A new `DensityOfSolidPhase` object.

classmethod `density_of_system()` → `tc_python.quantity.DensityOfSystem`

Creates a quantity representing the average density of the system [g/cm^3].

**Returns** A new `DensityOfSystem` object.

classmethod `distribution_of_component_of_phase` (**phase**: str, **component**: str) →

`tc_python.quantity.DistributionOfComponentOfPhase`

Creates a quantity representing the (molar) fraction of the specified component being present in the specified phase compared to the overall system [-]. This corresponds to the degree of segregation to that phase.

**Parameters**

- **phase** – The name of the phase
- **component** – The name of the component

**Returns** A new `DistributionOfComponentOfPhase` object.

classmethod `heat_per_gram()` → `tc_python.quantity.HeatPerGram`

Creates a quantity representing the total heat release from the liquidus temperature down to the current temperature [J/g].

**Note:** The total or apparent heat release during the solidification process consists of two parts: one is the so-called latent heat, i.e. heat due to the liquid -> solid phase transformation (`latent_heat_per_mole()` and `latent_heat_per_gram()`), and the other is the heat related to the specific heat of liquid and solid phases (`heat_per_mole()` and `heat_per_gram()`).

**Returns** A new `HeatPerGram` object.

classmethod `heat_per_mole()` → `tc_python.quantity.HeatPerMole`

Creates a quantity representing the total heat release from the liquidus temperature down to the current temperature [J/mol].

**Note:** The total or apparent heat release during the solidification process consists of two parts: one is the so-called latent heat, i.e. heat due to the liquid -> solid phase transformation (`latent_heat_per_mole()` and `latent_heat_per_gram()`), and the other is the heat related to the specific heat of liquid and solid phases (`heat_per_mole()` and `heat_per_gram()`).

**Returns** A new `HeatPerMole` object.

classmethod `latent_heat_per_gram()` → `tc_python.quantity.LatentHeatPerGram`

Creates a quantity representing the cumulated latent heat release from the liquidus temperature down to the current temperature [J/g].

**Note:** The total or apparent heat release during the solidification process consists of two parts: one is the so-called latent heat, i.e. heat due to the liquid -> solid phase transformation (`latent_heat_per_mole()` and `latent_heat_per_gram()`), and the other is the heat related to the specific heat of liquid and solid phases (`heat_per_mole()` and `heat_per_gram()`).

**Returns** A new `LatentHeatPerGram` object.
classmethod latent_heat_per_mole() → tc_python.quantity.LatentHeatPerMole
Creates a quantity representing the cumulated latent heat release from the liquidus temperature down to
the current temperature [J/mol].

Note: The total or apparent heat release during the solidification process consists of two parts: one is the so-called latent heat, i.e. heat due to the liquid -> solid phase transformation (latent_heat_per_mole() and latent_heat_per_gram()), and the other is the heat related to the specific heat of liquid and solid phases (heat_per_mole() and heat_per_gram()).

Returns A new LatentHeatPerMole object.

classmethod mass_fraction_of_a_solid_phase(phase: str) → tc_python.quantity.MassFractionOfASolidPhase
Creates a quantity representing the mass fraction of a solid phase.

Parameters phase – The name of the phase or ALL_PHASES to choose all solid phases

Returns A new MassFractionOfASolidPhase object.

classmethod mass_fraction_of_all_liquid() → tc_python.quantity.MassFractionOfAllLiquid
Creates a quantity representing the total mass fraction of all the liquid phase.

Returns A new MassFractionOfAllLiquid object.

classmethod mass_fraction_of_all_solid_phases() → tc_python.quantity.MassFractionOfAllSolidPhases
Creates a quantity representing the total mass fraction of all solid phases.

Returns A new MassFractionOfAllSolidPhase object.

classmethod molar_volume_of_phase(phase: str) → tc_python.quantity.MolarVolumeOfPhase
Creates a quantity representing the molar volume of a phase [m^3/mol].

Parameters phase – The name of the phase or ALL_PHASES to choose all phases

Returns A new MolarVolumeOfPhase object.

classmethod molar_volume_of_system() → tc_python.quantity.MolarVolumeOfSystem
Creates a quantity representing the molar volume of the system [m^3/mol].

Returns A new MolarVolumeOfSystem object.

classmethod mole_fraction_of_a_solid_phase(phase: str) → tc_python.quantity.MoleFractionOfASolidPhase
Creates a quantity representing the molar fraction of a solid phase.

Parameters phase – The name of the phase or ALL_PHASES to choose all solid phases

Returns A new MoleFractionOfASolidPhase object.

classmethod mole_fraction_of_all_liquid() → tc_python.quantity.MoleFractionOfAllLiquid
Creates a quantity representing the total molar fraction of all the liquid phase.

Returns A new MoleFractionOfAllLiquid object.

classmethod mole_fraction_of_all_solid_phases() → tc_python.quantity.MoleFractionOfAllSolidPhases
Creates a quantity representing the total molar fraction of all solid phases.

Returns A new MoleFractionOfAllSolidPhases object.
classmethod site_fraction_of_component_in_phase(phase: str, component: str, sub_lattice_ordinal_no: int = 0) -> tc_python.quantity.SiteFractionOfComponentInPhase

Creates a quantity representing the site fractions [-].

Parameters

- **phase** – The name of the phase, use ALL_PHASES to choose all stable phases
- **component** – The name of the component, use ALL_COMPONENTS to choose all components
- **sub_lattice_ordinal_no** – The ordinal number (i.e. 1, 2, ...) of the sublattice of interest, use None to choose all sublattices

Note: Detailed information about the sublattices can be obtained by getting the Phase object of a phase from the System object using tc_python.system.System.get_phase_in_system. For each phase the sublattices are obtained by using tc_python.system.Phase.get_sublattices. The order in the returned list is equivalent to the sublattice ordinal number expected, but note that the ordinal numbers start with 1.

Returns A new SiteFractionOfComponentInPhase object.

classmethod temperature() -> tc_python.quantity.Temperature

Creates a quantity representing the temperature [K].

Returns A new Temperature object.

class tc_python.quantity_factory.ThermodynamicQuantity
Bases: tc_python.quantity.AbstractQuantity

Factory class providing quantities used for defining equilibrium calculations (single equilibrium, property and phase diagrams, ...) and their results.

Note: In this factory class only the most common quantities are defined, you can always use the Console Mode syntax strings in the respective methods as an alternative (for example: “NPM(*)”).

classmethod activity_of_component(component: str, use_ser: bool = False) -> tc_python.quantity.ActivityOfComponent

Creates a quantity representing the activity of a component [-].

Parameters

- **component** – The name of the component, use ALL_COMPONENTS to choose all components
- **use_ser** – Use Stable-Element-Reference(SER). The user-defined reference state is used if this setting is set to False.

Returns A new ActivityOfComponent object.


Creates a quantity representing the chemical diffusion coefficient of a phase [m²/s].
Parameters

- **phase** – The name of the phase
- **diffusing_element** – The diffusing element
- **gradient_element** – The gradient element
- **reference_element** – The reference element (for example “Fe” in a steel)

**Returns** A new ChemicalDiffusionCoefficient object.

```python
classmethod chemical_potential_of_component(component: str, use_ser: bool = False) ->
tc_python.quantity.ChemicalPotentialOfComponent
```

Creates a quantity representing the chemical potential of a component [J].

Parameters

- **component** – The name of the component, use `ALL_COMPONENTS` to choose all components
- **use_ser** – Use Stable-Element-Reference(SER). The user-defined reference state is used if this setting is set to `False`.

**Returns** A new ChemicalPotentialOfComponent object.

```python
classmethod composition_of_phase_as_mole_fraction(phase: str, component: str = 'All') ->
tc_python.quantity.CompositionOfPhaseAsMoleFraction
```

Creates a quantity representing the composition of a phase [mole-fraction].

Parameters

- **phase** – The name of the phase, use `ALL_PHASES` to choose all stable phases
- **component** – The name of the component, use `ALL_COMPONENTS` to choose all components

**Returns** A new CompositionOfPhaseAsMoleFraction object.

```python
classmethod composition_of_phase_as_weight_fraction(phase: str, component: str) ->
tc_python.quantity.CompositionOfPhaseAsWeightFraction
```

Creates a quantity representing the composition of a phase [weight-fraction].

Parameters

- **phase** – The name of the phase, use `ALL_PHASES` to choose all stable phases
- **component** – The name of the component, use `ALL_COMPONENTS` to choose all components

**Returns** A new CompositionOfPhaseAsWeightFraction object.

```python
classmethod gibbs_energy_of_a_phase(phase: str, use_ser: bool = False) ->
tc_python.quantity.GibbsEnergyOfAPhase
```

Creates a quantity representing the Gibbs energy of a phase [J].

Parameters

- **phase** – The name of the phase or `ALL_PHASES` to choose all phases
- **use_ser** – Use Stable-Element-Reference(SER). The user-defined reference state will be used when this setting is set to False.

**Returns** A new GibbsEnergyOfAPhase object.

5.5. Module “quantity_factory”
class method mass_fraction_of_a_component(component: str) → tc_python.quantity.MassFractionOfAComponent

Creates a quantity representing the mass fraction of a component.

Parameters component – The name of the component or ALL_COMPONENTS to choose all components.

Returns A new MassFractionOfAComponent object.

class method mass_fraction_of_a_phase(phase: str) → tc_python.quantity.MassFractionOfAPhase

Creates a quantity representing the mass fraction of a phase.

Parameters phase – The name of the phase or ALL_PHASES to choose all phases.

Returns A new MassFractionOfAPhase object.

class method mole_fraction_of_a_component(component: str) → tc_python.quantity.MoleFractionOfAComponent

Creates a quantity representing the mole fraction of a component.

Parameters component – The name of the component or ALL_COMPONENTS to choose all components.

Returns A new MoleFractionOfAComponent object.

class method mole_fraction_of_a_phase(phase: str) → tc_python.quantity.MoleFractionOfAPhase

Creates a quantity representing the mole fraction of a phase.

Parameters phase – The name of the phase or ALL_PHASES to choose all phases.

Returns A new MoleFractionOfAPhase object.

class method normalized_driving_force_of_a_phase(phase: str) → tc_python.quantity.NormalizedDrivingForceOfAPhase

Creates a quantity representing normalized driving force of a phase [-].

Warning: A driving force calculation requires that the respective phase has been set to the state DORMANT. The parameter All is only reasonable if all phases have been set to that state.

Parameters phase – The name of the phase or ALL_PHASES to choose all phases.

Returns A new DrivingForceOfAPhase object.

class method pressure() → tc_python.quantity.Pressure

Creates a quantity representing the pressure [Pa].

Returns A new Pressure object.

class method system_size() → tc_python.quantity.SystemSize

Creates a quantity representing the system size [mol].

Returns A new SystemSize object.

class method temperature() → tc_python.quantity.Temperature

Creates a quantity representing the temperature [K].

Returns A new Temperature object.

class method tracer_diffusion_coefficient(phase: str, diffusing_element: str) → tc_python.quantity.TracerDiffusionCoefficient

Creates a quantity representing tracer diffusion coefficient of a phase [m^2/s].
Parameters

- **phase** – The name of the phase
- **diffusing_element** – The diffusing element

Returns A new TracerDiffusionCoefficient object.

```python
classmethod u_fraction_of_a_component(component: str) -> tc_python.quantity.UFractionOfAComponent
```

Creates a quantity representing the u-fraction of a component.

Parameters **component** – The name of the component

Returns A new UFractionOfAComponent object.

```python
classmethod user_defined_function(expression: str) -> tc_python.quantity.Function
```

Creates a quantity representing a user-defined function.

Parameters **expression** – The function expression

Returns A new Function object

```python
classmethod volume_fraction_of_a_phase(phase: str) -> tc_python.quantity.VolumeFractionOfAPhase
```

Creates a quantity representing the volume fraction of a phase.

Parameters **phase** – The name of the phase or ALL_PHASES to choose all phases

Returns A new VolumeFractionOfAPhase object.

---

### 5.6 Module “utils”

```python
class tc_python.utils.CompositionUnit(value)
    Bases: enum.Enum
    The composition unit.
    MASS_FRACTION = 2
        Mass fraction.
    MASS_PERCENT = 3
        Mass percent.
    MOLE_FRACTION = 0
        Mole fraction.
    MOLE_PERCENT = 1
        Mole percent.

class tc_python.utils.ConversionUnit(value)
    Bases: enum.Enum
    The composition unit used in a conversion.
    MOLE_FRACTION = 0
        Mole fraction.
    MOLE_PERCENT = 1
        Mole percent.
    WEIGHT_FRACTION = 2
        Weight fraction.
```
**WEIGHT_PERCENT = 3**

Weight percent.

class tc_python.utils.InterfacePosition(value)

Bases: enum.Enum

The position of an interface relative to its region. Only used for diffusion simulations.

**LOWER = 0**

The interface is on the lower side of its region.

**UPPER = 1**

The interface is on the upper side of its region.

class tc_python.utils.ResultValueGroup(result_line_group_java)

Bases: object

A x-y-dataset representing a line data calculation result (i.e. a Thermo-Calc quantity 1 vs. quantity 2).

| Warning: | Depending on the calculator, the dataset might contain NaN-values to separate the data between different subsets. |

**Variables**

- **label** – a str describing what the data corresponds to
- **x** – list of floats representing the first quantity (“x-axis”)
- **y** – list of floats representing the second quantity (“y-axis”)

get_label() → str

Accessor for the line label :return the line label

get_x() → List[float]

Accessor for the x-values :return the x values

get_y() → List[float]

Accessor for the y-values :return the y values

**class tc_python.utils.TemperatureProfile**

Bases: object

Represents a time-temperature profile used by non-isothermal calculations.

| Note: | The total simulation time can differ from the defined temperature profile. Constant temperature is assumed for any timepoint after the end of the defined profile. |

add_time_temperature(time: float, temperature: float)

Adds a time-temperature point to the non-isothermal temperature profile.

**Parameters**

- **time** – The time [s]
- **temperature** – The temperature [K]

**Returns** This TemperatureProfile object
5.7 Module “propertymodel_sdk”

class tc_python.propertymodel_sdk.CCTResult (quantity_id: str, description: str)
   Bases: tc_python.propertymodel_sdk.ResultQuantity

Represents a Continuous Cooling (CCT) result.

Parameters
   • quantity_id – The id of this result
   • description – The description of this result

add_time_temperature (time_temperature_id: str, description: str)
   Adds a time-temperature pair to the result.

Parameters
   • time_temperature_id – The id of the time-temperature pair
   • description – The description of the time-temperature pair

temperature_suffix = ' (T)'
   The temperature suffix of a CCTResult

time_suffix = ' (t)'
   The time suffix of a CCTResult

class tc_python.propertymodel_sdk.CCTResultValues (cooling_rate: float = -1.0,
   cooling_rate_start_temperature: float = -1.0, cooling_rate_end_temperature: float = -1.0)
   Bases: object

Represents Continuous Cooling (CCT) result values.

Parameters
   • cooling_rate – The cooling rate [K/s]
   • cooling_rate_start_temperature – The start temperature of cooling [K]
   • cooling_rate_end_temperature – The end temperature of cooling [K]

set_result_time_temperature (time_temperature_id: str, time: float, temperature: float)
   Sets a time-temperature pair of the result.

Parameters
   • time_temperature_id – The id of the time-temperature pair
   • time – The time [s]
   • temperature – The temperature [K]

class tc_python.propertymodel_sdk.CalculationContext (system: tc_python.system.System,
   model_utils=None)
   Bases: object

Represents the interface of the Property Model with the Thermo-Calc application and the rest of the TC-Python functionality.

Parameters
• **system** – The system object of this calculation

• **model_utils** – The model utils object

**get_argument_ids()** → Set[str]

Returns a list with the arguments of the models. Including arguments created from dynamic parameters.

**get_dependent_component()** → str

Obtains the dependent component from the UI

**Note:** The dependent component is that which has no composition specified explicitly, typically this is the major element of the material (such as Fe, Al, Ni, ...)

**Returns** The dependent component

**get_mass_fractions()** → Dict[str, float]

Obtains the current composition from the UI as mass-fraction.

**Note:** In case of stepping over one or multiple axis, the returned data will represent the composition at the current step.

**Returns** The composition (key: component, value: content) [mass-fraction]

**get_mass_percents()** → Dict[str, float]

Obtains the current composition from the UI in mass-percent.

**Note:** In case of stepping over one or multiple axis, the returned data will represent the composition at the current step.

**Returns** The composition (key: component, value: content) [mass-percent]

**get_mole_fractions()** → Dict[str, float]

Obtains the current composition from the UI as mole-fraction.

**Note:** In case of stepping over one or multiple axis, the returned data will represent the composition at the current step.

**Returns** The composition (key: component, value: content) [mole-fraction]

**get_mole_percents()** → Dict[str, float]

Obtains the current composition from the UI in mole-percent.

**Note:** In case of stepping over one or multiple axis, the returned data will represent the composition at the current step.

**Returns** The composition (key: component, value: content) [mole-percent]
get_temperature() → float
Obtains the current temperature from the UI.

Returns The temperature [K]

get_ui_boolean_value(component_id: str) → bool
Obtains the value from the specified checkbox UI component.

Parameters component_id – Id of the checkbox

Returns The setting of the checkbox

get_ui_condition_list(component_id: str) → tc_python.propertymodel_sdk.ConditionListEntry
Used to get the selected condition from components of type UIConditionListComponent.

Parameters component_id: Id of the list UI component

Returns The selected condition

get_ui_float_value(component_id: str) → float
Obtains the value from the specified UI component.

Parameters component_id – Id of the UI component

Returns The value

get_ui_list_value(component_id: str) → str
Obtains the selected entry from a UI component list. If a special element (such as ANY, NONE, ...) is selected, the corresponding locale-independent placeholder is provided.

Parameters component_id – Id of the list UI component

Returns The selected entry

get_ui_string_value(component_id: str) → str
Obtains the selected entry from a UI component text field.

Parameters component_id – Id of the string UI component

Returns The selected entry

get_ui_temperature_value(component_id: str) → float
Obtains the temperature from the specified temperature UI component.

Parameters component_id – Id of the temperature UI component

Returns The temperature [K], note that input unit of the UI is specified in the model panel. If required, the temperature is automatically converted to K.

set_result_cct_values(quantity_id: str, r: tc_python.propertymodel_sdk.CCTResultValues)
Sets the value of a previously defined result quantity (of type CCTResultValues) for further usage in the Thermo-Calc application for plotting, etc.

Parameters

• quantity_id – unique id of the result quantity
• r – the CCTResultValues to be set

set_result_quantity_value(quantity_id: str, value: float, parameter: str = '')
Sets the value of a previously defined result quantity for further usage in the Thermo-Calc application for plotting, etc.

Parameters

Note: Any result quantity that remains unset is automatically set to NaN.
• **quantity_id** – Unique id of the result quantity
• **parameter** – Use if result is parameterized. f.i. “per phase”
• **value** – The value to be set

```python
def set_single_equilibrium_result(quantity_id: str, r: tc_python.single_equilibrium.SingleEquilibriumResult)```

```python
class tc_python.propertymodel_sdk.ConditionListEntry
    Bases: object

    Used in combination with components of type UIConditionListComponent.
    Contains the element, if the selected condition is a composition
    Contains the Console Mode syntax of the selected condition
    Contains the unit of the selected condition
```

```python
class tc_python.propertymodel_sdk.PropertyModel(_locale: str = 'en-US')
    Bases: object

    The abstract base class for all property models.
```

**Note:** Every Property Model needs to implement most of the abstract methods of this class. However, some abstract methods are optional and should only be implemented if required.

**Note:** If overwriting the constructor in a Property Model, the constructor of the implemented class must have the identical signature and should pass the parameters to this base class constructor.

**Tip:** It is possible to switch off internal INFO-log messages coming from the calculation engine by changing the log-level on the TC-Python log object like this: `logging.getLogger(“tc_python”).setLevel(logging.ERROR)`.

**Parameters**

- `_locale` – The locale to be used, this is an internal parameter and is of no meaning to the end-user

**Variables**

- `logger` – logger object that is connected to the Thermo-Calc UI (`INFO-` and `WARN-ING-level will be printed as `INFO`, `ERROR`-level as `ERROR`), it can be accessed like this: `self.logger.info("Some message")`

```python
@abstractmethod
def add_button_callback(component_id: str, ui_component_ids: Dict[str, int]) → List[tc_python.propertymodel_sdk.UIComponent]
```

Implement this method if you have one or more UI components on which you called `UIComponent.enable_add_button()`, which adds a + button next to the component.

This method will be executed when you press any such + button.

This method is typically used to add more UI components dynamically and the method must return a list of the UI components to be added.

**This method can optionally be implemented by a Property Model.**

**Parameters**

- `component_id` – The id of the UI component next to the pressed + button
- `ui_component_ids` – dict with the id:s and the index of the current ui components of the model. Including arguments created in previous calls to add_button_callback.

**Returns** A list of `UIComponent` objects to be added
abstract after_evaluations()
Called by the Thermo-Calc application immediately after the last model evaluation (using the method `PropertyModel.evaluate_model()`). Use this method for any required cleanup.

This method can optionally be implemented by a Property Model.

abstract before_evaluations(context: tc_python.propertymodel_sdk.CalculationContext)
Called by the Thermo-Calc application immediately before the first model evaluation (using the method `PropertyModel.evaluate_model()`). Use this method for any required preparations.

This method can optionally be implemented by a Property Model.

Parameters
context – The calculation context

abstract evaluate_model(context: tc_python.propertymodel_sdk.CalculationContext)
Called by the Thermo-Calc application when the model should be actually calculated. This is the main method of the Property Model that contains the actual calculation code.

This method needs to be implemented by all property models.

Parameters
context – The calculation context, this provides access to the Thermo-Calc application and all other TC-Python modules

abstract get_license_key() → str
Provides the license key of the model.

This method can optionally be implemented by a Property Model.

abstract provide_calculation_result_quantities() → List[tc_python.propertymodel_sdk.ResultQuantity]
Called by the Thermo-Calc application when the model should provide its result quantity objects.

This method needs to be implemented by all property models.

Returns
Result quantity objects of the model (to be filled later with results in the method `PropertyModel.evaluate_model()`)

abstract provide_model_category() → List[str]
Called by the Thermo-Calc application when the model should provide its category (shown in the Thermo-Calc model tree).

This method needs to be implemented by all property models.

Returns
Category of the model, it may be present in several categories

abstract provide_model_description() → str
Called by the Thermo-Calc application when the model should provide its detailed description.

This method needs to be implemented by all property models.

Returns
Description text for the model

abstract provide_model_name() → str
Called by the Thermo-Calc application when the model should provide its name (shown in the Thermo-Calc model tree).

This method needs to be implemented by all property models.

Returns
Name of the model

abstract provide_model_parameters() → Dict[str, float]
Called by the Thermo-Calc application when the model should provide all model parameters and their current values.

This method can optionally be implemented by a Property Model.
Note: These are internal variables of the Property Model that are intended to be modified from the outside. Typically this is used to adjust their values in a optimizer during the development of the model.

**Returns** The model parameter ids and their current values [unit according to the parameter meaning]

```python
abstract provide_ui_panel_components() → List[tc_python.propertymodel_sdk.UIComponent]
```

Called by the Thermo-Calc application when the model should provide its UI components for the model panel to be plotted. This happens also whenever a model gets checked in the model tree.

**This method needs to be implemented by all property models.**

**Returns** Model UI panel components in the order to be presented in the model panel

```python
abstract remove_button_callback(component_id: str, ui_component_ids: Dict[str, int]) → List[str]
```

Implement this method if you have one or more UI components on which you called UIComponent.enable_remove_button(), which adds a - button next to the component.

This method will be executed when you press any such - button.

This method is typically used to remove UI components dynamically and the method must return a list of the ids of the components that are going to be removed.

**This method can optionally be implemented by a Property Model.**

**Parameters**

- **component_id** – the id of the UI component next to the pressed - button
- **ui_component_ids** – list with the id:s and the index of the current ui components of the model. Including arguments created in previous calls to add_button_callback.

**Returns** a list of UI component ids that are required to be removed

```python
abstract set_model_parameter(model_parameter_id: str, value: float)
```

Called by the Thermo-Calc application when a model parameter should be reset.

**This method can optionally be implemented by a Property Model.**

**Note:** These are internal variables of the Property Model that are intended to be modified from the outside. Typically this is used to adjust their values in a optimizer during the development of the model.

**Parameters**

- **model_parameter_id** – The parameter id
- **value** – The value [unit according to the parameter meaning]

```python
class tc_python.propertymodel_sdk.ResultQuantity(quantity_id: str, description: str, quantity_type: tc_python.propertymodel_sdk.ResultQuantityType)
```

**Bases:** object

Defines a calculation result quantity of a Property Model that is identified by a unique id.

**Parameters**

- **quantity_id** – Unique id of the quantity
• description – Description of the quantity (shown in the Thermo-Calc UI)

• quantity_type – Type of the quantity (defines the unit)

get_description() → str
Obtains the description of the quantity.

Returns Description of the quantity

get_id() → str
Obtains the id of the quantity.

Returns Unique id of the quantity

get_type() → tc_python.propertymodel_sdk.ResultQuantityType
Obtains the type of quantity.

Returns Type of the quantity

class tc_python.propertymodel_sdk.ResultQuantityType(value)
Bases: enum.Enum
Defining the type of a result quantity.

CCT_QUANTITY = 5
A cct quantity

ENERGY_QUANTITY = 2
An energy quantity

GENERAL_QUANTITY = 0
A general quantity

LENGTH_QUANTITY = 7
A length in quantity

SINGLE_EQUILIBRIUM_QUANTITY = 6
A cct quantity

STRENGTH_QUANTITY = 8
A strength quantity

SURFACE_ENERGY_QUANTITY = 3
A surface energy quantity

TEMPERATURE_QUANTITY = 1
A temperature quantity

TIME_QUANTITY = 4
A time quantity

class tc_python.propertymodel_sdk.SpecialListMarkers
Bases: object
Placeholders for special list elements that are locale-dependent. They will be provided by UI list components if a special marker has been selected.

ANY_LIST_MARKER = 'ANY'
Marker that represents “Any”

NONE_LIST_MARKER = 'NONE'
Marker that represents “None”

class tc_python.propertymodel_sdk.UIBooleanComponent(component_id: str, name: str, description: str, setting: bool)
Bases: tc_python.propertymodel_sdk.UIComponent

5.7. Module “propertymodel_sdk”
Checkbox UI component of the model panel.

Parameters

- **component_id** – Unique id of the component
- **name** – Name of the component, will be presented in the model panel
- **description** – Additional description of the component
- **setting** – Initial setting of the checkbox

`connect_component_visibility(dependent_component_id: str)`
Connects the visibility of any other UI component of the model panel to the value of this boolean component.

Parameters

- **dependent_component_id** – Id of the UI element to be dependent on this boolean component

`enable_add_button()`
Adds a + button to the right of the UI component.

Returns

This UI component

`enable_remove_button()`
Adds a - button to the right of the UI component.

Returns

This UI component

`get_dependent_components() → List[str]`
Obtains a list containing all UI elements currently connected regarding their visibility.

Returns

A list with the component id of all UI elements currently connected

`get_setting() → bool`
Obtains the setting of the checkbox.

Returns

The setting of the checkbox

`remove_component_visibility(dependent_component_id: str)`
Removes the visibility connection to a UI component that has been previously connected.

Parameters

- **dependent_component_id** – Id of the previously connection UI element

`set_index(index: int = -1)`
Sets the position in the graphical user interface.

Parameters

- **index** – The position

Returns

This UI component

```python
class tc_python.propertymodel_sdk.UIComponent(component_id: str, name: str, description: str)
Bases: object
```

Abstract Base class for all UI components of the model panel.

Never make an instance of UIComponent, always use the sub-classes. For instance UIStringComponent.

Parameters

- **component_id** – Unique id of the component
- **name** – Name of the component, will be presented in the model panel
- **description** – Additional description of the component
get_description() → str
Obtains the additional description of the component.

Returns Additional description of the component

get_id() → str
Obtains the unique id of the component.

Returns Unique id of the component

get_name() → str
Obtains the name of the component.

Returns Name of the component, will be presented in the model panel

class tc_python.propertymodel_sdk.UIConditionListComponent(component_id: str, name: str, description: str)
Bases: tc_python.propertymodel_sdk.UIComponent
System condition list UI component of the model panel.

Parameters

- component_id – Unique id of the component
- name – Name of the component, will be presented in the model panel
- description – Additional description of the component

class tc_python.propertymodel_sdk.UIFloatComponent(component_id: str, name: str, description: str, value: float)
Bases: tc_python.propertymodel_sdk.UIComponent
General real value text field UI component of the model panel.

Parameters

- component_id – Unique id of the component
- name – Name of the component, will be presented in the model panel
- description – Additional description of the component
- value – Initial setting of the text field

enable_add_button()
Adds a + button to the right of the UI component.

Returns This UI component

enable_remove_button()
Adds a - button to the right of the UI component.

Returns this UI component

get_value() → float
Obtains the setting of the text field.

Returns The setting of the text field

set_index(index: int = -1)
Sets the position in the graphical user interface.

Parameters index – The position

Returns This UI component
**class** `tc_python.propertymodel_sdk.UIGeneralListComponent` *(component_id: str, name: str, description: str, content: List[Tuple[str, str]], selected_entry: str = '')*

**Bases:** `tc_python.propertymodel_sdk.UIComponent`

General list UI component of the model panel that can contain any strings.

**Parameters**

- **component_id** – Unique id of the component
- **name** – Name of the component, will be presented in the model panel
- **description** – Additional description of the component
- **content** – Entries of the list, they need to contain a locale-independent id and a localized content string, for example: `[('ENTRY_1_ID', 'entry 1'), ('ENTRY_2_ID', 'entry 2')]`
- **selected_entry** – Entry to be initially selected. If omitted, by default the first element is selected.

**connect_component_visibility** *(dependent_component_id: str, selected_item_to_set_visible: str)*

Connects the visibility of any other UI component of the model panel to the selection of a certain entry of the list.

**Parameters**

- **dependent_component_id** – Id of the UI element to be dependent on the chosen element
- **selected_item_to_set_visible** – Entry (locale independent id) of the list to be chosen to set the dependent component visible

**enable_add_button**()

Adds a + button to the right of the UI component.

**Returns** This UI component

**enable_remove_button**()

Adds a - button to the right of the UI component.

**Returns** This UI component

**get_content**() → List[Tuple[str, str]]

Obtains the entries of the list.

**Returns** Entries of the list, they need to contain a locale-independent id and a localized content string, for example: `[('ENTRY_1_ID', 'entry 1'), ('ENTRY_2_ID', 'entry 2')]`

**get_dependent_components**() → Dict[str, List[str]]

Obtains a dictionary containing all UI elements currently connected regarding their visibility.

**Returns** All UI elements currently connected (key: dependent component id, value: required list entries to set it visible)

**get_selected_entry**() → str

Obtains the initially selected entry.

**Returns** Initially selected entry. If empty, the first element is selected.

**remove_component_visibility** *(dependent_component_id: str)*

Removes the visibility connection to a UI component that has been previously connected.
Parameters **dependent_component_id** – Id of the previously connection UI element

**set_index** *(index: int = -1)*

Sets the position in the graphical user interface.

**Parameters** **index** – The position

**Returns** This UI component

class tc_python.propertymodel_sdk.UIPhaseListComponent *(component_id: str, name: str, description: str, default_phase: str = '', any_marker_setting: bool = False)*

**Bases:** tc_python.propertymodel_sdk.UIComponent

Phase list UI component of the model panel.

**Parameters**

- **component_id** – Unique id of the component
- **name** – Name of the component, will be presented in the model panel
- **description** – Additional description of the component
- **default_phase** – Default phase, if omitted no default phase is chosen and only initially the first element of the list is selected. If an ANY-marker is added, this is chosen as the default element.
- **any_marker_setting** – Defines if an entry “ANY PHASE” should be added to the phase list, if set to true this overrides any default phase setting

**enable_add_button**

Adds a + button to the right of the UI component.

**Returns** This UI component

**enable_remove_button**

Adds a - button to the right of the UI component.

**Returns** This UI component

**get_any_marker_setting** *(→ bool)*

Obtains the setting if any entry “ANY PHASE” is added to the phase list.

**Returns** If an entry “ANY PHASE” is added to the phase list, if set to true this overrides any default phase setting

**get_default_phase** *(→ str)*

Obtains the default phase.

**Returns** Default phase, if omitted no default phase is chosen and only initially the first element of the list is selected. If an ANY-marker is added, this is chosen as the default element.

**set_index** *(index: int = -1)*

Sets the position in the graphical user interface.

**Parameters** **index** – The position

**Returns** This UI component

class tc_python.propertymodel_sdk.UIStringComponent *(component_id: str, name: str, description: str, string: str)*

**Bases:** tc_python.propertymodel_sdk.UIComponent
General text field UI component of the model panel.

Parameters

- `component_id` – Unique id of the component
- `name` – Name of the component, will be presented in the model panel
- `description` – Additional description of the component
- `string` – Initial setting of the text field

`enable_add_button()`
Adds a + button to the right of the UI component.

 Returns  This UI component

`enable_remove_button()`
Adds a - button to the right of the UI component.

 Returns  This UI component

`get_value()` → `str`
Obtains the setting of the text field.

 Returns  The setting of the text field

`set_index(index: int = -1)`
Sets the position in the graphical user interface.

 Parameters  index – The position

 Returns  This UI component

```python
class tc_python.propertymodel_sdk.UITemperatureComponent:
    component_id: str,
    name: str,
    description: str,
    temp: float

Bases: tc_python.propertymodel_sdk.UIComponent
```

Temperature value text field UI component of the model panel.

Parameters

- `component_id` – Unique id of the component
- `name` – Name of the component, will be presented in the model panel
- `description` – Additional description of the component
- `temp` – Initial temperature to be set in the text field (unit defined by the user in the Thermo-Calc system)

`enable_add_button()`
Adds a + button to the right of the UI component.

 Returns  This UI component

`enable_remove_button()`
Adds a - button to the right of the UI component.

 Returns  This UI component

`get_temp()` → `float`
Obtains the temperature set in the text field.

 Returns  The temperature to be set in the text field (unit defined by the user in the Thermo-Calc system)
**set_index** *(index: int = -1)*
Sets the position in the graphical user interface.

**Parameters**
- **index** – The position

**Returns**
This UI component

```python
tc_python.propertymodel_sdk.create_boolen_ui_component(component_id: str, name: str, description: str, initial_setting: bool) → tc_python.propertymodel_sdk.UIBooleanComponent
```

Creates a UI checkbox component for a boolean value. The value of that component can later be accessed during the model evaluation.

**Parameters**
- **component_id** – Unique id of the component
- **name** – Name of the component, will be presented in the model panel
- **description** – Additional description of the component
- **initial_setting** – Initial setting of the checkbox

**Returns**
The created component

```python
tc_python.propertymodel_sdk.create_condition_list_ui_component(component_id: str, name: str, description: str) → tc_python.propertymodel_sdk.UIConditionListComponent
```

Creates a UI list component for all conditions defined in the system. The value of that component can later be accessed during the model evaluation.

**Parameters**
- **component_id** – Unique id of the component
- **name** – Name of the component, will be presented in the model panel
- **description** – Additional description of the component

**Returns**
The created component

```python
tc_python.propertymodel_sdk.create_energy_quantity(quantity_id: str, description: str) → tc_python.propertymodel_sdk.ResultQuantity
```

Creates a UI energy result quantity (in J). When the model is evaluated, a value can be added to the quantity and it will be used to transfer the result to the Thermo-Calc plot engine.

**Parameters**
- **quantity_id** – Unique id of the result quantity
- **description** – Additional description of the result quantity

**Returns**
The created result quantity

```python
tc_python.propertymodel_sdk.create_float_ui_component(component_id: str, name: str, description: str, value: float) → tc_python.propertymodel_sdk.UIFloatComponent
```

Creates a UI text field component for a real number. The value of that component can later be accessed during the model evaluation.

**Parameters**

• component_id – Unique id of the component
• name – Name of the component, will be presented in the model panel
• description – Additional description of the component
• value – Initial setting of the text field

Returns  The created component

tc_python.propertymodel_sdk.create_general_quantity(quantity_id: str, description: str) -> tc_python.propertymodel_sdk.ResultQuantity

Creates a general result quantity that can contain any type of result (without a unit). When the model is evaluated, a value can be added to the quantity and it will be used to transfer the result to the Thermo-Calc plot engine.

Parameters
• quantity_id – Unique id of the result quantity
• description – Additional description of the result quantity

Returns  The created result quantity

tc_python.propertymodel_sdk.create_length_quantity(quantity_id: str, description: str) -> tc_python.propertymodel_sdk.ResultQuantity

Creates a length result quantity. When the model is evaluated, a value can be added to the quantity and it will be used to transfer the result to the Thermo-Calc plot engine.

Parameters
• quantity_id – Unique id of the result quantity
• description – Additional description of the result quantity

Returns  The created result quantity

tc_python.propertymodel_sdk.create_list_ui_component(component_id: str, name: str, description: str, entry_list: List[Tuple[str, str]], selected_entry: str = '') -> tc_python.propertymodel_sdk.UIGeneralListComponent

Creates a UI list component for string entries. The value of that component can later be accessed during the model evaluation.

Parameters
• component_id – Unique id of the component
• name – Name of the component, will be presented in the model panel
• description – Additional description of the component
• entry_list – Entries of the list, they need to contain a locale-independent id and a localized content string, for example: ["ENTRY_1_ID", "entry 1"], (ENTRY_2_ID", "entry 2")
• selected_entry – Entry to be initially selected. If omitted, by default the first element is selected.

Returns  The created component
tc_python.propertymodel_sdk.create_phase_list_ui_component(component_id: str,
name: str, description: str, default_phase: str = '', any_marker: bool = False) → tc_python.propertymodel_sdk.UIPhaseListComponent

Creates a UI list component for all phases defined in the system. It is possible to select a default phase that is supposed to be the expected phase selection for that list. The value of that component can later be accessed during the model evaluation.

A default phase is the phase that is initially selected and re-selected as soon as a currently selected phase is removed. If the default phase is not available, a “NONE”-marker will be created and used instead of the default phase. A typical use case for the default phase setting is a phase list that expects to contain the LIQUID-phase of a system.

Parameters

- **component_id** – Unique id of the component
- **name** – Name of the component, will be presented in the model panel
- **description** – Additional description of the component
- **default_phase** – Default phase, if omitted no default phase is chosen and only initially the first element of the list is selected. **If an ANY-marker is added, this is chosen as the default element.**
- **any_marker** – Defines if an entry “ANY PHASE” should be added to the phase list, if set to true this overrides any default phase setting

Returns The created component

5.7.1. Module “propertymodel_sdk”

5.7. Module “propertymodel_sdk”

tc_python.propertymodel_sdk.create_strength_quantity(quantity_id: str, description: str) → tc_python.propertymodel_sdk.ResultQuantity

Creates a strength result quantity. When the model is evaluated, a value can be added to the quantity and it will be used to transfer the result to the Thermo-Calc plot engine.

Parameters

- **quantity_id** – Unique id of the result quantity
- **description** – Additional description of the result quantity

Returns The created result quantity

5.7. Module “propertymodel_sdk”

tc_python.propertymodel_sdk.create_string_ui_component(component_id: str,
name: str, description: str, string: str) → tc_python.propertymodel_sdk.UIStringComponent

Creates a UI text field component. The value of that component can later be accessed during the model evaluation.

Parameters

- **component_id** – Unique id of the component
- **name** – Name of the component, will be presented in the model panel
- **description** – Additional description of the component
- **string** – Initial setting of the text field
Returns The created component

tc_python.propertymodel_sdk.create_surface_energy_quantity(quantity_id: str, description: str) → tc_python.propertymodel_sdk.ResultQuantity

Creates an energy result quantity (in J). When the model is evaluated, a value can be added to the quantity and it will be used to transfer the result to the Thermo-Calc plot engine.

Parameters

- **quantity_id** – Unique id of the result quantity
- **description** – Additional description of the result quantity

Returns The created result quantity

tc_python.propertymodel_sdk.create_temperature_quantity(quantity_id: str, description: str) → tc_python.propertymodel_sdk.ResultQuantity

Creates a temperature result quantity (in K). When the model is evaluated, a value can be added to the quantity and it will be used to transfer the result to the Thermo-Calc plot engine.

Parameters

- **quantity_id** – Unique id of the result quantity
- **description** – Additional description of the result quantity

Returns The created result quantity

tc_python.propertymodel_sdk.create_temperature_ui_component(component_id: str, name: str, description: str, initial_temp: float) → tc_python.propertymodel_sdk.UITemperature

Creates a UI text field component for a temperature value. The value of that component can later be accessed during the model evaluation.

Parameters

- **component_id** – Unique id of the component
- **name** – Name of the component, will be presented in the model panel
- **description** – Additional description of the component
- **initial_temp** – Initial temperature to be set in the text field. (The unit of initial_temp is Kelvin. The value in the text field will be automatically converted using the unit chosen by the user.)

Returns The created component

tc_python.propertymodel_sdk.create_time_quantity(quantity_id: str, description: str) → tc_python.propertymodel_sdk.ResultQuantity

Creates a time result quantity (in s). When the model is evaluated, a value can be added to the quantity and it will be used to transfer the result to the Thermo-Calc plot engine.

Parameters

- **quantity_id** – Unique id of the result quantity
- **description** – Additional description of the result quantity

Returns The created result quantity
5.8 Module “exceptions”

```python
exception tc_python.exceptions.APIServerException
    Bases: tc_python.exceptions.GeneralException
    An exception that occurred during the communication with the API-server. It is normally not related to an error in the user program.

exception tc_python.exceptions.CalculationException
    Bases: tc_python.exceptions.TCException
    An exception that occurred during a calculation.

exception tc_python.exceptions.ComponentNotExistingException
    Bases: tc_python.exceptions.GeneralException
    The selected component is not existing.

exception tc_python.exceptions.DatabaseException
    Bases: tc_python.exceptions.CalculationException
    Error loading a thermodynamic or kinetic database, typically due to a misspelled database name or a database missing in the system.

exception tc_python.exceptions.DegreesOfFreedomNotZeroException
    Bases: tc_python.exceptions.CalculationException
    The degrees of freedom in the system are not zero, i.e. not all required conditions have been defined. Please check the conditions given in the exception message.

exception tc_python.exceptions.EquilibriumException
    Bases: tc_python.exceptions.CalculationException
    An equilibrium calculation has failed, this might happen due to inappropriate conditions or a very difficult problem that can not be solved.

exception tc_python.exceptions.GeneralCalculationException
    Bases: tc_python.exceptions.CalculationException
    General error occurring while a calculation is performed.

exception tc_python.exceptions.GeneralException
    Bases: tc_python.exceptions.TCException
    A general exception that might occur in different situations.

exception tc_python.exceptions.InvalidCalculationConfigurationException
    Bases: tc_python.exceptions.CalculationException
    Thrown when errors are detected in the configuration of the calculation.

exception tc_python.exceptions.InvalidCalculationStateException
    Bases: tc_python.exceptions.CalculationException
    Trying to access an invalid calculation object that was invalidated by calling invalidate on it.

exception tc_python.exceptions.InvalidNumberOfResultGroupsException
    Bases: tc_python.exceptions.ResultException
    A calculation result contains several result groups, which is not supported for the used method.

exception tc_python.exceptions.InvalidResultConfigurationException
    Bases: tc_python.exceptions.ResultException
    A calculation result configuration is invalid.
```
exception tc_python.exceptions.InvalidResultStateException
Bases: tc_python.exceptions.CalculationException

Trying to access an invalid result (for example a SingleEquilibriumTempResult object that got already invalidated by condition changes or a result that was invalidated by calling invalidate on it).

exception tc_python.exceptions.LicenseException
Bases: tc_python.exceptions.GeneralException

No valid license for the API or any Thermo-Calc product used by it found.

exception tc_python.exceptions.NoDataForPhaseException
Bases: tc_python.exceptions.ResultException

There is no result data available for a selected phase.

exception tc_python.exceptionsNotAllowedOperationException
Bases: tc_python.exceptions.CalculationException

The called method or operation is not allowed in the current mode of operation (i.e. debug or production mode). Production mode means that the Property Model is only present as an *.py,encrypted-file, while in debug mode it is available as *.py-file. Certain methods for obtaining internal model parameters are not available for encrypted models.

exception tc_python.exceptions.PhaseNotExistingException
Bases: tc_python.exceptions.GeneralException

The selected phase is not existing, so no data can be provided for it.

exception tc_python.exceptions.ResultException
Bases: tc_python.exceptions.TCException

An exception that occurred during the configuration of a calculation result.

exception tc_python.exceptions.SyntaxException
Bases: tc_python.exceptions.CalculationException

Syntax error in a Console Mode expression.

exception tc_python.exceptions.TCException
Bases: Exception

The root exception of TC-Python.

exception tc_python.exceptions.UnrecoverableCalculationException
Bases: tc_python.exceptions.CalculationException

The calculation reached a state where no further actions are possible, this happens most often due to a FORTRAN- hard crash in the API server backend.

Note: It is possible to catch that exception outside of the with-clause context and to continue by setting up a new context (i.e. by a new with TCPython() as session).

tc_python.exceptions.handle_exception(e)
5.9 Module “abstract_base”

class tc_python.abstract_base.AbstractCalculation (calculator)
    Bases: object

    Abstract base class for calculations.

get_configuration_as_string () → str
    Returns detailed information about the current state of the calculation object.

    Warning: The structure of the calculator objects is an implementation detail and might change between releases without notice. Therefore do not rely on the internal object structure.

get_system_data () → tc_python.abstract_base.SystemData
    Returns the content of the database for the currently loaded system. This can be used to modify the parameters and functions and to change the current system by using with_system_modifications().

    Note: Parameters can only be read from unencrypted (i.e. user) databases loaded as *.tdb-file.

    Returns The system data

invalidate ()
    Invalidates the object and frees the disk space used by it. This is only required if the disk space occupied by the object needs to be released during the calculation. No data can be retrieved from the object afterwards.

with_system_modifications (system_modifications: tc_python.abstract_base.SystemModifications)
    Updates the system of this calculator with the supplied system modification (containing new phase parameters and system functions).

    Note: This is only possible if the system has been read from unencrypted (i.e. user) databases loaded as a *.tdb-file.

    Parameters system_modifications – The system modification to be performed

    Returns

class tc_python.abstract_base.AbstractResult (result)
    Bases: object

    Abstract base class for results. This can be used to query for specific values.

invalidate ()
    Invalidates the object and frees the disk space used by it. This is only required if the disk space occupied by the object needs to be released during the calculation. No data can be retrieved from the object afterwards.

class tc_python.abstract_base.PhaseParameter (parameter_name: Union[str, object])
    Bases: object

    Database phase parameter expression used by SystemModifications.set().

    Parameters parameter_name – The phase parameter name

get_intervals () → List[tc_python.abstract_base.TemperatureInterval]
    Returns the list of all defined intervals.
Returns The defined temperature intervals

get_lower_temperature_limit() → float
Returns the lower temperature limit.

Returns The lower temperature limit in K

get_name() → str
Returns the name of the phase parameter.

remove_all_intervals()
Removes all previously defined temperature intervals.

Returns This PhaseParameter object

remove_interval_with_upper_limit(upper_temperature_limit: float)
Removes a previously defined temperature interval with matching upper temperature limit.

If no such interval exists, an exception is thrown.

Returns This PhaseParameter object

set_expression_with_upper_limit(parameter_expression: str, upper_temperature_limit: float = 6000.0)
Adds/overwrites a parameter expression for a temperature interval.

Default value of the upper limit of the interval: 6000 K

Note: The lower temperature limit is either defined by the lower temperature limit given with PhaseParameter.set_lower_temperature_limit() or by the upper temperature limit of the adjacent interval.

Note: If there is an existing interval with exactly the same upper_temperature_limit, that interval is overwritten, otherwise the interval is added.

Parameters

• parameter_expression – The parameter expression, example: 
  +V34*T*LN(T)+V35*T**2+V36*T**(-1)+V37*T**3"

• upper_temperature_limit – The upper temperature limit for which the expression should be used

Returns This PhaseParameter object

set_interval(interval: tc_python.abstract_base.TemperatureInterval)
Adds/overwrites a temperature interval.

Note: The lower temperature limit is either defined by the lower temperature limit given with PhaseParameter.set_lower_temperature_limit() or by the upper temperature limit of the adjacent interval.
Note: If there is an existing interval with exactly the same `upper_temperature_limit`, that interval is overwritten, otherwise the interval is added.

Returns This `PhaseParameter` object

```python
set_lower_temperature_limit(lower_temperature_limit: float = 298.15)
```
Sets the lower temperature limit of the phase parameter.

Default: 298.15 K

Parameters `lower_temperature_limit` – The lower temperature limit in K

Returns This `PhaseParameter` object

```python
class tc_python.abstract_base.SystemData(system_data)
```
Bases: `object`

Provides information about the parameters and functions of a user database. The obtained objects can be used to modify the database using `with_system_modifications()` of all calculators.

Note: Parameters can only be read from unencrypted (i.e. `user`) databases loaded as *.tdb*-file.

```python
get_phase_parameter(parameter: str) \rightarrow tc_python.abstract_base.PhaseParameter
```
Returns a phase parameter.

Example:

```python
system_data.get_phase_parameter('G(HCP_A3,FE;0)')
```

Note: Parameters can only be read from unencrypted (i.e. `user`) databases loaded as a *, tdb-file.

Note: For details about the syntax search the Thermo-Calc help for GES (the name for the Gibbs Energy System module in Console Mode).

Parameters `parameter` – The name of the phase parameter (for example: “G(LIQUID,FE;0)”)

Returns The phase parameter

```python
get_phase_parameter_names() \rightarrow List[str]
```
Returns all phase parameters present in the current system.

Returns The list of phase parameters

```python
get_system_function(f: str) \rightarrow tc_python.abstract_base.SystemFunction
```
Returns a system function.

Note: The parameter ‘f’ was previously called ‘function’ but was renamed.

Example:

```python
system_data.get_system_function('GHSERCR')
```
Note: Functions can only be read from unencrypted (i.e. user) databases loaded as a *.tdb-file.

Note: For details about the syntax search the Thermo-Calc help for GES (the name for the Gibbs Energy System module in Console Mode).

Parameters \( f \) – The name of the system function (for example: “GHSERCR”)

Returns The system function

get_system_function_names() \( \rightarrow \) List[str]
Returns all system functions present in the current system.

Returns The list of system functions

class tc_python.abstract_base.SystemFunction (function_name: Union[str, object])
Bases: object
Database function expression used by SystemModifications.set().

Parameters function_name – The function name

get_intervals() \( \rightarrow \) List[tc_python.abstract_base.TemperatureInterval]
Returns the list of all defined intervals.

Returns The defined temperature intervals

get_lower_temperature_limit() \( \rightarrow \) float
Returns the lower temperature limit.

Returns The lower temperature limit in K

get_name() \( \rightarrow \) str
Returns the name of the system function.

Returns The name of the system function

remove_all_intervals()
Removes all previously defined temperature intervals.

Returns This SystemFunction object

remove_interval_with_upper_limit (upper_temperature_limit: float)
Removes a previously defined temperature interval with matching upper temperature limit.

If no such interval exists, an exception is thrown.

Returns This SystemFunction object

set_expression_with_upper_limit (function_expression: str, upper_temperature_limit: float = 6000.0)
Adds/overwrites a function expression for a temperature interval.

Default value of the upper limit of the interval: 6000 K

Note: The lower temperature limit is either defined by the lower temperature limit given with SystemFunction.set_lower_temperature_limit() or by the upper temperature limit of the adjacent interval.
Note: If there is an existing interval with exactly the same upper_temperature_limit, that interval is overwritten, otherwise the interval is added.

Parameters

- **function_expression** – The function expression, example:
  
  \[ +V34*T*LN(T)+V35*T**2+V36*T**(-1)+V37*T**3 \]

- **upper_temperature_limit** – The upper temperature limit for which the expression should be used

Returns This SystemFunction object

**set_interval (interval: tc_python.abstract_base.TemperatureInterval)**

 Adds/overwrites a temperature interval.

Note: The lower temperature limit is either defined by the lower temperature limit given with SystemFunction.set_lower_temperature_limit() or by the upper temperature limit of the adjacent interval.

Note: If there is an existing interval with exactly the same upper_temperature_limit, that interval is overwritten, otherwise the interval is added.

Returns This SystemFunction object

**set_lower_temperature_limit (lower_temperature_limit: float = 298.15)**

Sets the lower temperature limit of the system function.

Default: 298.15 K

Parameters **lower_temperature_limit** – The lower limit in K

Returns This SystemFunction object

**class tc_python.abstract_base.SystemModifications**

Bases: object

Functionality to modify a user database during a calculation by changing phase parameters and system functions.

The actual changes are only applied by using tc_python.abstract_base.AbstractCalculation.with_system_modifications() on a calculator object.

**run_ges_command (ges_command: str)**

Sends a GES-command. This is actually applied when running `with_system_modifications` on a calculator object.

Example: `run_ges_command(“AM-PH-DE FCC_A1 C_S 2 Fe:C”)` for adding a second composition set to the FCC_A1 phase with Fe as major constituent on first sublattice and C as major constituent on second sublattice.

Note: For details about the syntax search the Thermo-Calc help for GES (the name for the Gibbs Energy System module in Console Mode).
Note: It should not be necessary for most users to use this method, try to use the corresponding method implemented in the API instead.

Warning: As this method runs raw GES-commands directly in the engine, it may hang the program in case of spelling mistakes (e.g. forgotten parenthesis, ...).

Parameters ges_command – The GES-command (for example: “AM-PH-DE FCC_A1 C_S 2 Fe:C”)

Returns This SystemModifications object

```python
set(parameter_or_function: Union[tc_python.abstract_base.PhaseParameter,
       tc_python.abstract_base.SystemFunction])
```

Overwrites or creates a phase parameter or system function.

Example: `system_modifications.set(PhaseParameter('G(LIQUID,FE;0)').set_expression_with_upper_limit('+1.2*GFELIQ'))`

Example: `system_modifications.set(SystemFunction("DGDEF").set_expression_with_upper_limit('+10.0-R*T', 1000).set_expression_with_upper_limit('+20.0-R*T', 3000))`

Note: The old parameter/function is overwritten and any temperature intervals not defined are lost.

Note: Please consult the Thermo-Calc GES-system documentation for details about the syntax.

Returns This SystemModifications object

```python
class tc_python.abstract_base.TemperatureInterval(expression: Union[str, object],
       upper_temperature_limit: float)
```

Temperature interval expression used by the classes SystemFunction and PhaseParameter.

Parameters

- **expression** – The temperature function expressed in Thermo-Calc database syntax.
- **upper_temperature_limit** – The upper temperature limit in K

```python
get_expression() → str
```

Returns the function expression of this temperature interval.

Returns The temperature function expression

```python
get_upper_temperature_limit() → float
```

Returns the upper limit of this temperature interval.

Returns The upper temperature limit in K

```python
set_expression(expression: str)
```

Sets the function expression of this temperature interval.

Parameters expression – The temperature function expression

```python
set_upper_temperature_limit(upper_temperature_limit: float)
```

Sets the upper limit of this temperature interval.
Parameters `upper_temperature_limit` – The upper temperature limit in K
This section provides an FAQ for common problems that occur when using TC-Python.

6.1 Diagnostics script

If you have problems running TC-Python, run the diagnostics script below.

On Linux you can alternatively download the script directly into your current working directory by:

curl -O https://www2.thermocalc.com/downloads/support/diagnostics-py/tc-python-
↪ diagnostic-script-2022a.py

```python
###
Run this script when troubleshooting TC-Python

It is important to run this script EXACTLY the same way as you run your TC-Python
script
(In the same IDE, same project, same Python environment, same Jupyter notebook e.t.c)
###

version = '2022a'

print('Testing TC-Python version: ' + version)
print('Please make sure that the variable "version" above, matches the release that
-you want to test, if not change it and re-run this script."

# below this line, nothing needs to be manually updated.

import sys
print('')
print('Python version: (should be at least 3.5 and can NOT be older than 3.0)')
print(str(sys.version_info[0]) + '.' + str(sys.version_info[1]))
if sys.version_info[0] < 3 or sys.version_info[1] < 5:
    print('Wrong version of Python !!!!!')
print('')
print('Python executable path: (gives a hint about the used virtual / conda
-environment, in case of Anaconda the corresponding \n'
    'environment name can be found by running `conda env list` on the Anaconda
-command prompt, '
    'TC-Python must be installed into \nEACH separate environment used!')
print(sys.executable)
```

(continues on next page)
import os

print('')
print('Thermo-Calc ' + version + ' installation directory: (must be a valid path to a_
complete installation of ' + version + ')')
tc_env_variable = 'TC' + version[2:].upper() + '_HOME'

try:
    print(os.environ[tc_env_variable])
except:
    print('No Thermo-calc environment variable for ' + version + ' was found. (' + tc_
_env_variable + ')')

print('')
print('Url of license server: (if license server is NO-NET, you need a local license_
file)'

try:
    print(os.environ['LSHOST'])
except:
    print('No Thermo-calc license server url was found. (LSHOST)'

print('')
print('Path to local license file: (only necessary if not using license server)'

try:
    print(os.environ['LSERVRC'])
except:
    print('No path to local license file was found. (LSERVRC)'

import tc_python

numerical_version = version[:-1]

if version[-1] == 'a':
    numerical_version += '.1.*'

elif version[-1] == 'b':
    numerical_version += '.2.*'

print('')
print('TC-Python version: (needs to be ' + numerical_version + ')')
print(tc_python.__version__)

with tc_python.TCPython() as session:
    print('')
    print('Lists the databases: (should be a complete list of the installed databases_
that you have license for or do not require license)'
    print(session.get_databases())
6.2 “No module named tc_python” error on first usage

This problem occurs because your used Python interpreter cannot find the TC-Python package. We expect that you have installed the TC-Python package in your Python system interpreter following the instructions in the Installation Guide.

Normally the error message “No module named tc_python” is caused by unintentionally configuring a PyCharm project to use a so-called Virtual Environment. This happens unfortunately by default when creating a new PyCharm project with not changing the default settings.

Note: A Virtual Environment is basically a separate and completely independent copy of the system-wide Python interpreter. It does not contain any packages.

On Windows systems we recommend to use the Anaconda Python Distribution as Python interpreter. However, the instructions given here are valid for any operating system and distribution.

Since TC-Python 2018b we do recommend to not use Virtual Environments unless there is a reasonable use case for that.

There are two possible solutions to fix the problem:

1. The quick fix for your problem is to run

   ```bash
   pip install <path to the TC-Python folder>/TC_Python-<version>-py3-none-any.whl
   ```

   within the Terminal window of the opened PyCharm project. This Terminal window automatically runs within the Virtual Environment configured for the project (if any). You can see the name of the Virtual Environment at the beginning of each command prompt line (here it is called venv):

   ```
   Microsoft Windows [Version 10.0.16299.431]
   (c) 2017 Microsoft Corporation. All rights reserved.
   (venv) C:\Users\User\Documents>
   ```

   The command will consequently install TC-Python also within that Virtual Environment automatically. The Terminal window can be found at the bottom of the IDE. Note that it might be necessary to enable these buttons first by selecting the menu entry View→Tool Buttons.

2. The better fix is to change your project to use the system interpreter. This is described in detail in the section Fixing potential issues with the environment in Step 5 of the Installation Guide.

   ```
   It is recommendable to use that approach also for all your future projects.
   ```

Both fixes will only change the configuration of the opened project. Further useful information can be found in the section Python Virtual Environments.
6.3 “pip install” fails with “Failed to establish a new network connection” or similar

If `pip install` fails with a network related error (might also be “socket not available”, “retrying after connection broken”, ...) it is often due to the computer being behind a proxy-server, this is common in large organizations. Of course also the network connection might be broken.

TC-Python has dependencies to a few other packages:

- `py4j`
- `jproperties`
- `six` (transient dependency of `jproperties`)

1. The recommended approach is to simply use `pip`. It will resolve the dependencies automatically by downloading them from the `PyPI`-repository server (https://pypi.org). If your computer is located behind a proxy-server, the connection to the repository will fail. In that case it is necessary to configure `pip` with the detailed configuration of the proxy server:

   ```bash
   pip install -proxy user:password@proxy_ip:port py4j jproperties
   ```

2. Another alternative is to manually download the latest *.whl*-file of each dependency from the repository server (https://pypi.org -> Search projects) and to install it manually using:

   ```bash
   pip install py4j-#.#.#-py2.py3-none-any.whl
   ...
   ```

   The actual actual version number needs to be inserted into the file name. The downside of this approach is that updates to that package have to be fully manual also in the future. Additionally it is also necessary to install all transient dependencies in that way.
tc_python.abstract_base, 227
tc_python.batch_equilibrium, 48
tc_python.diffusion, 112
tc_python.entities, 186
tc_python.exceptions, 225
tc_python.material_to_material, 155
tc_python.precipitation, 52
tc_python.propertymodel, 150
tc_python.propertymodel_sdk, 209
tc_python.quantity_factory, 195
tc_python.scheil, 74
tc_python.server, 190
tc_python.single_equilibrium, 35
tc_python.step_or_map_diagrams, 85
tc_python.system, 178
tc_python.utils, 207